



Shaw Environmental, Inc.

11 Northeastern Boulevard
Salem, NH 03079-1953
603.870.4500
Fax: 603.870.4501

March 27, 2007
Project 101960

Mr. Joseph T. Martella, II
Rhode Island Department of Environmental Management
Office of Waste Management
235 Promenade Street
Providence, RI 02908-5767

**Re: Status Report: December 2006 - March 2007 Activities
Former Gorham Manufacturing Facility
333 Adelaide Avenue, Providence, RI
Site Remediation Case No. 97-030**

Dear Mr. Martella:

Shaw Environmental, Inc. (Shaw) has prepared this quarterly status report on behalf of Textron, Inc. (Textron). This status report is associated with the remediation of tetrachloroethene (PCE) contaminated groundwater at the former Gorham Manufacturing Facility at 333 Adelaide Avenue, Providence, Rhode Island (Figure 1).

PCE is the primary contaminant of concern for groundwater. As discussed in the Remedial Action Work Plan (RAWP) and subsequent revisions, the PCE source area in the vicinity of the former building W is the area of concern with a site-specific remedial goal of 7,700 micrograms per liter (ug/L). This area was treated using in-situ applications of sodium permanganate. Figure 2 shows the most recent treatment area.

This status report describes groundwater monitoring activities conducted in accordance with the proposed groundwater monitoring program submitted to the Rhode Island Department of Environmental Management (RIDEM) in February 2007 (Shaw – Groundwater Monitoring Program letter, dated February 1, 2007). This status report also describes activities conducted in accordance with a letter proposing to conduct a laboratory treatability study in the source area to evaluate enhanced bioremediation submitted to RIDEM in August 2006 (Shaw – Laboratory Treatability Study letter, dated August 21, 2006). Shaw collected soil and groundwater samples on December 6, 2006. The laboratory treatability testing results are included in this report.

In addition, Textron has installed and sampled two new wells (MW-220S and MW-221S) (Figure 1) in accordance with a letter submitted to RIDEM in February 2007 (Textron – Parcel A Well Installations, dated February 9, 2007). The purpose of this activity is to provide groundwater monitoring points between the treatment area and the retail building that will enhance our understanding of current groundwater conditions in the area.

TREATABILITY STUDY

Field Activities

Soil and groundwater samples were collected from the apparent PCE source area for a bioaugmentation treatability study conducted to evaluate present subsurface conditions for bioremediation. On December 6, 2006, two soil borings (SB-219A and SB-219B) were advanced in the source area based on recent soil and groundwater data collected during the subsurface investigation conducted in April 2006. Discrete soil samples were collected from one soil boring, SB-219A, at two depth intervals where the highest PCE concentrations were detected in soil samples collected from soil boring, SB-210 in April 2006. A soil sample was collected from the 27 to 31 feet bgs, and a soil sample was collected from 43 to 47 feet bgs. The locations of the soil borings and subsequent well installations are shown on Figure 1. The boring logs are contained in Appendix A. The drilling services were performed by Pine and Swallow Associates (Pine) of Groton, MA and overseen by a Shaw Environmental field geologist.

Prior to drilling, each subsurface location was cleared to a depth of 5 feet below ground surface (bgs) by manual digging. Soil samples from each boring were collected at four foot intervals. Soil samples were collected via dedicated plastic sample liners using a VibraDrill H641 drill rig. Soil samples were collected for geologic logging and field screening for the presence of volatile organic vapors and residual source VOCs. The selected sample intervals identified above were sent to the laboratory for microcosm testing. The soil samples used for testing were collected under anaerobic conditions, which required that the sample be contained within the dedicated plastic liner, capped at each end immediately upon retrieval, sealed at each cap with tape and wrapped in plastic bags for shipment to the laboratory. The samples collected for the treatability study were submitted to Shaw's treatability laboratory in Lawrenceville, New Jersey for testing.

Each boring was completed as a well for possible future remediation use. Each monitoring well consisted of 1.32-inch nominal steel casing and a 4-foot section of well screen consisting of 2-inch by 0.015-inch vertical slots. Further details on the screen interval at each boring are provided in the table provided below.

| Treatability Study Well Installations | | | |
|--|-------------------------------------|------------------------|--|
| Well ID | Location | Screen Interval | SB-210 PCE Concentration (ug/kg) |
| MW-219A | Approximately 2 west of SB-210 | 43 to 47 ft. bgs | 4,400,000 (at SB-210, 43 to 46 feet bgs) |
| MW-219B | Approximately 2 feet east of SB-210 | 27 to 31 ft. bgs | 77,000,000 (at SB-210, 30-31 feet bgs) |

Soil samples collected from each interval (with the exception of the intervals sent to the laboratory for the treatability study) were placed in jars for headspace screening and screened for volatile organic vapors using a Mini Rae 2000 photoionization detector (PID) equipped with a 10.6 electron-volt lamp. PID readings ranged from 7.9 to 9,999 parts per million by volume (ppmv). The highest PID screening values were detected at SB-219A. Generally, higher PID values were measured at the water table and within the saturated zone for each boring.

In addition, soil samples collected from each interval (with the exception of the intervals sent to the laboratory for the treatability study) were screened for the presence of residual source VOCs by use of the OilScreenSoil (Sudan IV)TM dye test kit. Soil was placed in a Sudan IV test bottle to the line indicated on the bottle and hot water was added to the test bottle to the line indicated on the bottle. The test bottle containing soil and hot water was shaken until the red Azo dye completely dissolved in the contents of the test bottle. The test indicated the presence of total petroleum hydrocarbon (TPH) or other compound, at a concentration greater than 500 ppm, in soil boring SB-219A at 34-39 feet bgs. The test did not indicate the presence of dense non-aqueous phase liquid (DNAPL).

Groundwater was collected from an existing Site well in the source area (MW-101D) for use in the treatability study. The groundwater was collected via a modified low flow methodology. MW-101D was purged for 30 to 40 minutes prior to collecting a sample. Field parameters were recorded and included: oxidation/reduction potential (ORP); dissolved oxygen (DO); pH, temperature; specific conductance (SC); and turbidity. Initial and final water level measurements were recorded.

The groundwater samples were submitted to Shaw's treatability laboratory in Lawrenceville, NJ for analysis of volatile organic compounds (VOCs) by EPA Method 8260B, total and dissolved iron and manganese, reduced gases - methane, ethane, ethene (M/E/E) by EPA Method 8015 and anions by EPA Method 300. In addition, a

groundwater sample was collected from MW-101D for microcosm testing. All samples were collected under a nitrogen blanket to minimize exposure to atmospheric air and the sample bottle was completely filled in order to eliminate the headspace within the sample bottle, which was capped with a Teflon-lined cap. A discussion of the laboratory treatability study is provided below and a copy of the laboratory analytical reports are included in Appendix B.

Laboratory Treatability Study

The objective of the treatability study was to assess the effectiveness of bioaugmentation for treatment of PCE in soil and groundwater collected from the former Gorham Manufacturing Facility. The study was performed between December 2006 and February 2007. Detailed descriptions of the methodologies, results, and conclusions are presented below.

Groundwater used for the laboratory treatability study was collected from monitoring well MW-101D; soil used was collected from an adjacent location at a depth interval of approximately 27-31 ft bgs. These sample locations are located within the presumed PCE source area.

Microcosms were prepared in glass serum bottles (approximate volume, 160 mL). All microcosm preparation and sampling was performed in an anaerobic chamber. Thirty grams of homogenized Site soil and 135-mL of Site groundwater were added to each of the bottles. A total of 25 bottles were prepared. The bottles were sealed with Teflon®-lined butyl rubber stoppers and crimp caps.

Bottles were incubated with gentle shaking overnight at 15°C, after which the aqueous phase was analyzed for volatile organic compounds (VOCs) via EPA Method 8260. Aqueous samples were analyzed with a 24-hour turnaround time to determine whether PCE levels were representative of the elevated concentrations observed in the field. Since the results indicated PCE levels of only 3-6 mg/L in the groundwater sample, it was determined that a contaminant spike of PCE was needed so that PCE concentrations in the microcosms more representative of Site conditions. PCE was added to all microcosms to achieve a final PCE concentration of approximately 20 mg/L.

Four sets of microcosm treatments were prepared in triplicate as follows:

Treatment 1: KILLED CONTROL: These treatments were amended with a formaldehyde solution (final concentration in groundwater approximately 1.5% by volume) to inhibit microbial activity, and were used to evaluate abiotic loss of VOCs;

Treatment 2: LIVE CONTROL: This treatment did not receive any amendments except for deionized water (to maintain the same liquid volume in the other treatments). This treatment served as a control to monitor VOC loss in the absence of any amendments;

Treatment 3: BIOAUGMENTATION 1: Shaw's SDC-9 culture was used as the bacterial inoculum. Nutrient solution and yeast extract were added to ensure that the bacteria were not limited in nitrogen, phosphorus, or other trace nutrients. Bottles also were amended with lactate to serve as the electron donor (final lactate concentration of 1,000 mg/L). This treatment was used to evaluate the effects of anaerobic bioaugmentation on contaminant biodegradation.

Treatment 4: BIOAUGMENTATION 2: Shaw's SDC-9 culture was used as the bacterial inoculum. Nutrient solution and yeast extract were also added to ensure that the bacteria were not limited in nitrogen, phosphorus, or other trace nutrients. Bottles also were amended with an emulsified vegetable oil (EVO) to serve as the electron donor (final EVO concentration of 1,000 mg/L). This treatment was used to evaluate the effects of anaerobic bioaugmentation on contaminant biodegradation.

A parallel set of bottles were also prepared for pH and ORP monitoring, and extra bottles were prepared for a final soil analysis (if needed) at the completion of the study.

Microcosm bottles were incubated with gentle shaking at 15°C. At each sampling event, microcosm bottles were removed from the shaker and placed in the anaerobic chamber. Sufficient time was allowed for the soils to settle so that the supernatant groundwater could be sampled. Samples were collected from all treatments at t= 0, 8, 15, 30, and 51 days. At each of these sampling events, groundwater was analyzed for VOCs and reduced gases. In addition, one bottle from each treatment was analyzed for volatile fatty acids, anions, pH, and ORP at each sampling event (equal sample volume were collected from the other bottles and treatments to maintain equal groundwater volumes in all the treatments). Glass beads were placed in the bottles after sampling to maintain zero headspace.

Additional lactate was added to Treatment 3 at t=7 days. Treatments 3 and 4 were re-inoculated with SDC-9 at t= 33 days.

Treatability Study Results

Results of the VOC analyses are presented in Table 1. Only results for PCE are shown, since concentrations of the expected dechlorination daughter products (trichloroethene, cis-1,2-dichloroethene, and vinyl chloride) were below the analytical detection limit (approximately 4 mg/L) throughout the study. Comparison among the controls and bioaugmented treatments indicate that addition of the bioaugmentation amendments did not result in a measurable reduction in PCE concentrations (i.e., no PCE degradation was observed). The increasing trend in PCE concentrations among the four treatments between days 8 and 51 is likely due to PCE desorption from aquifer solids and/or dissolution of non-aqueous phase PCE liquids.

Reduced gas concentrations (methane and ethane) for each treatment are shown in Table 2; ethene concentrations were below the analytical detection limit of 2 µg/L throughout

the study. Results indicate that only trace levels of ethane were present in the bioaugmented samples. These ethane levels were approximately equal to the ethane levels in the controls, and were orders of magnitude below the stoichiometric concentrations expected for complete dechlorination of PCE. This result is consistent with the absence of any measurable PCE degradation, and indicates that PCE dechlorination was not occurring at a measurable rate in the microcosms.

In addition, methane concentrations in the bioaugmented treatments were approximately equal to methane concentrations in the controls, indicating that methanogenic activity (which typically accompanies dechlorination using SDC-9) in the amended treatments was negligible.

Sulfate concentrations are shown in Table 3. No decreases in sulfate were observed in any of the treatments, indicating that biological activity was insufficient to attain sulfate-reducing conditions (sulfate reduction is typically observed prior to PCE dechlorination).

Volatile fatty acids are shown in Table 4. Results indicate that lactate (i.e., electron donor) levels were sufficiently high throughout the study in the lactate-amended bioaugmentation treatment. However, the absence of substantial quantities of fermentation daughter products (relative to the lactate dosage) suggests that substantial lactate fermentation was not occurring.

ORP and pH levels are shown in Table 5. The positive ORP values are consistent with the observed lack of sulfate and PCE reduction, as well as the absence of substantial lactate fermentation. The pH levels generally range between 6 and 8, which are typically suitable for PCE dechlorination using SDC-9.

Treatability Study Conclusions

Overall results of this laboratory treatability study indicate that bioaugmentation is not effective for treating PCE in Site soil and groundwater. The VOC, reduced gas, sulfate, volatile fatty acid, and ORP data were all consistent, indicating that biological activity in the bioaugmented samples was limited, and that reducing conditions needed for PCE dechlorination were not attained.

Results of this treatability study indicate that bioaugmentation is likely not a feasible remedial option for PCE in the source area. As with any laboratory microcosm testing, results should always be interpreted with the understanding that the relatively small closed systems used in the study may not completely capture the appropriate diversity of mechanisms present at the field scale.

GROUNDWATER MONITORING AND WELL INSTALLATIONS

Groundwater Sampling and Monitoring Activities

On February 17, 2007, twenty (20) groundwater samples (including one duplicate) were collected for analysis for volatile organic compounds (VOCs) (EPA Method 8260B) and two (2) groundwater samples (including one duplicate) were collected for analysis for total petroleum hydrocarbon (TPH) (EPA Method 8015B). A duplicate sample was collected for VOC analysis from MW-101S and a duplicate sample was collected from CW-6 for TPH analysis. Groundwater samples were collected by first purging approximately three well volumes from each well and then collecting a sample in a dedicated bailer. Groundwater samples were delivered to AMRO Environmental Laboratories Corporation in Merrimack, New Hampshire for analysis.

Field measurements were also taken from the wells sampled in the treatment area and included oxidation/reduction potential (ORP), dissolved oxygen (DO), pH, temperature, and specific conductance (SC). Groundwater elevation measurements were also collected from both the treatment area wells and the compliance wells. These results are presented in Tables 6 and 7.

Well Installations

On February 22, 2007, Shaw oversaw Technical Drilling Services, Inc (TDS) of Sterling, MA in the installation of two additional groundwater monitoring wells (MW-220S and MW-221S). Prior to drilling, each subsurface location was cleared to a depth of 5 feet below ground surface (bgs) by manual digging.

The monitoring wells were advanced via 4.25 inch hollow stem augers, using a track mounted CME-75 hollow stem auger drill rig. Split spoon samples were collected at 5 foot intervals beginning at 5 feet below ground surface. The split spoon samples were logged in the field using Unified Soil Classification System (USCS). Each split spoon was collected for jar headspace screening. Headspace was collected with a Mini Rae 2000[®] Photo Ionization Detector (PID) that was field calibrated with Isobutylene Balance Air to 100 parts per million.

During drilling elevated PID readings were noted in the boring for MW-221S. PID readings ranged from 3.7 to 70 ppm. Dark black staining on soil was noted at approximately 25 to 27 feet bgs and staining and a sheen was observed on soil at approximately 30 to 32 feet bgs. During drilling for well MW-220S PID readings were low and no staining or sheens were noted. A septic odor was noted a 10 feet bgs.

MW-220S and MW-221S were advanced approximately 7 feet into the water table (32 feet below ground surface) The wells were completed with 2 inch PVC, and screened with 0.010 inch slotted screen from 22 to 32 feet bgs. Solid PVC riser extended to the ground surface. The annulus of the screen interval was then backfilled with silica sand to 20 feet below ground surface (2 feet above the screened interval). Approximately 2 feet

of Bentonite chips were then placed above the sand pack to ensure a proper seal. The wells were then finished with a 6 inch flush mount, bolt down roadboxes set in concrete. The boring logs are contained in Appendix A. Both wells are located within 30 feet of the edge of the building (Figure 1)

All associated waste from the drilling activities (i.e. soil cuttings) were drummed, labeled, and staged on Site. During the well installation (2) 55 gallon DOT drums of soil were produced. All waste generated during this phase of field activities (i.e. soil cuttings and purge water) will be transported off Site, as hazardous waste, and disposed of accordingly.

Well Development, Sampling, and Surveying

On February 26, 2007 Shaw returned to the Site to purge and develop the newly installed wells. The wells were gauged to determine the depth to water, possible presence of non-aqueous phase liquid (NAPL), depth to bottom, and total water column thickness. NAPL was not detected in either well. Approximately 25 gallons were pumped from each well using a 2 inch stainless steel Grundfos® pump, until turbidity levels were visually clear. The purge water was containerized in (1) 55 gallon DOT drum labeled, and staged onsite.

On March 1, 2007 Shaw returned to the Site to sample the newly installed wells (MW-220S and MW-221S) two additional wells, MW-109D and GZA-3, and survey in the new well locations. Each well was sampled via purge method. Prior to purging, the wells were gauged to determine the depth to water, NAPL, depth to bottom, and total water column thickness. Water column thickness was used to calculate the water volume within the well. Using a dedicated polyethylene bailer and dedicated foot valve and tubing (to prevent cross contamination) a minimum of three well volumes were evacuated from each well prior to collecting the sample. MW-109D, MW-220S and MW-221S were all sampled for VOCs (method EPA 8260B), in addition MW-221S was sampled for TPH (method EPA 8015B). Well GZA-3 was sampled for total dissolved lead, (method EPA SW 846). (Note: since well GZA-5 has been destroyed and well GZA-6 was not accessible, results from wells GZA-3 and MW-109D were utilized for Mashapaug Pond compliance evaluation purposes during this sampling event.) The sample for lead analysis was field filtered and acid preserved. All samples was preserved and transported on ice under a valid chain of custody back to the Salem, NH office, where a representative from AMRO Environmental Laboratories Corp. (Merrimack, NH), delivered the samples to the laboratory for analysis.

SUMMARY OF GROUNDWATER ANALYTICAL DATA

A summary of all the analytical data associated with the groundwater sampling conducted in February and March 2007 is contained in Table 8. A copy of the laboratory analytical report is attached as Appendix B of this report. The PCE concentrations found in wells MW-101S, MW-101D, MW-202D, MW-202S, MW-207D, and MW-207S are currently at or above the treatment goal of 7,700 ug/L.

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A summary of the compliance well results is contained in Table 9. The results for the compliance wells indicate that exceedances occurred for wells MW-209D (PCE), MW-218D (PCE, TCE, and 1,1-DCE), and well MW-218S (PCE) .

FUTURE ACTIVITIES

The next quarterly sampling event is scheduled to be conducted in May 2007.

If you have any questions, please contact Ed Van Doren at (603) 870-4530.

Sincerely,

SHAW ENVIRONMENTAL, INC.

Edward P. Van Doren, PE, LSP
Project Manager

cc: Craig Roy, RIDEM OWR
Greg Simpson, Textron
Dave McCabe, Textron
Jamieson Schiff, Textron
Dave Heislein, MACTEC
Thomas Dellar, City of Providence
Jeff Morgan, Stop & Shop
Ronald Ruth, Sherin and Lodgen

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Attachments:

Figure 1 – Site Plan

Figure 2 – Injection Well Locations

Table 1 – Measured PCE Aqueous Concentrations

Table 2 – Measured Reduced Gas Concentrations

Table 3 – Measured Sulfate Concentrations

Table 4 – Measured Volatile Fatty Acid Concentrations

Table 5 – ORP and pH Readings

Table 6 – Summary Field Parameters

Table 7 – Water Table Elevations

Table 8 – Groundwater Analytical Results

Table 9 – Compliance Wells Analytical Results

Appendix A – Boring Logs

Appendix B – Laboratory Analytical Report: Bioaugmentation

Appendix C – Laboratory Analytical Report: Groundwater Sampling

Mr. Joseph T. Martella, II

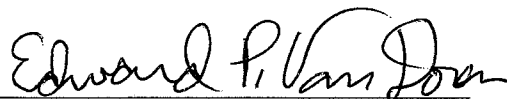
March 27, 2007

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CERTIFICATIONS

The following certifications are provided pursuant to Rule 9.19 of the Remediation Regulations:

I, Edward P. Van Doren, as an authorized representative of Shaw Environmental, Inc. and the person responsible for the preparation of this Status Report dated March 27, 2007, certify that the information contained in this report is complete and accurate to the best of my knowledge.



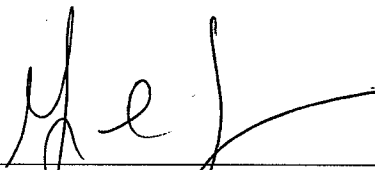
Edward P. Van Doren, PE, LSP
Project Manager

04/02/2007

Date:

We, Textron, Inc., as the party responsible for submittal of this Status Report, certify that this report is a complete and accurate representation of the contaminated site and the release, and contains all known facts surrounding the release, to the best of our knowledge.

Certification on behalf of Textron Inc.





Gregory L. Simpson
Project Manager

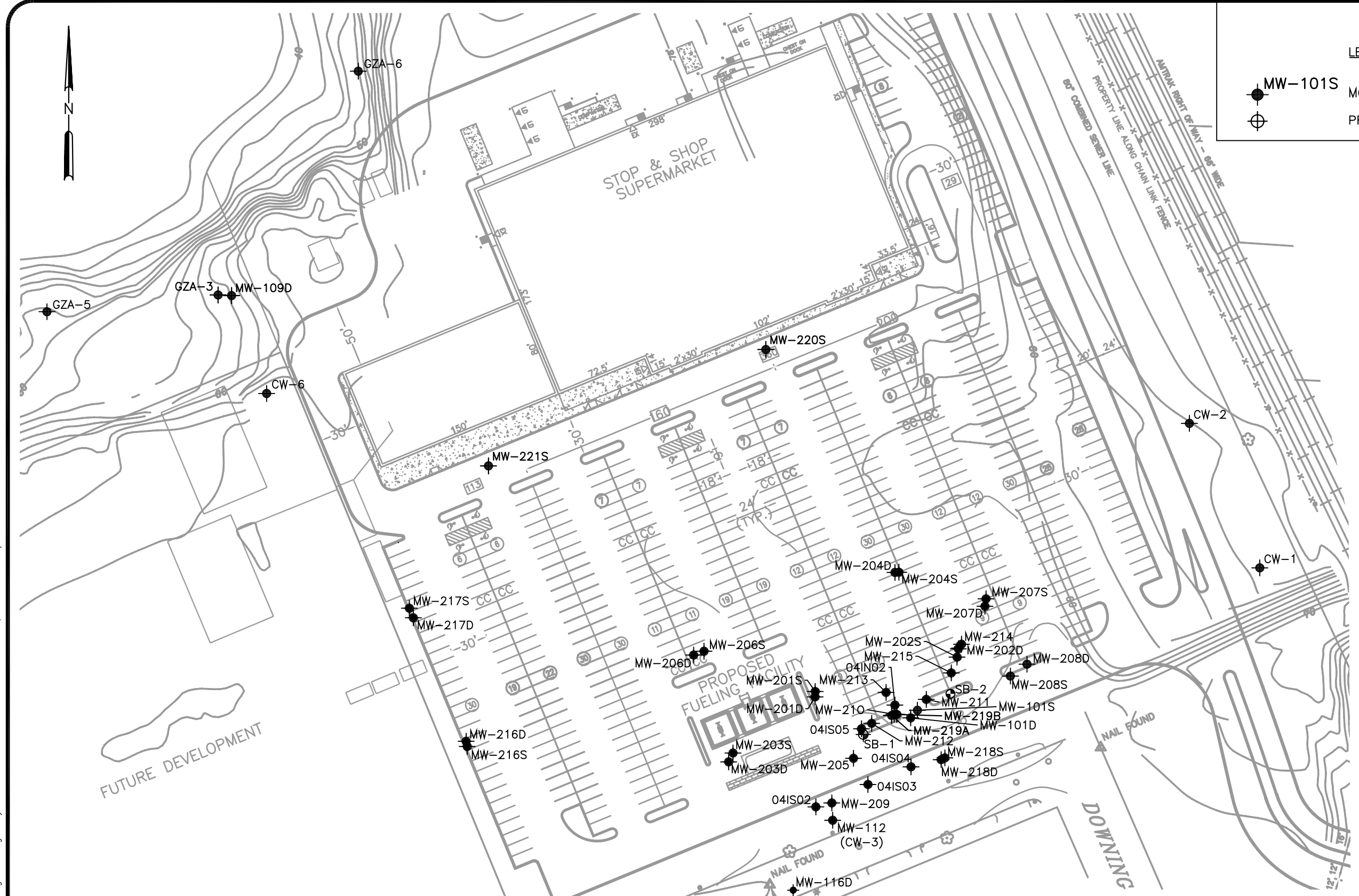
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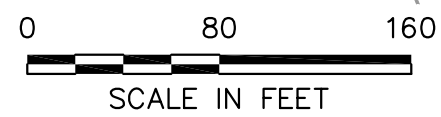
FIGURES

LEGEND

-  MW-101S MONITORING WELL
-  PROPOSED WELL LOCATION



File: N:\dwg\Gorham\antgf-14.dwg Layout: SP User: James.O'Donnell Mar 08, 2007 - 5:31pm
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|-------------|---------|
| DATE | 1/30/07 |
| DWN | J.O'D. |
| APP | |
| REV | |
| PROJECT NO. | 101960 |

FIGURE 1
TEXTRON PROVIDENCE
 333 ADELAIDE AVENUE
 PROVIDENCE, RHODE ISLAND
SITE PLAN

TABLES

Table 1
Measured PCE Aqueous Concentrations

| Days | Killed Control | Live Control | Bioaug. - Lactate | Bioaug.-EVO |
|------|----------------|--------------|-------------------|-------------|
| 0 | 4.6 ± 2.0 | 3.7 ± 0.82 | 3.9 ± 1.7 | 3.1 ± 0.78 |
| 8 | 17 ± 1.2 | 11 ± 5.7 | 14 ± 3.6 | 7.2 ± 6.0 |
| 15 | 20 ± 1.0 | 15 ± 5.0 | 20 ± 6.0 | 18 ± 2.1 |
| 30 | 27 ± 1.0 | 19 ± 3.1 | 26 ± 5.3 | 15 ± 2.8 |
| 51 | 24 ± 2.5 | 19 ± 2.5 | 23 ± 2.9 | 20 ± 5.1 |

NOTES:

Average values among triplicate samples are shown (± standard deviation).

PCE spike after 24 hours

Visibly outlying data were discarded

Concentrations are in mg/L

Table 2
Measured Reduced Gas Concentrations

Methane

| Days | Killed Control | Live Control | Bioaug. - Lactate | Bioaug.-EVO |
|------|----------------|----------------|-------------------|----------------|
| 0 | NS | NS | NS | NS |
| 8 | 0.019 ± 0.0083 | 0.023 ± 0.0051 | 0.017 ± 0.0079 | 0.015 ± 0.0030 |
| 15 | 0.018 ± 0.0017 | 0.027 ± 0.0043 | 0.023 ± 0.0016 | 0.014 ± 0.0012 |
| 30 | 0.021 ± 0.0015 | 0.023 ± 0.0018 | 0.016 ± 0.0058 | 0.018 ± 0.0014 |
| 51 | 0.017 ± 0.0073 | 0.025 ± 0.0014 | 0.020 ± 0.0031 | 0.020 ± 0.0081 |

Ethane

| Days | Killed Control | Live Control | Bioaug. - Lactate | Bioaug.-EVO |
|------|-----------------|-----------------|-------------------|-----------------|
| 0 | NS | NS | NS | NS |
| 8 | 0.0077 ± 0.0004 | 0.0096 ± 0.0006 | 0.010 ± 0.0020 | 0.011 ± 0.0010 |
| 15 | 0.0070 ± 0.0006 | 0.0090 ± 0.0023 | 0.012 ± 0.0037 | 0.0087 ± 0.0014 |
| 30 | 0.0021 ± 0.0001 | 0.0021 ± 0.0001 | 0.0022 ± 0.0002 | 0.0046 ± 0.0010 |
| 51 | 0.0014 ± 0.0012 | 0.0021 ± 0.0007 | 0.0022 ± 0.0001 | 0.0025 ± 0.0022 |

NOTES

Average values among triplicate samples are shown (± standard deviation).

NS=not sampled

Concentrations are in mg/L

Table 3
Measured Sulfate Concentrations

| Days | Killed Control | Live Control | Bioaug. - Lactate | Bioaug.-EVO |
|------|----------------|--------------|-------------------|-------------|
| 0 | NS | NS | NS | NS |
| 8 | 32.4 | 39.5 | 44.2 | 45.5 |
| 15 | 33.1 | 39.1 | 46.1 | 46.1 |
| 30 | 39.9 | 43.5 | 42.3 | 45.5 |
| 51 | 42.2 | 42.8 | 46 | 42.3 |

NOTES

Samples were collected from a single bottle for each treatment.

NS=not sampled

Concentrations are in mg/L

Table 4

Measured Volatile Fatty Acid Concentrations

| Treatment | Days | Lactate | Acetate | Propionate | Formate | Butyrate | Pyruvate | Valeric Acid |
|------------------------|------|---------|---------|------------|---------|----------|----------|--------------|
| Killed Control | 0 | NS | NS | NS | NS | NS | NS | NS |
| | 8 | <1.0 | <1.0 | <1.0 | 177 | <1.0 | <1.0 | <1.0 |
| | 15 | <10.0 | 1.98 | <10.0 | 193 | <10.0 | <10.0 | <10.0 |
| | 30 | <10.0 | 2.54 | <10.0 | 178 | <10.0 | <10.0 | <10.0 |
| | 51 | <10.0 | 2.27 | <10.0 | 192 | <10.0 | <10.0 | <10.0 |
| Live Control | 0 | NS | NS | NS | NS | NS | NS | NS |
| | 8 | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| | 15 | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| | 30 | <1.0 | 0.41 | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| | 51 | <1.0 | 0.59 | <1.0 | <1.0 | <1.0 | <1.0 | <1.0 |
| Bioaug.-Lactate | 0 | NS | NS | NS | NS | NS | NS | NS |
| | 8 | 1370 | 51.5 | 22.5 | 4.5 | <1.0 | <1.0 | <1.0 |
| | 15 | 2620 | 53.2 | 20.5 | <50.0 | <50.0 | <50.0 | <50.0 |
| | 30 | 3080 | 139 | 28.9 | 16.8 | <50.0 | <50.0 | <50.0 |
| | 51 | 2740 | 166 | 51.6 | 23.1 | <50.0 | <50.0 | <50.0 |
| Bioaug. - EVO | 0 | NS | NS | NS | NS | NS | NS | NS |
| | 8 | 31.0 | 33.4 | 24.1 | 8.01 | <1.0 | <1.0 | <1.0 |
| | 15 | <1.0 | 55.4 | 14.1 | <1.0 | <1.0 | <1.0 | <1.0 |
| | 30 | <10.0 | 107 | 66.8 | 16.9 | 18.0 | 3.64 | <10.0 |
| | 51 | <10.0 | 130 | 82.5 | 16.3 | 25.8 | 2.71 | <10.0 |

NOTES

Samples were collected from a single bottle for each treatment.

NS=not sampled

Concentrations are in mg/L

Formate detections in the Killed Control are likely formaldehyde, which was used as the microbial inhibitor

Table 5**ORP and pH Readings**

| Treatment | Days | pH | ORP |
|-----------------------------|-------------|-----------|------------|
| Killed Control | 0 | NS | NS |
| | 8 | 5.44 | 167 |
| | 15 | 6.80 | 132 |
| | 30 | 5.90 | 129 |
| | 51 | 6.74 | 102 |
| Live Control | 0 | NS | NS |
| | 8 | 6.82 | 178 |
| | 15 | 7.00 | 116 |
| | 30 | 7.25 | 130 |
| | 51 | 8.33 | 131 |
| Bioaug.- Lactate | 0 | NS | NS |
| | 8 | 6.87 | 157 |
| | 15 | 6.74 | 118 |
| | 30 | 6.97 | 90 |
| | 51 | 7.29 | 52 |
| Bioaug. - EVO | 0 | NS | NS |
| | 8 | 6.95 | 162 |
| | 15 | 6.88 | 87 |
| | 30 | 7.00 | 107 |
| | 51 | 7.46 | 70 |

NOTES

Samples were collected from a single bottle for each treatment.

NS=not sampled

pH values in standard units

ORP values in millivolts

**Table 6
Well Field Parameters
Textron Gorham
Providence, Rhode Island**

| WELL ID | DATE | pH | Temperature (deg.c) | Conductivity (ms/cm) | Dissolved Oxygen (mg/l) | Oxidation Reduction Potential (mv) |
|----------------|-------------|-----------|--------------------------------|---------------------------------|--|---|
| MW-101D | 2/17/2007 | 6.32 | 15.20 | 1.475 | 1.30 | 104.4 |
| MW-101S | 2/17/2007 | 6.27 | 15.13 | 0.689 | 1.18 | 61.3 |
| MW-112 | 2/17/2007 | 6.85 | 13.01 | 0.290 | 5.23 | 183.4 |
| MW-116D | 2/17/2007 | 5.61 | 11.63 | 0.133 | 2.82 | 299.0 |
| MW-116S | 2/17/2007 | 5.94 | 12.51 | 0.134 | 4.23 | 298.1 |
| MW-201D | 2/17/2007 | 7.04 | 14.94 | 1.181 | 1.57 | 4.8 |
| MW-202D | 2/17/2007 | 5.83 | 15.48 | 0.923 | 1.39 | 306.1 |
| MW-202S | 2/17/2007 | 5.71 | 15.52 | 0.744 | 1.73 | 259.4 |
| MW-207D | 2/17/2007 | 6.06 | 15.52 | 1.205 | 1.20 | 294.9 |
| MW-207S | 2/17/2007 | 6.41 | 16.38 | 1.124 | 1.24 | 219.3 |
| MW-209D | 2/17/2007 | 6.98 | 13.60 | 0.521 | 2.55 | 128.3 |
| MW-216D | 2/17/2007 | 6.34 | 13.43 | 0.290 | 1.30 | -6.5 |
| MW-216S | 2/17/2007 | 6.55 | 13.57 | 0.577 | 1.02 | -103.0 |
| MW-217D | 2/17/2007 | 6.51 | 13.84 | 0.330 | 2.43 | -53.7 |
| MW-217S | 2/17/2007 | 6.55 | 13.97 | 0.578 | 2.89 | 28.5 |
| MW-218D | 2/17/2007 | 6.37 | 13.88 | 0.446 | 2.32 | 131.5 |
| MW-218S | 2/17/2007 | 6.11 | 14.55 | 0.715 | 2.25 | 123.5 |

Note
C° = degrees Celsius
ms/cm = microsiemens per centimeter
mg/l = milligrams per liter
mV = milli volts

**Table 7
Groundwater Elevations
Textron Gorham
Providence, Rhode Island**

| Well ID | Date | Reference Elevation (Feet) | Depth to Water (Feet) | LNAPL Thickness (Feet) | Groundwater Elevation (Feet) |
|----------------|-------------|---|--------------------------------------|---------------------------------------|---|
| MW-101D | 2/17/2007 | 98.91 | 24.52 | -- | 74.39 |
| MW-101S | 2/17/2007 | 98.90 | 24.55 | -- | 74.35 |
| MW-109D | 3/1/2007 | 90.03 | 19.01 | -- | 71.02 |
| MW-112 | 2/17/2007 | 100.63 | 26.77 | -- | 73.86 |
| MW-116D | 2/17/2007 | 98.92 | 24.41 | -- | 74.51 |
| MW-116S | 2/17/2007 | 99.40 | 24.51 | -- | 74.89 |
| MW-201D | 2/17/2007 | 98.80 | 24.36 | -- | 74.44 |
| MW-202D | 2/17/2007 | 98.17 | 23.85 | -- | 74.32 |
| MW-202S | 2/17/2007 | 98.06 | 23.81 | -- | 74.25 |
| MW-207D | 2/17/2007 | 98.18 | 23.95 | -- | 74.23 |
| MW-207S | 2/17/2007 | 98.28 | 24.02 | -- | 74.26 |
| MW-209D | 2/17/2007 | 100.47 | 25.94 | -- | 74.53 |
| MW-210 | 3/1/2007 | 98.59 | 24.95 | -- | 73.64 |
| MW-213 | 3/1/2007 | 98.89 | 24.88 | -- | 74.01 |
| MW-214 | 3/1/2007 | 98.50 | 24.32 | -- | 74.18 |
| MW-215 | 3/1/2007 | 98.77 | 24.60 | -- | 74.17 |
| MW-216D | 2/17/2007 | 98.69 | 25.12 | -- | 73.57 |
| MW-216S | 2/17/2007 | 99.58 | 25.11 | -- | 74.47 |
| MW-217D | 2/17/2007 | 98.65 | 24.59 | -- | 74.06 |
| MW-217S | 2/17/2007 | 98.71 | 24.60 | -- | 74.11 |
| MW-218D | 2/17/2007 | 99.67 | 25.30 | -- | 74.37 |
| MW-218S | 2/17/2007 | 99.61 | 25.25 | -- | 74.36 |
| MW-219S | 3/1/2007 | 99.29 | 25.04 | -- | 74.25 |
| MW-219D | 3/1/2007 | 99.17 | 24.94 | -- | 74.23 |
| MW-220S | 3/1/2007 | 99.41 | 25.28 | -- | 74.13 |
| MW-221S | 3/1/2007 | 98.92 | 25.39 | -- | 73.53 |
| CW-1 | 2/17/2007 | 99.53 | 25.43 | -- | 74.10 |
| CW-2 | 2/17/2007 | 98.86 | 24.56 | -- | 74.30 |
| CW-6 | 2/17/2007 | 99.52 | 26.11 | -- | 73.41 |

Notes:
Groundwater elevations are based on an arbitrary reference datum established for the site.

Table 8
Groundwater Analytical Results
Textron/Gorham
Providence, Rhode Island

| Sample Identifier Date Sampled Constituent | CW-01 2/17/2007 Primary | CW-02 2/17/2007 Primary | CW-06 2/17/2007 Primary | CW-06 2/17/2007 Duplicate 1 | MW-101D 2/17/2007 Primary | MW-101S 2/17/2007 Primary | MW-101S 2/17/2007 Duplicate 1 | MW-109D 3/1/2007 Primary | MW-112 2/17/2007 Primary | MW-116D 2/17/2007 Primary | MW-116S 2/17/2007 Primary |
|--|-------------------------------|-------------------------------|-------------------------------|-----------------------------------|---------------------------------|---------------------------------|-------------------------------------|--------------------------------|--------------------------------|---------------------------------|---------------------------------|
| Method 8260 (ug/l) | | | | | | | | | | | |
| 1,1,1-Trichloroethane | <200 | <2 | *** | *** | <20 | <200 | <200 | <2 | <2 | <2 | <2 |
| 1,1-Dichloroethane | <200 | <2 | *** | *** | <20 | <200 | <200 | <2 | <2 | <2 | <2 |
| 1,1-Dichloroethene | 180 | <1 | *** | *** | 11 | <100 | <100 | <1 | <1 | <1 | <1 |
| 1,2,4-Trimethylbenzene | <200 | <2 | *** | *** | <20 | <200 | <200 | <2 | <2 | <2 | <2 |
| 1,3,5-Trimethylbenzene | <200 | <2 | *** | *** | <20 | <200 | <200 | <2 | <2 | <2 | <2 |
| 4-Isopropyltoluene | <200 | <2 | *** | *** | <20 | <200 | <200 | <2 | <2 | <2 | <2 |
| Benzene | <100 | <1 | *** | *** | <10 | <100 | <100 | <1 | <1 | <1 | <1 |
| Bromodichloromethane | <200 | <2 | *** | *** | <20 | <200 | <200 | <2 | <2 | 4.3 | 4.2 |
| Carbon tetrachloride | <200 | <2 | *** | *** | 32 | <200 | <200 | <2 | <2 | <2 | <2 |
| Chloroethane | <500 | <5 | *** | *** | <50 | <500 | <500 | <5 | <5 | <5 | <5 |
| Chloroform | <200 | <2 | *** | *** | 20 | <200 | <200 | <2 | <2 | 39 | 32 |
| cis-1,2-Dichloroethene | 350 | <2 | *** | *** | 280 | 510 | 500 | <2 | <2 | <2 | <2 |
| Dichlorodifluoromethane | <500 | <5 | *** | *** | <50 | <500 | <500 | <5 | <5 | <5 | <5 |
| Ethylbenzene | <200 | <2 | *** | *** | <20 | <200 | <200 | <2 | <2 | <2 | <2 |
| m/p-xylene | <200 | <2 | *** | *** | <20 | <200 | <200 | <2 | <2 | <2 | <2 |
| Methyltert-butylether | <200 | <2 | *** | *** | <20 | <200 | <200 | <2 | <2 | <2 | <2 |
| Naphthalene | <500 | <5 | *** | *** | <50 | <500 | <500 | <5 | <5 | <5 | <5 |
| n-Propylbenzene | <200 | <2 | *** | *** | <20 | <200 | <200 | <2 | <2 | <2 | <2 |
| o-Xylene | <200 | <2 | *** | *** | <20 | <200 | <200 | <2 | <2 | <2 | <2 |
| Tetrachloroethene | <200 | <2 | *** | *** | 19000D | 45000D | 44000D | <2 | 42 | <2 | <2 |
| Tetrahydrofuran | <1000 | <10 | *** | *** | <100 | <1000 | <1000 | <10 | <10 | <10 | <10 |
| Toluene | <200 | <2 | *** | *** | <20 | <200 | <200 | <2 | <2 | <2 | <2 |
| Trichloroethene | 5900 | <2 | *** | *** | 160 | <200 | <200 | <2 | 3.9 | <2 | <2 |
| Trichlorofluoromethane | <200 | <2 | *** | *** | <20 | <200 | <200 | <2 | <2 | <2 | <2 |
| Vinyl chloride | <200 | <2 | *** | *** | 30 | 520 | 460 | <2 | <2 | <2 | <2 |
| Xylene (total) | <200 | <2 | *** | *** | <20 | <200 | <200 | <2 | <2 | <2 | <2 |
| TPH (mg/l) | *** | *** | 8.5 | 8.6 | *** | *** | *** | *** | *** | *** | *** |
| Unidentified TPH | *** | *** | *** | *** | *** | *** | *** | *** | *** | *** | *** |
| Metals 6010B (ug/l) | *** | *** | *** | *** | *** | *** | *** | *** | *** | *** | *** |
| Dissolved Lead | *** | *** | *** | *** | *** | *** | *** | *** | *** | *** | *** |

Notes:
D = Result reported is from a diluted sample
< = Less than the laboratory reporting limit
ug/l = Micro grams per liter, parts per billion
mg/l = Milligrams per liter, parts per million
TPH = Total Petroleum Hydrocarbons
*** = Not analyzed for.

Table 8
Groundwater Analytical Results
Textron/Gorham
Providence, Rhode Island

| Sample Identifier Date Sampled Constituent | MW-201D 2/17/2007 Primary | MW-202D 2/17/2007 Primary | MW-202S 2/17/2007 Primary | MW-207D 2/17/2007 Primary | MW-207S 2/17/2007 Primary | MW-209D 2/17/2007 Primary | MW-216D 2/17/2007 Primary | MW-216S 2/17/2007 Primary | MW-217D 2/17/2007 Primary | MW-217S 2/17/2007 Primary | MW-218D 2/17/2007 Primary | MW-218S 2/17/2007 Primary | MW-220S 3/1/2007 Primary | MW-221S 3/1/2007 Primary | GZA-3 3/1/2007 Primary |
|--|---------------------------------|---------------------------------|---------------------------------|---------------------------------|---------------------------------|---------------------------------|---------------------------------|---------------------------------|---------------------------------|---------------------------------|---------------------------------|---------------------------------|--------------------------------|--------------------------------|------------------------------|
| Method 8260 (ug/l) | | | | | | | | | | | | | | | |
| 1,1,1-Trichloroethane | <200 | <200 | <200 | <200 | <200 | <20 | <2 | <2 | <2 | <2 | <20 | <20 | 210 | 25 | --- |
| 1,1-Dichloroethane | <200 | <200 | <200 | <200 | <200 | <20 | <2 | <2 | <2 | <2 | <20 | <20 | 290 | 140 | --- |
| 1,1-Dichloroethene | <100 | <100 | <100 | <100 | <100 | <10 | <1 | <1 | <1 | <1 | 23 | <10 | 11 | 1.2 | --- |
| 1,2,4-Trimethylbenzene | <200 | <200 | <200 | <200 | <200 | <20 | <2 | 13 | <2 | <2 | <20 | <20 | <20 | 15 | --- |
| 1,3,5-Trimethylbenzene | <200 | <200 | <200 | <200 | <200 | <20 | <2 | 10 | <2 | <2 | <20 | <20 | <20 | 6.1 | --- |
| 4-Isopropyltoluene | <200 | <200 | <200 | <200 | <200 | <20 | <2 | 2.9 | <2 | <2 | <20 | <20 | <20 | <2 | --- |
| Benzene | <100 | <100 | <100 | <100 | <100 | <10 | <1 | <1 | <1 | <1 | <10 | <10 | <10 | 2.6 | --- |
| Bromodichloromethane | <200 | <200 | <200 | <200 | <200 | <20 | <2 | <2 | <2 | <2 | <20 | <20 | <20 | <2 | --- |
| Carbon tetrachloride | <200 | <200 | <200 | <200 | <200 | <20 | <2 | <2 | <2 | <2 | <20 | <20 | <20 | <2 | --- |
| Chloroethane | <500 | <500 | <500 | <500 | <500 | <50 | <5 | <5 | <5 | <5 | <50 | <50 | <50 | 26 | --- |
| Chloroform | <200 | <200 | <200 | <200 | <200 | <20 | <2 | <2 | <2 | <2 | <20 | <20 | <20 | <2 | --- |
| cis-1,2-Dichloroethene | <200 | <200 | <200 | <200 | <200 | <20 | <2 | 140 | 60 | 9.8 | 28 | 650 | 240 | 8.5 | --- |
| Dichlorodifluoromethane | <500 | <500 | <500 | <500 | <500 | <50 | <5 | <5 | <5 | <5 | <50 | <50 | <50 | 5.1 | --- |
| Ethylbenzene | <200 | <200 | <200 | <200 | <200 | <20 | <2 | 2.6 | <2 | <2 | <20 | <20 | <20 | 6.7 | --- |
| m/p-xylene | <200 | <200 | <200 | <200 | <200 | <20 | <2 | 7 | <2 | <2 | <20 | <20 | <20 | 14 | --- |
| Methyltert-butylether | <200 | <200 | <200 | <200 | <200 | <20 | <2 | <2 | <2 | <2 | <20 | <20 | <20 | <2 | --- |
| Naphthalene | <500 | <500 | <500 | <500 | <500 | <50 | <5 | 17 | <5 | <5 | <50 | <50 | <50 | 9 | --- |
| n-Propylbenzene | <200 | <200 | <200 | <200 | <200 | <20 | <2 | <2 | <2 | <2 | <20 | <20 | <20 | 2.4 | --- |
| o-Xylene | <200 | <200 | <200 | <200 | <200 | <20 | <2 | 8.9 | <2 | <2 | <20 | <20 | <20 | 10 | --- |
| Tetrachloroethene | 7600 | 29000 | 51000D | 7700 | 12000 | 430 | <2 | <2 | <2 | 23 | 600 | 370 | <20 | 4.5 | --- |
| Tetrahydrofuran | <1000 | <1000 | <1000 | <1000 | <1000 | <100 | <10 | <10 | <10 | <10 | <100 | <100 | <100 | 29 | --- |
| Toluene | <200 | <200 | <200 | <200 | <200 | <20 | <2 | 3.4 | <2 | <2 | <20 | <20 | <20 | 3.1 | --- |
| Trichloroethene | 970 | <200 | <200 | <200 | <200 | 110 | 6.4 | <2 | 75 | <2 | 840 | 49 | <20 | 12 | --- |
| Trichlorofluoromethane | <200 | <200 | <200 | <200 | <200 | <20 | 3.1 | <2 | <2 | 2.2 | <20 | <20 | <20 | <2 | --- |
| Vinyl chloride | <200 | <200 | <200 | <200 | <200 | <20 | <2 | <2 | <2 | <2 | <20 | <20 | <20 | 22 | --- |
| Xylene (total) | <200 | <200 | <200 | <200 | <200 | <20 | <2 | 16 | <2 | <2 | <20 | <20 | <20 | 24 | --- |
| TPH (mg/l) | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | 35 | --- |
| Unidentified TPH | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- |
| Metals 6010B (ug/l) | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | 12 |
| Dissolved Lead | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- | --- |

Notes:
D = Result reported is from a dilute
< = Less than the laboratory report
ug/l = Micro grams per liter, parts pe
mg/l = Milligrams per liter, parts per
TPH = Total Petroleum Hydrocarbon
--- = Not analyzed for.

Table 9
Compliance Wells Analytical Results
February/March 2007
Former Gorham
Manufacturing Facility
Providence, Rhode Island

| Mashapaug Pond Compliance Wells | | | |
|--|-----------------|------------------|-----------------------------|
| Sample ID | GZA-3 | MW-109D | Compliance |
| Date Collected | 3/1/2007 | 3/1/2007 | Standard¹ |
| CONSTITUENT | | Duplicate | |
| Metals (mg/L) | | | |
| Lead | 0.012 | NA | 0.03 |
| VOCs (ug/L) | | | |
| 1,1,1-Trichloroethane | NA | <2 | 50,000 |
| 1,1-Dichloroethane | NA | <2 | 50,000 |
| 1,1-Dichloroethene | NA | <1 | 50,000 |
| Chloroform | NA | <2 | 10,000 |
| cis-1,2-Dichloroethene | NA | <2 | 50,000 |
| Tetrachloroethene | NA | <2 | 5,000 |
| Trichloroethene | NA | <2 | 20,000 |

| TPH Remediation Area Well | | | |
|----------------------------------|------------------|------------------|-----------------------------|
| Sample ID | CW-6 | CW-6 | Compliance |
| Date Collected | 2/17/2007 | 2/17/2007 | Standard¹ |
| CONSTITUENT | | Duplicate | |
| TPH (mg/L) | 8.5 | 8.6 | 20 |

| Sewer Interceptor Area Wells | | | |
|-------------------------------------|------------------|------------------|-----------------------------|
| Sample ID | CW-1 | CW-2 | Compliance |
| Date Collected | 2/17/2007 | 2/17/2007 | Standard² |
| CONSTITUENT | | | |
| VOCs (ug/L) | | | |
| 1,1-Dichloroethene | 180 | <1 | 23,000 |
| cis-1,2-Dichloroethene | 350 | <2 | 69,000 |
| Trichloroethene | 5,900 | <2 | 87,000 |

| Adelaide Avenue Well | | | | | |
|-----------------------------|------------------|------------------|------------------|------------------|-----------------------------|
| Sample ID | MW-112 | MW-209D | MW-218D | MW-218S | Compliance |
| Date Collected | 2/17/2007 | 2/17/2007 | 2/17/2007 | 2/17/2007 | Standard³ |
| CONSTITUENT | | | | | |
| VOCs (ug/L) | | | | | |
| cis-1,2-Dichloroethene | <2 | <20 | 28 | 650 | 2,400 |
| 1,1-Dichloroethene | <1 | <10 | 23 | <10 | 7 |
| Methyl tert-butyl ether | 17 | <20 | <20 | <20 | 5,000 |
| Tetrachloroethene | 42 | 430 | 600 | 370 | 150 |
| Trichloroethene | 3.9 | 110 | 840 | 49 | 540 |

Notes:

1. These Site specific compliance standards were taken from the approved RAWP dated April 1, 2001 and/or the RIDEM Remediation Regulations.
2. These compliance standards taken from Table 5 - Upper Concentration Limits for GB Groundwater, RIDEM Remediation Regulations.
3. These compliance standards taken from Table 4 - GB Groundwater Objectives, RIDEM Remediation Regulations.

mg/L - milligrams per liter

ug/L - micrograms per liter

< - compound was not detected below the laboratory reporting limit, concentration shown is the reporting limit.

VOCs - volatile organic compounds

TPH - total petroleum hydrocarbons

NA - Indicates that the analysis was not performed.

APPENDIX A

Boring Logs


Drilling Log

Monitoring Well **SB219A/MW219A**

Page: 1 of 2

Project Former Gorham Manufacturing Facility Owner Textron, Inc.
 Location 333 Adelaide Avenue, Providence, RI Proj. No. 101960
 Surface Elev. NA Total Hole Depth 47.5 ft. North _____ East _____
 Top of Casing NA Water Level Initial ▽ 24.0 ft. Static ▽ 24.0 ft. Diameter 2 in.
 Screen: Dia 1.32 in. Length 4 ft. Type/Size Steel/Vertical Slot/0.0015 in.
 Casing: Dia 1.32 in. Length 43 ft. Type Steel
 Fill Material Concrete, Native Rig/Core Vibra Drill/H641
 Drill Co. Pine & Swallow Method Vibratory Drill
 Driller Mike Conlin Log By J. Danieli Date 12/6/06 Permit # NA
 Checked By _____ License No. _____

COMMENTS
 Soil samples collected at 27-31' and 43-47' below surface grade were sent to the laboratory for microcosm study.
 bsg = Below surface grade

| Depth (ft.) | Well Completion | PID (ppm) | Sample ID % Recovery | Blow Count Recovery | Graphic Log | USCS Class. | Description |
|-------------|---|-----------|----------------------|---------------------|-------------|-------------|--|
| | | | | | | | (Color, Texture, Structure) Geologic Descriptions are Based on the USCS. |
| 0 |  | | | | | | Asphalt Hand clear (0-5' below surface grade) |
| 2 | | | | | | | See Boring Log SB-219B for soil description for 0-27 feet below surface grade. |
| 4 | | | | | | | |
| 6 | | | | | | | |
| 8 | | | | | | | |
| 10 | | | | | | | |
| 12 | | | | | | | |
| 14 | | | | | | | |
| 16 | | | | | | | |
| 18 | | | | | | | |
| 20 | | | | | | | |
| 22 | | | | | | | |
| 24 | | | | | | | |

SHAW_COMMERCIAL Rev: 6/12/02 TEXTRON PROVIDENCE.GPJ IT_CORP.GDT 12/20/06

Project Former Gorham Manufacturing Facility Owner Textron, Inc.

Location 333 Adelaide Avenue, Providence, RI Proj. No. 101960

| Depth (ft.) | Well Completion | PID (ppm) | Sample ID % Recovery | Blow Count Recovery | Graphic Log | USCS Class. | Description (Color, Texture, Structure) Geologic Descriptions are Based on the USCS. |
|-------------|-----------------|-----------|----------------------|---------------------|-------------|-------------|--|
| | | | | | | | <i>Continued</i> |
| 26 | | | | | | | |
| 28 | | | | | | SP | Dark grey, wet, medium SAND in nose of sampler (solvent odor noted) 27-31' bsg: Soil sample sent to lab for analysis in acetate sleeve under anaerobic conditions. Acetate sleeve was not cut open for soil screening or logging. |
| 30 | | | | | | | |
| 32 | | 9990 | S-8 90% | | | SP | Grey, wet, loose, medium SAND, layer of dark grey, sand and gravel (solvent odor noted) |
| 34 | | | | | | SW | Brown, medium SAND |
| 36 | | 9999 | S-9 90% | | | SP | Grey, wet, loose, medium SAND, layers of orange-brown, medium sand and grey fine sand with silt (solvent odor noted) (Note: Trace DNAPL detected, <500 ppm, with oil screen soil (Sudan IV) screening kit at 34-39' bsg) |
| 38 | | | | | | | |
| 40 | | | | | | SP | Orange brown to grey, wet, loose, medium SAND |
| 42 | | 583 | S-10 90% | | | ML | Brown to grey, wet, compact SILT |
| 44 | | | | | | SP | Grey, wet, medium SAND |
| 46 | | | S-11 90% | | | SP | Grey, wet, medium SAND in nose of sampler 43-47' bsg: Soil sample sent to lab for analysis in acetate sleeve under anaerobic conditions. Acetate sleeve was not cut open for soil screening or logging. |
| 48 | | | | | | | End of exploration at 47.5 feet below surface grade. Sump (0.5') installed in bottom of well. |
| 50 | | | | | | | |
| 52 | | | | | | | |
| 54 | | | | | | | |
| 56 | | | | | | | |
| 58 | | | | | | | |

SHAW_COMMERCIAL Rev: 6/12/02 TEXTRON PROVIDENCE.GPJ IT_CORP.GDT 12/20/06

Project Former Gorham Manufacturing Facility Owner Textron, Inc.
 Location 333 Adelaide Avenue, Providence, RI Proj. No. 101960
 Surface Elev. NA Total Hole Depth 31.5 ft. North _____ East _____
 Top of Casing NA Water Level Initial NA Static NA Diameter 2 in.
 Screen: Dia 1.32 in. Length 4 ft. Type/Size Steel/Vertical Slot/0.0015 in.
 Casing: Dia 1.32 in. Length 27 ft. Type Steel
 Fill Material Concrete, Native Rig/Core Vibra Drill/H641
 Drill Co. Pine & Swallow Method Vibratory Drill
 Driller Mike Conlin Log By J. Danieli Date 12/6/06 Permit # NA
 Checked By _____ License No. _____

COMMENTS
 No soil samples collected.
 NR = No recovery

| Depth (ft.) | Well Completion | PID (ppm) | Sample ID % Recovery | Blow Count Recovery | Graphic Log | USCS Class. | Description (Color, Texture, Structure) Geologic Descriptions are Based on the USCS. |
|-------------|-----------------|-----------|----------------------|---------------------|-------------|-------------|---|
| 0 | | | | | | | Asphalt |
| 0-5 | | | | | | | Hand clear (0-5' below surface grade) |
| 5 | | 16.3 | S-1 100% | | | SW | Dark brown, dry, loose, coarse to medium SAND, some gravel, black silt layers over light brown, medium sand |
| 10 | | 12.6 | S-2 100% | | | SW | Tan, loose, dry, medium SAND, over dark brown to brown, medium sand, pieces of wood, trace gravel |
| 15 | | 7.9 | S-3 100% | | | SW | Tan, loose, dry, medium SAND, over dark brown to brown, medium sand, trace gravel |
| 20 | | 11.8 | S-4 90% | | | SW | Tan, dry, loose, medium SAND, some gravel, layers of dark brown, slightly dense, sand with gravel |
| 25 | | 14.1 | S-5 100% | | | SW | Brown, dry, loose, medium SAND, some gravel and layers of dark brown sand and gravel, black staining, over grey, medium sand with some gravel |
| 25 | | 198 | S-6 90% | | | SW | Brown to grey, dry, loose, medium SAND, some gravel |
| 25 | | | | | | SM ML | Olive grey, moist to wet, fine SAND, some silt Grey, wet, SILT (solvent odor noted) |
| 30 | | NR | S-7 0% | | | | No recovery |
| 31.5 | | | | | | | End of exploration at 31.5 feet below surface grade. Sump (0.5') installed in bottom of well. |
| 35 | | | | | | | |
| 40 | | | | | | | |

SHAW_COMMERCIAL Rev. 6/12/02 TEXTRON PROVIDENCE.GPJ IT_CORP.GDT 12/20/06

APPENDIX B

Laboratory Analytical Report: Bioaugmentation



17 Princess Rd
Lawrenceville, New Jersey 08648
Tel: 609/895-5370
Fax: 609/895-1858

Limited Chemistry Deliverables

**Prepared for
Gorham Textron**

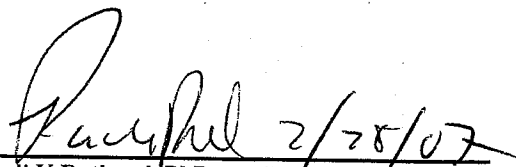
**Lab ID
7543**

Project Number: 101960 02000000

**Samples Received
8-Feb-07**

**Reported
28-Feb-07**

NJDEP Certified Lab 11001


Randi K Rothmel, PhD Date
Laboratory Director

1.0 Chain of Custody

Lab ID
7543

Client:

Date Received:

Shaw E&I Analytical and Treatability Laboratories Internal Chain of Custody

Pg _____ of _____

| Sample ID | Parameter | Bottle Type | Preservative | Date/Time Removed | Relinquishing Custodian Initials | Receiving Analyst Initials | Date/Time Returned | Receiving Custodian Initials | Relinquishing Analyst Initials |
|-----------|-----------|-------------|--------------|-------------------|----------------------------------|----------------------------|--------------------|------------------------------|--------------------------------|
| 7543-1 | VOCs | 1-10-1 | | 02-08-07 9:00 | AM | AZ | 02-08-07 1700 | PM | AZ |
| 2 | | | | | | | | | |
| 3 | | | | | | | | | |
| 4 | | | | | | | | | |
| 5 | | | | | | | | | |
| 6 | | | | | | | | | |
| 7 | | | | | | | | | |
| 8 | | | | | | | | | |
| 9 | | | | | | | | | |
| 10 | | | | | | | | | |
| 11 | | | | | | | | | |
| 7543-12 | VOCs | 1-10-1 | | 02-08-07 9:00 | | AZ | 02-08-07 1700 | | AZ |
| 7543-1 | METALS | 1-10-1 | | 2/12/07 9:00 | PM | PM | 2/12/07 1730 | PM | PM |
| 2 | | | | | | | | | |
| 3 | | | | | | | | | |
| 4 | | | | | | | | | |
| 5 | | | | | | | | | |
| 6 | | | | | | | | | |
| 7 | | | | | | | | | |
| 8 | | | | | | | | | |
| 9 | | | | | | | | | |
| 10 | | | | | | | | | |
| 11 | | | | | | | | | |
| 7543-12 | METALS | 1-10-1 | | | | | | | |
| 7543-13 | ANIONS | 1-10-1 | | 2/2/07 1700 | PM | PM | 2/2/07 1600 | PM | PM |
| 14 | | | | | | | | | |
| 15 | | | | | | | | | |
| 7543-16 | ANIONS | 1-10-1 | | | | | | | |
| 7543-17 | VFA2 | 1-10-1 | | 2/14/07 1200 | PM | PM | 2/14/07 1600 | PM | PM |
| 18 | | | | | | | | | |
| 15 | | | | | | | | | |
| 7543-16 | VFA2 | 1-10-1 | | | | | | | |

2.0 Sample Results

| Sample Information | | | |
|--------------------|------------------|---------------|------------|
| Lab ID | 7543-1 | Date Sampled | 02/08/2007 |
| Sample ID | Killed Control A | Date Received | 02/08/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 02/12/2007 | 8.26 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 02/12/2007 | 2.0 | U | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 02/12/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(2) Not available as a certified parameter under the NJDEP lab certification program.

() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|------------------|---------------|------------|
| Lab ID | 7543-2 | Date Sampled | 02/08/2007 |
| Sample ID | Killed Control B | Date Received | 02/08/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 02/12/2007 | 20.3 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 02/12/2007 | 2.07 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 02/12/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|------------------|---------------|------------|
| Lab ID | 7543-3 | Date Sampled | 02/08/2007 |
| Sample ID | Killed Control C | Date Received | 02/08/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 02/12/2007 | 21.4 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 02/12/2007 | 2.17 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 02/12/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|----------------|---------------|------------|
| Lab ID | 7543-4 | Date Sampled | 02/08/2007 |
| Sample ID | Live Control A | Date Received | 02/08/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 02/12/2007 | 25.6 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 02/12/2007 | 1.47 | J | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 02/12/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|----------------|---------------|------------|
| Lab ID | 7543-5 | Date Sampled | 02/08/2007 |
| Sample ID | Live Control B | Date Received | 02/08/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 02/12/2007 | 25.4 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 02/12/2007 | 1.94 | J | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 02/12/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

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(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|----------------|---------------|------------|
| Lab ID | 7543-6 | Date Sampled | 02/08/2007 |
| Sample ID | Live Control C | Date Received | 02/08/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 02/12/2007 | 23.0 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 02/12/2007 | 2.90 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 02/12/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7543-7 | Date Sampled | 02/08/2007 |
| Sample ID | Bioaug 1 L A | Date Received | 02/08/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 02/12/2007 | 23.8 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 02/12/2007 | 2.02 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 02/12/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7543-8 | Date Sampled | 02/08/2007 |
| Sample ID | Bioaug 1 L B | Date Received | 02/08/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 02/12/2007 | 17.8 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 02/12/2007 | 2.21 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 02/12/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

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(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7543-9 | Date Sampled | 02/08/2007 |
| Sample ID | Bioaug 1 L C | Date Received | 02/08/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 02/12/2007 | 19.2 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 02/12/2007 | 2.30 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 02/12/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7543-10 | Date Sampled | 02/08/2007 |
| Sample ID | Bioaug 2 E A | Date Received | 02/08/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|----------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concent ration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 02/12/2007 | 25.2 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 02/12/2007 | 3.47 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 02/12/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7543-11 | Date Sampled | 02/08/2007 |
| Sample ID | Bioaug 2 E B | Date Received | 02/08/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 02/12/2007 | 23.2 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 02/12/2007 | 4.09 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 02/12/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7543-12 | Date Sampled | 02/08/2007 |
| Sample ID | Bioaug 2 E C | Date Received | 02/08/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|----------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concent ration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 02/12/2007 | 10.2 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 02/12/2007 | 2.0 | U | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 02/12/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|------------------|---------------|------------|
| Lab ID | 7543-13 | Date Sampled | 02/08/2007 |
| Sample ID | Killed Control D | Date Received | 02/08/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-----------------------|---------------|---------------|------------------|-------|------|------|-----------------|-------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Chloride | 02/21/2007 | 73.5 | D | mg/L | 0.5 | 0.07 | 5 | EPA 300.0 |
| Nitrite as N | 02/21/2007 | 0.5 | U | mg/L | 0.5 | 0.08 | 5 | EPA 300.0 |
| Sulfate as SO4 | 02/21/2007 | 42.2 | D | mg/L | 0.5 | 0.09 | 5 | EPA 300.0 |
| Nitrate as N | 02/21/2007 | 0.52 | D | mg/L | 0.5 | 0.05 | 5 | EPA 300.0 |
| Phosphate as P, ortho | 02/21/2007 | 0.5 | U | mg/L | 0.5 | 0.11 | 5 | EPA 300.0 |
| Lactic Acid (2) | 02/14/2007 | 10.0 | U | mg/L | 10.0 | 2.97 | 10 | EPA 300m |
| Acetic Acid (2) | 02/14/2007 | 2.27 | JD | mg/L | 10.0 | 1.05 | 10 | EPA 300m |
| Propionic Acid (2) | 02/14/2007 | 10.0 | U | mg/L | 10.0 | 1.16 | 10 | EPA 300m |
| Formic Acid (2) | 02/14/2007 | 192 | D* | mg/L | 10.0 | 0.81 | 10 | EPA 300m |
| Butyric Acid (2) | 02/14/2007 | 10.0 | U | mg/L | 10.0 | 2.73 | 10 | EPA 300m |
| Pyruvic Acid (2) | 02/14/2007 | 10.0 | U | mg/L | 10.0 | 1.38 | 10 | EPA 300m |
| Valeric Acid (2) | 02/14/2007 | 10.0 | U | mg/L | 10.0 | 1.53 | 10 | EPA 300m |

* probable formaldehyde -- not formic acid -- high level seen only in killed?

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(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|----------------|---------------|------------|
| Lab ID | 7543-14 | Date Sampled | 02/08/2007 |
| Sample ID | Live Control D | Date Received | 02/08/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-----------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|-------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Chloride | 02/21/2007 | 72.7 | D | mg/L | 0.5 | 0.07 | 5 | EPA 300.0 |
| Nitrite as N | 02/21/2007 | 1.3 | D | mg/L | 0.5 | 0.08 | 5 | EPA 300.0 |
| Sulfate as SO4 | 02/21/2007 | 42.8 | D | mg/L | 0.5 | 0.09 | 5 | EPA 300.0 |
| Nitrate as N | 02/21/2007 | 0.5 | U | mg/L | 0.5 | 0.05 | 5 | EPA 300.0 |
| Phosphate as P, ortho | 02/21/2007 | 0.5 | U | mg/L | 0.5 | 0.11 | 5 | EPA 300.0 |
| Lactic Acid (2) | 02/14/2007 | 1.0 | U | mg/L | 1.0 | 0.30 | 1 | EPA 300m |
| Acetic Acid (2) | 02/14/2007 | 0.59 | J | mg/L | 1.0 | 0.11 | 1 | EPA 300m |
| Propionic Acid (2) | 02/14/2007 | 1.0 | U | mg/L | 1.0 | 0.12 | 1 | EPA 300m |
| Formic Acid (2) | 02/14/2007 | 1.0 | U | mg/L | 1.0 | 0.08 | 1 | EPA 300m |
| Butyric Acid (2) | 02/14/2007 | 1.0 | U | mg/L | 1.0 | 0.27 | 1 | EPA 300m |
| Pyruvic Acid (2) | 02/14/2007 | 1.0 | U | mg/L | 1.0 | 0.14 | 1 | EPA 300m |
| Valeric Acid (2) | 02/14/2007 | 1.0 | U | mg/L | 1.0 | 0.15 | 1 | EPA 300m |

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(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7543-15 | Date Sampled | 02/08/2007 |
| Sample ID | Bioaug 1 L D | Date Received | 02/08/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-----------------------|---------------|---------------|---------------------|-------|------|-------|-----------------|-------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Chloride | 02/21/2007 | 72.1 | D | mg/L | 0.5 | 0.07 | 5 | EPA 300.0 |
| Nitrite as N | 02/21/2007 | 0.5 | U | mg/L | 0.5 | 0.08 | 5 | EPA 300.0 |
| Sulfate as SO4 | 02/21/2007 | 46.0 | D | mg/L | 0.5 | 0.09 | 5 | EPA 300.0 |
| Nitrate as N | 02/21/2007 | 0.5 | U | mg/L | 0.5 | 0.05 | 5 | EPA 300.0 |
| Phosphate as P, ortho | 02/21/2007 | 46.5 | D | mg/L | 0.5 | 0.11 | 5 | EPA 300.0 |
| Lactic Acid (2) | 02/14/2007 | 2,740 | D | mg/L | 50.0 | 14.85 | 50 | EPA 300m |
| Acetic Acid (2) | 02/14/2007 | 166 | D | mg/L | 50.0 | 5.25 | 50 | EPA 300m |
| Propionic Acid (2) | 02/14/2007 | 51.6 | D | mg/L | 50.0 | 5.80 | 50 | EPA 300m |
| Formic Acid (2) | 02/14/2007 | 23.1 | JD | mg/L | 50.0 | 4.05 | 50 | EPA 300m |
| Butyric Acid (2) | 02/14/2007 | 50.0 | U | mg/L | 50.0 | 13.65 | 50 | EPA 300m |
| Pyruvic Acid (2) | 02/14/2007 | 50.0 | U | mg/L | 50.0 | 6.90 | 50 | EPA 300m |
| Valeric Acid (2) | 02/14/2007 | 50.0 | U | mg/L | 50.0 | 7.65 | 50 | EPA 300m |

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(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7543-16 | Date Sampled | 02/08/2007 |
| Sample ID | Bioaug 2 E D | Date Received | 02/08/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-----------------------|---------------|---------------|------------------|-------|------|------|-----------------|-------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Chloride | 02/21/2007 | 74.2 | D | mg/L | 0.5 | 0.07 | 5 | EPA 300.0 |
| Nitrite as N | 02/21/2007 | 0.5 | U | mg/L | 0.5 | 0.08 | 5 | EPA 300.0 |
| Sulfate as SO4 | 02/21/2007 | 42.3 | D | mg/L | 0.5 | 0.09 | 5 | EPA 300.0 |
| Nitrate as N | 02/21/2007 | 0.5 | U | mg/L | 0.5 | 0.05 | 5 | EPA 300.0 |
| Phosphate as P, ortho | 02/21/2007 | 44.8 | D | mg/L | 0.5 | 0.11 | 5 | EPA 300.0 |
| Lactic Acid (2) | 02/14/2007 | 10.0 | U | mg/L | 10.0 | 2.97 | 10 | EPA 300m |
| Acetic Acid (2) | 02/14/2007 | 130 | D | mg/L | 10.0 | 1.05 | 10 | EPA 300m |
| Propionic Acid (2) | 02/14/2007 | 82.5 | D | mg/L | 10.0 | 1.16 | 10 | EPA 300m |
| Formic Acid (2) | 02/14/2007 | 16.3 | D | mg/L | 10.0 | 0.81 | 10 | EPA 300m |
| Butyric Acid (2) | 02/14/2007 | 25.8 | D | mg/L | 10.0 | 2.73 | 10 | EPA 300m |
| Pyruvic Acid (2) | 02/14/2007 | 2.71 | JD | mg/L | 10.0 | 1.38 | 10 | EPA 300m |
| Valeric Acid (2) | 02/14/2007 | 10.0 | U | mg/L | 10.0 | 1.53 | 10 | EPA 300m |

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(E) Estimated value beyond linear range



17 Princess Rd
Lawrenceville, New Jersey 08648
Tel: 609/895-5370
Fax: 609/895-1858

Limited Chemistry Deliverables

Prepared for
Gorham Textron


Lab ID
7519

Project Number: 101960 02000000

Samples Received
18-Jan-07

Reported
31-Jan-07

NJDEP Certified Lab 11001

 1/31/07
Randi K Rothmel, PhD Date
Laboratory Director

1.0 Chain of Custody

Lawrenceville, NJ 08648
609-895-5340 / 609-895-1858

CHAIN OF CUSTODY

Ref. Document #

Project Number/Cost code: 101960 / 02

Project Name / Location: Gorham Textron

Send bill to:

Project Contact: Charles Schaefer, Chuck Condee

(Name & phone #)

Send Report To: Chuck Condee

Phone/Fax Number:

Address:

City/State:

Shipment date:

Lab Destination:

Lab Contact Name / ph. #:

Sample's Name(s): Sheryl Streeter

| Lab No. | Sample ID Number | Sample Description |
|---------|------------------|--------------------|
| 1 | Killed A | |
| 2 | Killed B | |
| 3 | Killed C | |
| 4 | Live A | |
| 5 | Live B | |
| 6 | Live C | |
| 7 | Bioaug 1 LA | |
| 8 | Bioaug 1 LB | |
| 9 | Bioaug 1 LC | |
| 10 | Bioaug 2 EA | |
| 11 | Bioaug 2 EB | |
| 12 | Bioaug 2 EC | |
| 13 | Killed D | |
| 14 | Live D | |
| 15 | Bioaug 1 LD | |
| 16 | Bioaug 2 ED | |

Collection Information

| Date | Time | Q/C |
|----------|-------|-----|
| 01/18/07 | 12:00 | |

| Lab No. | Sample ID Number | Sample Description | Matrix | # of Containers | Container Type | Preservative | | | | | VFA | Antons | MER | VOC | Turn Around Time Requested |
|---------|------------------|--------------------|--------|-----------------|----------------|--------------|------|------------------|--------------------------------|------|-----|--------|-----|-----|----------------------------|
| | | | | | | HCL | NaOH | HNO ₃ | H ₂ SO ₄ | None | | | | | |
| 1 | Killed A | | Aq | 1 | 10mL | X | | | | | | | X | | Std |
| 2 | Killed B | | Aq | 1 | 3.75mL | | | | | X | | | X | | Std |
| 3 | Killed C | | Aq | 1 | 10mL | X | | | | | | | X | | Std |
| 4 | Live A | | Aq | 1 | 3.75mL | | | | | X | | | X | | Std |
| 5 | Live B | | Aq | 1 | 10mL | X | | | | | | | X | | Std |
| 6 | Live C | | Aq | 1 | 3.75mL | | | | | X | | | X | | Std |
| 7 | Bioaug 1 LA | | Aq | 1 | 10mL | X | | | | | | | X | | Std |
| 8 | Bioaug 1 LB | | Aq | 1 | 10mL | X | | | | | | | X | | Std |
| 9 | Bioaug 1 LC | | Aq | 1 | 3.75mL | | | | | X | | | X | | Std |
| 10 | Bioaug 2 EA | | Aq | 1 | 3.75mL | | | | | X | | | X | | Std |
| 11 | Bioaug 2 EB | | Aq | 1 | 10mL | X | | | | | | | X | | Std |
| 12 | Bioaug 2 EC | | Aq | 1 | 10mL | X | | | | | | | X | | Std |
| 13 | Killed D | | Aq | 1 | 3.75mL | | | | | X | | | X | | Std |
| 14 | Live D | | Aq | 1 | 15mL | | | | | X | | | X | | Std |
| 15 | Bioaug 1 LD | | Aq | 1 | 15mL | | | | | X | | | X | | Std |
| 16 | Bioaug 2 ED | | Aq | 1 | 15mL | | | | | X | | | X | | Std |

Special Instructions:

MEE bottles hold 5 mL with 3.75 mL sample inside.

QC Codes

C = Composite G = Grab

Known Waste Stream Circles:

RCRA PCB/dioxin PAH/oil

QC/Data Package Level Required:

II III IV NJ EDD GIS EDD Preliminary data

Relinquished By: 

Date: 1-18-07

Time: 10:30

Relinquished By:

Date:

Time:

Received By: 

Date: 1/20/07

Time: 10:30

Received By:

Date:

Time:

Level II = data summary + basic QC
Level III = New Jersey QC reduced deliverable
Level IV = Full deliverable CLP package

Cooler temperature upon arrival at Lab:

2.0 Sample Results

| Sample Information | | | |
|--------------------|------------------|---------------|------------|
| Lab ID | 7519-1 | Date Sampled | 01/18/2007 |
| Sample ID | Killed Control A | Date Received | 01/18/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/29/2007 | 19.6 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/29/2007 | 2.10 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/29/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

Shaw Environmental NJDEP certified Lab ID 11001.

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(2) Not available as a certified parameter under the NJDEP lab certification program.

() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|------------------|---------------|------------|
| Lab ID | 75119-2 | Date Sampled | 01/18/2007 |
| Sample ID | Killed Control B | Date Received | 01/18/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/29/2007 | 20.8 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/29/2007 | 2.13 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/29/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|------------------|---------------|------------|
| Lab ID | 7519-3 | Date Sampled | 01/18/2007 |
| Sample ID | Killed Control C | Date Received | 01/18/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/29/2007 | 22.5 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/29/2007 | 1.96 | J | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/29/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|----------------|---------------|------------|
| Lab ID | 7519-4 | Date Sampled | 01/18/2007 |
| Sample ID | Live Control A | Date Received | 01/18/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/29/2007 | 21.6 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/29/2007 | 2.00 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/29/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|----------------|---------------|------------|
| Lab ID | 7519-5 | Date Sampled | 01/18/2007 |
| Sample ID | Live Control B | Date Received | 01/18/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/29/2007 | 25.0 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/29/2007 | 2.10 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/29/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|----------------|---------------|------------|
| Lab ID | 7519-6 | Date Sampled | 01/18/2007 |
| Sample ID | Live Control C | Date Received | 01/18/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/29/2007 | 22.5 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/29/2007 | 2.14 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/29/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7519-7 | Date Sampled | 01/18/2007 |
| Sample ID | Bioaug 1 L A | Date Received | 01/18/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/29/2007 | 21.9 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/29/2007 | 2.10 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/29/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7519-8 | Date Sampled | 01/18/2007 |
| Sample ID | Bioaug 1 L B | Date Received | 01/18/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/29/2007 | 10.4 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/29/2007 | 2.41 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/29/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

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(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7519-9 | Date Sampled | 01/18/2007 |
| Sample ID | Bioaug 1 L C | Date Received | 01/18/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/29/2007 | 16.1 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/29/2007 | 2.11 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/29/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7519-10 | Date Sampled | 01/18/2007 |
| Sample ID | Bioaug 2 E A | Date Received | 01/18/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/29/2007 | 15.9 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/29/2007 | 4.26 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/29/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7519-11 | Date Sampled | 01/18/2007 |
| Sample ID | Bioaug 2 E B | Date Received | 01/18/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/29/2007 | 18.1 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/29/2007 | 3.81 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/29/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|---------------------------|--------------|---------------|------------|
| Lab ID | 7519-12 | Date Sampled | 01/18/2007 |
| Sample ID | Bioaug 2 E C | Date Received | 01/18/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|--------------------------|----------------------|----------------------|-------------------------|--------------|------------|------------|------------------------|--------------------|
| <i>Parameter</i> | <i>Date Analyzed</i> | <i>Concentration</i> | <i>Qual (see below)</i> | <i>Units</i> | <i>PQL</i> | <i>MDL</i> | <i>Dilution Factor</i> | <i>Method Code</i> |
| Methane (2) | 01/29/2007 | 18.6 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/29/2007 | 5.75 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/29/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|------------------|---------------|------------|
| Lab ID | 7519-13 | Date Sampled | 01/18/2007 |
| Sample ID | Killed Control D | Date Received | 01/18/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-----------------------|---------------|---------------|---------------------|-------|------|------|-----------------|-------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Chloride | 01/19/2007 | 72.1 | D | mg/L | 0.5 | 0.07 | 5 | EPA 300.0 |
| Nitrite as N | 01/19/2007 | 0.5 | U | mg/L | 0.5 | 0.08 | 5 | EPA 300.0 |
| Sulfate as SO4 | 01/19/2007 | 39.9 | D | mg/L | 0.5 | 0.09 | 5 | EPA 300.0 |
| Nitrate as N | 01/19/2007 | 0.72 | D | mg/L | 0.5 | 0.05 | 5 | EPA 300.0 |
| Phosphate as P, ortho | 01/19/2007 | 0.5 | U | mg/L | 0.5 | 0.11 | 5 | EPA 300.0 |
| Lactic Acid (2) | 01/20/2007 | 10.0 | U | mg/L | 10.0 | 2.97 | 10 | EPA 300m |
| Acetic Acid (2) | 01/20/2007 | 2.54 | JD | mg/L | 10.0 | 1.05 | 10 | EPA 300m |
| Propionic Acid (2) | 01/20/2007 | 10.0 | U | mg/L | 10.0 | 1.16 | 10 | EPA 300m |
| Formic Acid (2) | 01/20/2007 | 178 | D* | mg/L | 10.0 | 0.81 | 10 | EPA 300m |
| Butyric Acid (2) | 01/20/2007 | 10.0 | U | mg/L | 10.0 | 2.73 | 10 | EPA 300m |
| Pyruvic Acid (2) | 01/20/2007 | 10.0 | U | mg/L | 10.0 | 1.38 | 10 | EPA 300m |
| Valeric Acid (2) | 01/20/2007 | 10.0 | U | mg/L | 10.0 | 1.53 | 10 | EPA 300m |

* probable formaldehyde -- not formic acid -- high level seen only in killed?

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| Sample Information | | | |
|--------------------|----------------|---------------|------------|
| Lab ID | 7519-14 | Date Sampled | 01/18/2007 |
| Sample ID | Live Control D | Date Received | 01/18/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-----------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|-------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Chloride | 01/19/2007 | 73.5 | D | mg/L | 0.5 | 0.07 | 5 | EPA 300.0 |
| Nitrite as N | 01/19/2007 | 1.3 | D | mg/L | 0.5 | 0.08 | 5 | EPA 300.0 |
| Sulfate as SO4 | 01/19/2007 | 43.5 | D | mg/L | 0.5 | 0.09 | 5 | EPA 300.0 |
| Nitrate as N | 01/19/2007 | 0.11 | JD | mg/L | 0.5 | 0.05 | 5 | EPA 300.0 |
| Phosphate as P, ortho | 01/19/2007 | 0.5 | U | mg/L | 0.5 | 0.11 | 5 | EPA 300.0 |
| Lactic Acid (2) | 01/20/2007 | 1.0 | U | mg/L | 1.0 | 0.30 | 1 | EPA 300m |
| Acetic Acid (2) | 01/20/2007 | 0.41 | J | mg/L | 1.0 | 0.11 | 1 | EPA 300m |
| Propionic Acid (2) | 01/20/2007 | 1.0 | U | mg/L | 1.0 | 0.12 | 1 | EPA 300m |
| Formic Acid (2) | 01/20/2007 | 1.0 | U | mg/L | 1.0 | 0.08 | 1 | EPA 300m |
| Butyric Acid (2) | 01/20/2007 | 1.0 | U | mg/L | 1.0 | 0.27 | 1 | EPA 300m |
| Pyruvic Acid (2) | 01/20/2007 | 1.0 | U | mg/L | 1.0 | 0.14 | 1 | EPA 300m |
| Valeric Acid (2) | 01/20/2007 | 1.0 | U | mg/L | 1.0 | 0.15 | 1 | EPA 300m |

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(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7519-15 | Date Sampled | 01/18/2007 |
| Sample ID | Bioaug 1 L D | Date Received | 01/18/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-----------------------|---------------|---------------|---------------------|-------|------|-------|-----------------|-------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Chloride | 01/19/2007 | 71.8 | D | mg/L | 0.5 | 0.07 | 5 | EPA 300.0 |
| Nitrite as N | 01/19/2007 | 0.5 | U | mg/L | 0.5 | 0.08 | 5 | EPA 300.0 |
| Sulfate as SO4 | 01/19/2007 | 42.3 | D | mg/L | 0.5 | 0.09 | 5 | EPA 300.0 |
| Nitrate as N | 01/19/2007 | 0.5 | U | mg/L | 0.5 | 0.05 | 5 | EPA 300.0 |
| Phosphate as P, ortho | 01/19/2007 | 50.3 | D | mg/L | 0.5 | 0.11 | 5 | EPA 300.0 |
| Lactic Acid (2) | 01/20/2007 | 3,080 | D | mg/L | 50.0 | 14.85 | 50 | EPA 300m |
| Acetic Acid (2) | 01/20/2007 | 139 | D | mg/L | 50.0 | 5.25 | 50 | EPA 300m |
| Propionic Acid (2) | 01/20/2007 | 28.9 | JD | mg/L | 50.0 | 5.80 | 50 | EPA 300m |
| Formic Acid (2) | 01/20/2007 | 16.8 | JD | mg/L | 50.0 | 4.05 | 50 | EPA 300m |
| Butyric Acid (2) | 01/20/2007 | 50.0 | U | mg/L | 50.0 | 13.65 | 50 | EPA 300m |
| Pyruvic Acid (2) | 01/20/2007 | 50.0 | U | mg/L | 50.0 | 6.90 | 50 | EPA 300m |
| Valeric Acid (2) | 01/20/2007 | 50.0 | U | mg/L | 50.0 | 7.65 | 50 | EPA 300m |

Shaw Environmental NJDEP certified Lab ID 11001.

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(2) Not available as a certified parameter under the NJDEP lab certification program.

() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7519-16 | Date Sampled | 01/18/2007 |
| Sample ID | Bioaug 2 E D | Date Received | 01/18/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-----------------------|---------------|---------------|---------------------|-------|------|------|-----------------|-------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Chloride | 01/19/2007 | 74.8 | D | mg/L | 0.5 | 0.07 | 5 | EPA 300.0 |
| Nitrite as N | 01/19/2007 | 0.5 | U | mg/L | 0.5 | 0.08 | 5 | EPA 300.0 |
| Sulfate as SO4 | 01/19/2007 | 45.5 | D | mg/L | 0.5 | 0.09 | 5 | EPA 300.0 |
| Nitrate as N | 01/19/2007 | 0.5 | U | mg/L | 0.5 | 0.05 | 5 | EPA 300.0 |
| Phosphate as P, ortho | 01/19/2007 | 48.4 | D | mg/L | 0.5 | 0.11 | 5 | EPA 300.0 |
| Lactic Acid (2) | 01/20/2007 | 10.0 | U | mg/L | 10.0 | 2.97 | 10 | EPA 300m |
| Acetic Acid (2) | 01/20/2007 | 107 | D | mg/L | 10.0 | 1.05 | 10 | EPA 300m |
| Propionic Acid (2) | 01/20/2007 | 66.8 | D | mg/L | 10.0 | 1.16 | 10 | EPA 300m |
| Formic Acid (2) | 01/20/2007 | 16.9 | D | mg/L | 10.0 | 0.81 | 10 | EPA 300m |
| Butyric Acid (2) | 01/20/2007 | 18.0 | D | mg/L | 10.0 | 2.73 | 10 | EPA 300m |
| Pyruvic Acid (2) | 01/20/2007 | 3.64 | JD | mg/L | 10.0 | 1.38 | 10 | EPA 300m |
| Valeric Acid (2) | 01/20/2007 | 10.0 | U | mg/L | 10.0 | 1.53 | 10 | EPA 300m |

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(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range



17 Princess Rd
Lawrenceville, New Jersey 08648
Tel: 609/895-5370
Fax: 609/895-1858

Limited Chemistry Deliverables

**Prepared for
Gorham Textron**

**Lab ID
7504**

Project Number: 101960 02000000

**Samples Received
3-Jan-07**

**Reported
18-Jan-07**

NJDEP Certified Lab 11001

Randi K Rothmel
Randi K Rothmel, PhD
Laboratory Director

1/18/07
Date

1.0 Chain of Custody

CHAIN OF CUSTODY

Ref. Document #

Page 1 of 1

Shaw Environmental and Infrastructure Inc.

7504

Project Number/Cost code: 101960 / 02

Project Name / Location: Garham Textron

Send bill to:

Project Contact: Charles Schaefer, Chuck Condee
(Name & phone #)

Send Report To: Chuck Condee

Phone/Fax Number:

Address:

City/State:

Shipment date:

Lab Destination:

Lab Contact Name / ph. #:

Sampler's Name(s): Sheryl Stregler

| Lab No. | Sample ID Number | Sample Description | Collection Information | | | # of containers | Container Type | Preservative | | | | MeOH | VOC | MFF | Turn Around Time Requested | |
|---------|------------------|--------------------|------------------------|------|-----|-----------------|----------------|--------------|------|------------------|--------------------------------|------|-----|-----|----------------------------|------|
| | | | Date | Time | G/C | | | HCL | NaOH | HNO ₃ | H ₂ SO ₄ | | | | | None |
| 1 | Killed Control A | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | | | | |
| | " " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | | | | |
| 2 | Killed Control B | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | | | | |
| | " " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | | | | |
| 3 | Killed Control C | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | | | | |
| | " " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | | | | |
| 4 | Live Control A | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | | | | |
| | " " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | | | | |
| 5 | Live Control B | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | | | | |
| | " " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | | | | |
| 6 | Live Control C | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | | | | |
| | " " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | | | | |
| 7 | Bioaug 1 L A | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | | | | |
| | " " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | | | | |
| 8 | Bioaug 1 L B | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | | | | |
| | " " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | | | | |
| 9 | Bioaug 1 L C | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | | | | |
| | " " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | | | | |
| 10 | Bioaug 2 E A | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | | | | |
| | " " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | | | | |
| 11 | Bioaug 2 E B | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | | | | |
| | " " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | | | | |

Special Instructions:

Killed controls contain 1.5% formaldehyde.

Known Waste Stream Circle:

RCRA PCB/dioxin PAH/loil
QC/Data Package Level Required: I II III IV NJ EDD GIS EDD Preliminary data

G/C Codes

C = Composite G = Grab

QC Package Codes

Level II = data summary + basic QC
Level III = New Jersey QC reduced deliverable
Level IV = Full deliverable CLP package

Cooler temperature upon arrival at Lab:

Received By:

Date: 1-3-07
Time: 1008

Sheryl Stregler

Received By:

Date:
Time:

Date: 01/03/07
Time: 10:30

Cooler temperature upon arrival at Lab:

Lab ID 7504-16

Client:

Date Received:

Shaw E&I Analytical and Treatability Laboratories Internal Chain of Custody

Pg _____ of _____

| Sample ID | Parameter | Bottle Type | Preservative | Date/Time Removed | Relinquishing Custodian Initials | Receiving Analyst Initials | Date/Time Returned | Receiving Custodian Initials | Relinquishing Analyst Initials |
|-----------|-----------|-------------|--------------|-------------------|----------------------------------|----------------------------|--------------------|------------------------------|--------------------------------|
| 7504-1 | VOC's | 1-10-1 | HCl | 01-06-07 9:2 | AS | AS | 01-06-07 174 | | AS |
| 2 | | | | | | | | | |
| 3 | | | | | | | | | |
| 4 | | | | | | | | | |
| 5 | | | | | | | | | |
| 6 | | | | | | | | | |
| 7 | | | | | | | | | |
| 8 | | | | | | | | | |
| 9 | | | | | | | | | |
| 10 | | | | | | | | | |
| 11 | | | | | | | | | |
| 7504-12 | VOC's | 1-10-1 | HCl | 01-06-07 9:2 | AS | AS | 01-06-07 174 | | AS |
| 7504-1 | MSEEP | 1-10-1 | None | 4/5/07 8:00 | PM | PM | 4/5/07 1500 | | PM |
| 2 | | | | | | | | | |
| 3 | | | | | | | | | |
| 4 | | | | | | | | | |
| 5 | | | | | | | | | |
| 6 | | | | | | | | | |
| 7 | | | | | | | | | |
| 8 | | | | | | | | | |
| 9 | | | | | | | | | |
| 10 | | | | | | | | | |
| 11 | | | | | | | | | |
| 7504-12 | MSEEP | 1-10-1 | None | | | | | | |
| 7504-13 | VFA's | 10-1 | None | 1/4/07 12:00 | PM | PM | 1/4/07 1800 | | PM |
| 14 | | 10-1 | | | | | | | |
| 15 | | 10-1 | | | | | | | |
| 7504-16 | VFA's | 10-1 | None | | | | | | |

2.0 Sample Results

| Sample Information | | | |
|--------------------|------------------|---------------|------------|
| Lab ID | 7504-1 | Date Sampled | 01/03/2007 |
| Sample ID | Killed Control A | Date Received | 01/03/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|----------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concent ration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/15/2007 | 16.0 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/15/2007 | 6.29 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/15/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|------------------|---------------|------------|
| Lab ID | 7504-2 | Date Sampled | 01/03/2007 |
| Sample ID | Killed Control B | Date Received | 01/03/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/15/2007 | 18.2 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/15/2007 | 7.40 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/15/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|------------------|---------------|------------|
| Lab ID | 7504-3 | Date Sampled | 01/03/2007 |
| Sample ID | Killed Control C | Date Received | 01/03/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|----------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concent ration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/15/2007 | 19.4 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/15/2007 | 7.26 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/15/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

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(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|----------------|---------------|------------|
| Lab ID | 7504-4 | Date Sampled | 01/03/2007 |
| Sample ID | Live Control A | Date Received | 01/03/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|----------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concent ration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/15/2007 | 22.4 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/15/2007 | 7.03 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/15/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range.

| Sample Information | | | |
|--------------------|----------------|---------------|------------|
| Lab ID | 7504-5 | Date Sampled | 01/03/2007 |
| Sample ID | Live Control B | Date Received | 01/03/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/15/2007 | 30.3 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/15/2007 | 8.32 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/15/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

Sample Information

| | | | |
|-----------|----------------|---------------|------------|
| Lab ID | 7504-6 | Date Sampled | 01/03/2007 |
| Sample ID | Live Control C | Date Received | 01/03/2007 |
| Matrix | Aqueous | | |

Limited Chemistry

| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
|-------------|---------------|---------------|------------------|-------|-----|------|-----------------|------------------|
| Methane (2) | 01/15/2007 | 29.5 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/15/2007 | 11.5 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/15/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7504-7 | Date Sampled | 01/03/2007 |
| Sample ID | Bioaug 1 L A | Date Received | 01/03/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/15/2007 | 21.9 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/15/2007 | 8.67 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/15/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7504-8 | Date Sampled | 01/03/2007 |
| Sample ID | Bioaug 1 L B | Date Received | 01/03/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/15/2007 | 25.0 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/15/2007 | 16.1 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/15/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7504-9 | Date Sampled | 01/03/2007 |
| Sample ID | Bioaug 1 L C | Date Received | 01/03/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/15/2007 | 23.4 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/15/2007 | 12.7 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/15/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7504-10 | Date Sampled | 01/03/2007 |
| Sample ID | Bioaug 2 E A | Date Received | 01/03/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/15/2007 | 14.6 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/15/2007 | 10.2 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/15/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7504-11 | Date Sampled | 01/03/2007 |
| Sample ID | Bioaug 2 E B | Date Received | 01/03/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/15/2007 | 14.2 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/15/2007 | 7.51 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/15/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7504-12 | Date Sampled | 01/03/2007 |
| Sample ID | Bioaug 2 E C | Date Received | 01/03/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/15/2007 | 12.3 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/15/2007 | 8.47 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/15/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|------------------|---------------|------------|
| Lab ID | 7504-13 | Date Sampled | 01/03/2007 |
| Sample ID | Killed Control D | Date Received | 01/03/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-----------------------|---------------|---------------|------------------|-------|------|------|-----------------|-------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Chloride | 01/05/2007 | 68.2 | D | mg/L | 1.0 | 0.13 | 10 | EPA 300.0 |
| Nitrite as N | 01/05/2007 | 1.0 | U | mg/L | 1.0 | 0.16 | 10 | EPA 300.0 |
| Sulfate as SO4 | 01/05/2007 | 33.1 | D | mg/L | 1.0 | 0.19 | 10 | EPA 300.0 |
| Nitrate as N | 01/05/2007 | 0.64 | JD | mg/L | 1.0 | 0.10 | 10 | EPA 300.0 |
| Phosphate as P, ortho | 01/05/2007 | 1.0 | U | mg/L | 1.0 | 0.22 | 10 | EPA 300.0 |
| Lactic Acid (2) | 01/04/2007 | 10.0 | U | mg/L | 10.0 | 2.97 | 10 | EPA 300m |
| Acetic Acid (2) | 01/04/2007 | 1.98 | JD | mg/L | 10.0 | 1.05 | 10 | EPA 300m |
| Propionic Acid (2) | 01/04/2007 | 10.0 | U | mg/L | 10.0 | 1.16 | 10 | EPA 300m |
| Formic Acid (2) | 01/04/2007 | 193 | D* | mg/L | 10.0 | 0.81 | 10 | EPA 300m |
| Butyric Acid (2) | 01/04/2007 | 10.0 | U | mg/L | 10.0 | 2.73 | 10 | EPA 300m |
| Pyruvic Acid (2) | 01/04/2007 | 10.0 | U | mg/L | 10.0 | 1.38 | 10 | EPA 300m |
| Valeric Acid (2) | 01/04/2007 | 10.0 | U | mg/L | 10.0 | 1.53 | 10 | EPA 300m |

* probable formaldehyde -- not formic acid -- high level seen only in killed?

Shaw Environmental NJDEP certified Lab ID 11001.

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|----------------|---------------|------------|
| Lab ID | 75042-14 | Date Sampled | 01/03/2007 |
| Sample ID | Live Control D | Date Received | 01/03/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-----------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|-------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Chloride | 01/05/2007 | 69.7 | D | mg/L | 0.5 | 0.07 | 5 | EPA 300.0 |
| Nitrite as N | 01/05/2007 | 0.5 | U | mg/L | 0.5 | 0.08 | 5 | EPA 300.0 |
| Sulfate as SO4 | 01/05/2007 | 39.1 | D | mg/L | 0.5 | 0.09 | 5 | EPA 300.0 |
| Nitrate as N | 01/05/2007 | 0.49 | JD | mg/L | 0.5 | 0.05 | 5 | EPA 300.0 |
| Phosphate as P, ortho | 01/05/2007 | 0.5 | U | mg/L | 0.5 | 0.11 | 5 | EPA 300.0 |
| Lactic Acid (2) | 01/04/2007 | 1.0 | U | mg/L | 1.0 | 0.30 | 1 | EPA 300m |
| Acetic Acid (2) | 01/04/2007 | 1.0 | U | mg/L | 1.0 | 0.11 | 1 | EPA 300m |
| Propionic Acid (2) | 01/04/2007 | 1.0 | U | mg/L | 1.0 | 0.12 | 1 | EPA 300m |
| Formic Acid (2) | 01/04/2007 | 1.0 | U | mg/L | 1.0 | 0.08 | 1 | EPA 300m |
| Butyric Acid (2) | 01/04/2007 | 1.0 | U | mg/L | 1.0 | 0.27 | 1 | EPA 300m |
| Pyruvic Acid (2) | 01/04/2007 | 1.0 | U | mg/L | 1.0 | 0.14 | 1 | EPA 300m |
| Valeric Acid (2) | 01/04/2007 | 1.0 | U | mg/L | 1.0 | 0.15 | 1 | EPA 300m |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7504-15 | Date Sampled | 01/03/2007 |
| Sample ID | Bioaug 1 L D | Date Received | 01/03/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-----------------------|---------------|---------------|------------------|-------|------|-------|-----------------|-------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Chloride | 01/05/2007 | 78.7 | D | mg/L | 0.5 | 0.07 | 5 | EPA 300.0 |
| Nitrite as N | 01/05/2007 | 0.5 | U | mg/L | 0.5 | 0.08 | 5 | EPA 300.0 |
| Sulfate as SO4 | 01/05/2007 | 46.1 | D | mg/L | 0.5 | 0.09 | 5 | EPA 300.0 |
| Nitrate as N | 01/05/2007 | 0.5 | U | mg/L | 0.5 | 0.05 | 5 | EPA 300.0 |
| Phosphate as P, ortho | 01/05/2007 | 41.9 | D | mg/L | 0.5 | 0.11 | 5 | EPA 300.0 |
| Lactic Acid (2) | 01/04/2007 | 2,620 | D | mg/L | 50.0 | 14.85 | 50 | EPA 300m |
| Acetic Acid (2) | 01/04/2007 | 53.2 | D | mg/L | 50.0 | 5.25 | 50 | EPA 300m |
| Propionic Acid (2) | 01/04/2007 | 20.5 | JD | mg/L | 50.0 | 5.80 | 50 | EPA 300m |
| Formic Acid (2) | 01/04/2007 | 50.0 | U | mg/L | 50.0 | 4.05 | 50 | EPA 300m |
| Butyric Acid (2) | 01/04/2007 | 50.0 | U | mg/L | 50.0 | 13.65 | 50 | EPA 300m |
| Pyruvic Acid (2) | 01/04/2007 | 50.0 | U | mg/L | 50.0 | 6.90 | 50 | EPA 300m |
| Valeric Acid (2) | 01/04/2007 | 50.0 | U | mg/L | 50.0 | 7.65 | 50 | EPA 300m |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7504-16 | Date Sampled | 01/03/2007 |
| Sample ID | Bioaug 2 E D | Date Received | 01/03/2007 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-----------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|-------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Chloride | 01/05/2007 | 71.4 | D | mg/L | 0.5 | 0.07 | 5 | EPA 300.0 |
| Nitrite as N | 01/05/2007 | 0.5 | U | mg/L | 0.5 | 0.08 | 5 | EPA 300.0 |
| Sulfate as SO4 | 01/05/2007 | 46.1 | D | mg/L | 0.5 | 0.09 | 5 | EPA 300.0 |
| Nitrate as N | 01/05/2007 | 0.5 | U | mg/L | 0.5 | 0.05 | 5 | EPA 300.0 |
| Phosphate as P, ortho | 01/05/2007 | 41.9 | D | mg/L | 0.5 | 0.11 | 5 | EPA 300.0 |
| Lactic Acid (2) | 01/04/2007 | 1.0 | U | mg/L | 1.0 | 0.30 | 1 | EPA 300m |
| Acetic Acid (2) | 01/04/2007 | 55.4 | | mg/L | 1.0 | 0.11 | 1 | EPA 300m |
| Propionic Acid (2) | 01/04/2007 | 22.5 | | mg/L | 1.0 | 0.12 | 1 | EPA 300m |
| Formic Acid (2) | 01/04/2007 | 14.1 | | mg/L | 1.0 | 0.08 | 1 | EPA 300m |
| Butyric Acid (2) | 01/04/2007 | 1.0 | U | mg/L | 1.0 | 0.27 | 1 | EPA 300m |
| Pyruvic Acid (2) | 01/04/2007 | 1.0 | U | mg/L | 1.0 | 0.14 | 1 | EPA 300m |
| Valeric Acid (2) | 01/04/2007 | 1.0 | U | mg/L | 1.0 | 0.15 | 1 | EPA 300m |

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(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

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(E) Estimated value beyond linear range



17 Princess Rd
Lawrenceville, New Jersey 08648
Tel: 609/895-5370
Fax: 609/895-1858

Limited Chemistry Deliverables

Prepared for
Gorham Textron

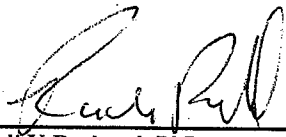
Lab ID
7502

Project Number: 101960 02000000

Samples Received
27-Dec-06

Reported
17-Jan-07

NJDEP Certified Lab 11001


Randi K Rothmel, PhD Date
Laboratory Director

1.0 Chain of Custody

CHAIN OF CUSTODY

Ref. Document #

Project Number/Cost code: 101960 / 02

Project Name / Location: Gorham Textron

Send bill to:

Project Contact: Charles Schaefer, Chuck Condee
(Name & phone #)

Send Report To: Chuck Condee

Phone/Fax Number:

Address:

City/State:

Shipment date:

Lab Destination:

Lab Contact Name / ph. #:

Sampler's Name(s): Sheryl Stregger

Sample ID Number

Sample Description

Collection Information

Date

Time

G/C

g

Matrix

of containers

Container type

HCL

NaOH

HNO₃

H₂SO₄

None

MeOH

Preservative

VOC

MEH

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* Note: PCE ~ 30 PPM!

Known Waste Stream Circle:

PCRA

PCB/dioxin

PAH/oil

QC/Data Package Level Required:

I

II

III

IV

NJ EDD

GIS EDD

Preliminary data

Flammable

Corrosive

Reactive

QC Codes

C = Composite

G = Grab

Relinquished By:

Date: 12-27-06

Time: 1027

Relinquished By:

Date:

Time:

Received By:

Date: 12/27

Time: 1230

Received By:

Date:

Time:

Level II = data summary + basic QC

Level III = New Jersey QC reduced deliverable

Level IV = Full deliverable CLP package

Cooler temperature upon arrival at Lab:



COC Continuation Page

COC Ref. Document #

702

Project Number/ cost code: 101960

02

Shipment Date:

Project Name / Location: Garham Textron

| Lab No. | Sample ID Number | Sample Description | Collection Information | | | # of containers | Container Type | Preservative | | | | | MEE | VOC | Anions | VFA | Turn Around Time Requested |
|---------|------------------|--------------------|------------------------|------|-----|-----------------|----------------|--------------|------|------------------|--------------------------------|------|-----|-----|--------|-----|----------------------------|
| | | | Date | Time | G/C | | | HCL | NaOH | HNO ₃ | H ₂ SO ₄ | None | | | | | |
| 12 | Bioaug 2 E C | " " | 12/27/06 | 7:00 | g | 1 | 10mL | X | | | | | | | | | |
| 13 | Killed Control D | | 12/27/06 | 7:00 | g | 1 | 5mL | | | | | X | | | | | |
| 14 | Live Control D | | 12/27/06 | 7:00 | g | 1 | 15mL | | | | | X | | X | X | | |
| 15 | Bioaug 1 L D | | 12/27/06 | 7:00 | g | 1 | 15mL | | | | | X | | X | X | | |
| 16 | Bioaug 2 E D | | 12/27/06 | 7:00 | g | 1 | 15mL | | | | | X | | X | X | | |

2.0 Sample Results

| Sample Information | | | |
|--------------------|------------------|---------------|------------|
| Lab ID | 7502-1 | Date Sampled | 12/27/2006 |
| Sample ID | Killed Control A | Date Received | 12/27/2006 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|----------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concent ration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/09/2007 | 11.7 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/09/2007 | 7.28 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/09/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|------------------|---------------|------------|
| Lab ID | 7502-2 | Date Sampled | 12/27/2006 |
| Sample ID | Killed Control B | Date Received | 12/27/2006 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|----------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concent ration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/09/2007 | 17.6 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/09/2007 | 8.13 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/09/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|------------------|---------------|------------|
| Lab ID | 7502-3 | Date Sampled | 12/27/2006 |
| Sample ID | Killed Control C | Date Received | 12/27/2006 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/09/2007 | 28.1 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/09/2007 | 7.54 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/09/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(D) Sample analyzed at indicated dilution

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(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|----------------|---------------|------------|
| Lab ID | 7502-4 | Date Sampled | 12/27/2006 |
| Sample ID | Live Control A | Date Received | 12/27/2006 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/09/2007 | 26.8 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/09/2007 | 9.60 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/09/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|----------------|---------------|------------|
| Lab ID | 7502-5 | Date Sampled | 12/27/2006 |
| Sample ID | Live Control B | Date Received | 12/27/2006 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|----------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concent ration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/09/2007 | 17.2 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/09/2007 | 9.02 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/09/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

Sample Information

| | | | |
|-----------|----------------|---------------|------------|
| Lab ID | 7502-6 | Date Sampled | 12/27/2006 |
| Sample ID | Live Control C | Date Received | 12/27/2006 |
| Matrix | Aqueous | | |

Limited Chemistry

| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
|-------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Methane (2) | 01/09/2007 | 25.6 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/09/2007 | 10.2 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/09/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7502-7 | Date Sampled | 12/27/2006 |
| Sample ID | Bioaug 1 L A | Date Received | 12/27/2006 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/09/2007 | 25.8 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/09/2007 | 12.6 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/09/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7502-8 | Date Sampled | 12/27/2006 |
| Sample ID | Bioaug 1 L B | Date Received | 12/27/2006 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/09/2007 | 12.4 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/09/2007 | 9.51 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/09/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7502-9 | Date Sampled | 12/27/2006 |
| Sample ID | Bioaug 1 L C | Date Received | 12/27/2006 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/09/2007 | 11.9 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/09/2007 | 8.85 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/09/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

Sample Information

| | | | |
|-----------|--------------|---------------|------------|
| Lab ID | 7502-10 | Date Sampled | 12/27/2006 |
| Sample ID | Bioaug 2 E A | Date Received | 12/27/2006 |
| Matrix | Aqueous | | |

Limited Chemistry

| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
|-------------|---------------|---------------|------------------|-------|-----|------|-----------------|------------------|
| Methane (2) | 01/09/2007 | 11.6 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/09/2007 | 10.7 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/09/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7502-11 | Date Sampled | 12/27/2006 |
| Sample ID | Bioaug 2 E B | Date Received | 12/27/2006 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/09/2007 | 16.2 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/09/2007 | 10.7 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/09/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

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() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7502-12 | Date Sampled | 12/27/2006 |
| Sample ID | Bioaug 2 E C | Date Received | 12/27/2006 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-------------------|---------------|---------------|------------------|-------|-----|------|-----------------|------------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Methane (2) | 01/09/2007 | 17.1 | | ug/L | 2.0 | 0.98 | 1 | EPA3810, RSK-175 |
| Ethane (2) | 01/09/2007 | 12.4 | | ug/L | 2.0 | 1.00 | 1 | EPA3810, RSK-175 |
| Ethene (2) | 01/09/2007 | 2.0 | U | ug/L | 2.0 | 0.37 | 1 | EPA3810, RSK-175 |

Shaw Environmental NJDEP certified Lab ID 11001.

(1) Not listed as a Shaw Certified parameters under the NJDEP lab certification program.

(2) Not available as a certified parameter under the NJDEP lab certification program.

() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|------------------|---------------|------------|
| Lab ID | 7502-13 | Date Sampled | 12/27/2006 |
| Sample ID | Killed Control D | Date Received | 12/27/2006 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-----------------------|---------------|---------------|------------------|-------|------|------|-----------------|-------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Chloride | 12/28/2006 | 70.7 | D | mg/L | 0.5 | 0.07 | 5 | EPA 300.0 |
| Nitrite as N | 12/28/2006 | 0.5 | U | mg/L | 0.5 | 0.08 | 5 | EPA 300.0 |
| Sulfate as SO4 | 12/28/2006 | 32.4 | D | mg/L | 0.5 | 0.09 | 5 | EPA 300.0 |
| Nitrate as N | 12/28/2006 | 0.59 | D | mg/L | 0.5 | 0.05 | 5 | EPA 300.0 |
| Phosphate as P, ortho | 12/28/2006 | 0.5 | U | mg/L | 0.5 | 0.11 | 5 | EPA 300.0 |
| Lactic Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.30 | 1 | EPA 300m |
| Acetic Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.11 | 1 | EPA 300m |
| Propionic Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.12 | 1 | EPA 300m |
| Formic Acid (2) | 12/28/2006 | 177 | D* | mg/L | 10.0 | 0.81 | 10 | EPA 300m |
| Butyric Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.27 | 1 | EPA 300m |
| Pyruvic Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.14 | 1 | EPA 300m |
| Valeric Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.15 | 1 | EPA 300m |

* probable formaldehyde -- not formic acid -- high level seen only in killed?

Shaw Environmental NJDEP certified Lab ID 11001.

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(2) Not available as a certified parameter under the NJDEP lab certification program.

() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|----------------|---------------|------------|
| Lab ID | 7502-14 | Date Sampled | 12/27/2006 |
| Sample ID | Live Control D | Date Received | 12/27/2006 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-----------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|-------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Chloride | 12/28/2006 | 71.10 | D | mg/L | 0.5 | 0.07 | 5 | EPA 300.0 |
| Nitrite as N | 12/28/2006 | 0.5 | U | mg/L | 0.5 | 0.08 | 5 | EPA 300.0 |
| Sulfate as SO4 | 12/28/2006 | 39.5 | D | mg/L | 0.5 | 0.09 | 5 | EPA 300.0 |
| Nitrate as N | 12/28/2006 | 0.47 | J | mg/L | 0.5 | 0.05 | 5 | EPA 300.0 |
| Phosphate as P, ortho | 12/28/2006 | 0.5 | U | mg/L | 0.5 | 0.11 | 5 | EPA 300.0 |
| Lactic Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.30 | 1 | EPA 300m |
| Acetic Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.11 | 1 | EPA 300m |
| Propionic Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.12 | 1 | EPA 300m |
| Formic Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.08 | 1 | EPA 300m |
| Butyric Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.27 | 1 | EPA 300m |
| Pyruvic Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.14 | 1 | EPA 300m |
| Valeric Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.15 | 1 | EPA 300m |

Shaw Environmental NJDEP certified Lab ID 11001.

(1) Not listed as a Shaw Certified parameters under the NJDEP lab certification program.

(2) Not available as a certified parameter under the NJDEP lab certification program.

() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7502-15 | Date Sampled | 12/27/2006 |
| Sample ID | Bioaug 1 L D | Date Received | 12/27/2006 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-----------------------|---------------|---------------|---------------------|-------|------|-------|-----------------|-------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Chloride | 12/28/2006 | 71.5 | D | mg/L | 0.5 | 0.07 | 5 | EPA 300.0 |
| Nitrite as N | 12/28/2006 | 0.5 | U | mg/L | 0.5 | 0.08 | 5 | EPA 300.0 |
| Sulfate as SO4 | 12/28/2006 | 44.2 | D | mg/L | 0.5 | 0.09 | 5 | EPA 300.0 |
| Nitrate as N | 12/28/2006 | 0.5 | U | mg/L | 0.5 | 0.05 | 5 | EPA 300.0 |
| Phosphate as P, ortho | 12/28/2006 | 40.2 | D | mg/L | 0.5 | 0.11 | 5 | EPA 300.0 |
| Lactic Acid (2) | 12/28/2006 | 1,370 | D | mg/L | 50.0 | 14.85 | 50 | EPA 300m |
| Acetic Acid (2) | 12/28/2006 | 51.5 | D | mg/L | 10.0 | 1.05 | 10 | EPA 300m |
| Propionic Acid (2) | 12/28/2006 | 22.5 | D | mg/L | 10.0 | 1.16 | 10 | EPA 300m |
| Formic Acid (2) | 12/28/2006 | 4.50 | | mg/L | 1.0 | 0.08 | 1 | EPA 300m |
| Butyric Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.27 | 1 | EPA 300m |
| Pyruvic Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.14 | 1 | EPA 300m |
| Valeric Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.15 | 1 | EPA 300m |

Shaw Environmental NJDEP certified Lab ID 11001.

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(2) Not available as a certified parameter under the NJDEP lab certification program.

() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range

| Sample Information | | | |
|--------------------|--------------|---------------|------------|
| Lab ID | 7502-16 | Date Sampled | 12/27/2006 |
| Sample ID | Bioaug 2 E D | Date Received | 12/27/2006 |
| Matrix | Aqueous | | |

| Limited Chemistry | | | | | | | | |
|-----------------------|---------------|---------------|---------------------|-------|-----|------|-----------------|-------------|
| Parameter | Date Analyzed | Concentration | Qual (see below) | Units | PQL | MDL | Dilution Factor | Method Code |
| Chloride | 12/28/2006 | 72.8 | D | mg/L | 0.5 | 0.07 | 5 | EPA 300.0 |
| Nitrite as N | 12/28/2006 | 0.5 | U | mg/L | 0.5 | 0.08 | 5 | EPA 300.0 |
| Sulfate as SO4 | 12/28/2006 | 45.5 | D | mg/L | 0.5 | 0.09 | 5 | EPA 300.0 |
| Nitrate as N | 12/28/2006 | 0.5 | U | mg/L | 0.5 | 0.05 | 5 | EPA 300.0 |
| Phosphate as P, ortho | 12/28/2006 | 40.6 | D | mg/L | 0.5 | 0.11 | 5 | EPA 300.0 |
| Lactic Acid (2) | 12/28/2006 | 31.0 | | mg/L | 1.0 | 0.30 | 1 | EPA 300m |
| Acetic Acid (2) | 12/28/2006 | 33.4 | | mg/L | 1.0 | 0.11 | 1 | EPA 300m |
| Propionic Acid (2) | 12/28/2006 | 24.1 | | mg/L | 1.0 | 0.12 | 1 | EPA 300m |
| Formic Acid (2) | 12/28/2006 | 8.01 | | mg/L | 1.0 | 0.08 | 1 | EPA 300m |
| Butyric Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.27 | 1 | EPA 300m |
| Pyruvic Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.14 | 1 | EPA 300m |
| Valeric Acid (2) | 12/28/2006 | 1.0 | U | mg/L | 1.0 | 0.15 | 1 | EPA 300m |

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(2) Not available as a certified parameter under the NJDEP lab certification program.

() no qualification - sample run undiluted

(U) Compound not detected above method practical quantitation limit.

(D) Sample analyzed at indicated dilution

(J) Estimated value above MDL and less than PQL

(E) Estimated value beyond linear range



17 Princess Rd
Lawrenceville, New Jersey 08648
Tel: 609/895-5370
Fax: 609/895-1858

Volatile Organic Compound Data Summary Package

Prepared for
Gorham Textron

Lab ID
7489

Project Number 101960 02000000

Samples Received
19-Dec-06

Reported
3-Jan-07

NJDEP Certified Lab 11001

A handwritten signature in black ink, appearing to read "Randi K Rothmel", is written over a horizontal line.

Randi K Rothmel, PhD

Laboratory Director

A handwritten date "1/3/07" is written in black ink to the right of the signature line.

Date

1.0 Chain of Custody

17 Princess Road

Lawrenceville, NJ 08648

609-895-5340 / 609-895-1858



Shaw Environmental and Infrastructure Inc.

7489

Project Contact: Charles Schaefer, Sheryl Strager
(Name & phone #)

Send Report To: Sheryl Strager

Phone/Fax Number:

Address:

City/State:

CHAIN OF CUSTODY

Ref. Document #

Project Number/Cost code: 101960 / 02

Project Name / Location: Gorham Texton

Send bill to:

Shipment date:

Lab Destination:

Lab Contact Name / ph. #:

Analyses Requested

Sampler's Name(s): Corinne Topoleski

| Lab No. | Sample ID Number | Sample Description |
|---------|------------------|--------------------|
| 1 | Killed Control A | |
| 2 | Killed Control B | |
| 3 | Killed Control C | |
| 4 | Live Control A | |
| 5 | Live Control B | |
| 6 | Live Control C | |
| 7 | Bioaug 1L A | |
| 8 | Bioaug 1L B | |
| 9 | Bioaug 1L C | |
| 10 | Bioaug 2E A | |
| 11 | Bioaug 2E B | |
| 12 | Bioaug 2E C | |

| Collection Information | Preservative | | | | | |
|------------------------|--------------|----------|------------------|--------------------------------|----------|----------|
| | HCL | NaOH | HNO ₃ | H ₂ SO ₄ | None | MeOH |
| Date | 12/19/06 | 12/19/06 | 12/19/06 | 12/19/06 | 12/19/06 | 12/19/06 |
| Time | 0700 | | | | | |
| G/C | G | G | G | G | G | G |
| Matrix | Aq | Aq | Aq | Aq | Aq | Aq |
| # of Containers | 1 | 1 | 1 | 1 | 1 | 1 |
| Container Type | 8 mL | 8 mL | 8 mL | 8 mL | 8 mL | 8 mL |

VOC 8260

RAPID RAPID RAPID RAPID RAPID RAPID RAPID RAPID RAPID RAPID RAPID

Special Instructions:

24 hour turnaround

Known Waste Stream Circle:

FCRA PCB/dioxin PAH/oil
QC/Data Package Level Required: I III III IV NJ EDD GIS EDD Preliminary data

G/C Codes
C = Composite
G = Grab

QC Package Codes

Relinquished By: Sheryl Strager

Date: 12.19.06
Time: 0753

Received By: [Signature]

Date: 12-19-06
Time: 8:25a

Relinquished By:

Date: _____
Time: _____

Received By:

Date: _____
Time: _____

Level II = data summary + basic QC
Level III = New Jersey QC reduced deliverable
Level IV = Full deliverable CLP package
Cooler temperature upon arrival at Lab:

Lab ID 7489-12

Client: _____

Date Received: _____

Shaw E&I Analytical and Treatability Laboratories Internal Chain of Custody

Pg _____ of _____

| Sample ID | Parameter | Bottle Type | Perservative | Date/Time Removed | Relinquishing Custodian Initials | Receiving Analyst Initials | Date/Time Returned | Receiving Custodian Initials | Relinquishing Analyst Initials |
|-----------|-----------|-------------|--------------|-------------------|----------------------------------|----------------------------|--------------------|------------------------------|--------------------------------|
| 7489-1 | VOC's | 1-10-1 | HCL | 12-20-06 9:50 | AS | AS | 12-20-06 17L | AS | AS |
| 2 | | | | | | | | | |
| 3 | | | | | | | | | |
| 4 | | | | | | | | | |
| 5 | | | | | | | | | |
| 6 | | | | | | | | | |
| 7 | | | | | | | | | |
| 8 | | | | | | | | | |
| 9 | | | | | | | | | |
| 10 | | | | | | | | | |
| 11 | | | | | | | | | |
| 7489-12 | VOC's | 1-10-1 | HCL | 12-20-06 9:50 | AS | AS | 12-20-06 17L | AS | AS |

2.0. Sample Summaries

Sample summaries are enclosed

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-1

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_1 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 9_1.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/20/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | CONCENTRATION UNITS: | |
|------------|---------------------------|----------------------|------|
| | | (ug/L or ug/Kg) | UG/L |
| 75-71-8 | Dichlorodifluoromethane | 520 | U |
| 74-87-3 | Chloromethane | 520 | U |
| 75-01-4 | Vinyl chloride | 520 | U |
| 74-83-9 | Bromomethane | 520 | U |
| 75-00-3 | Chloroethane | 520 | U |
| 75-69-4 | Trichlorofluoromethane | 520 | U |
| 75-35-4 | 1,1-Dichloroethene | 520 | U |
| 75-09-2 | Methylene chloride | 520 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 520 | U |
| 75-34-3 | 1,1-Dichloroethane | 520 | U |
| 594-20-7 | 2,2-Dichloropropane | 520 | U |
| 156-59-2 | cis-1,2-Dichloroethene | 520 | U |
| 74-97-5 | Bromochloromethane | 520 | U |
| 67-66-3 | Chloroform | 520 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 520 | U |
| 56-23-5 | Carbon tetrachloride | 520 | U |
| 563-58-6 | 1,1-Dichloropropene | 520 | U |
| 71-43-2 | Benzene | 520 | U |
| 107-06-2 | 1,2-Dichloroethane | 520 | U |
| 79-01-6 | Trichloroethene | 520 | U |
| 78-87-5 | 1,2-Dichloropropane | 520 | U |
| 74-95-3 | Dibromomethane | 520 | U |
| 75-27-4 | Bromodichloromethane | 520 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | 520 | U |
| 108-88-3 | Toluene | 520 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | 520 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 520 | U |
| 127-18-4 | Tetrachloroethene | 6600 | D |
| 142-28-9 | 1,3-Dichloropropane | 520 | U |
| 124-48-1 | Dibromochloromethane | 520 | U |
| 106-93-4 | 1,2-Dibromoethane | 520 | U |
| 108-90-7 | Chlorobenzene | 520 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 520 | U |
| 100-41-4 | Ethylbenzene | 520 | U |
| 1330-20-7 | Xylene (para & meta) | 520 | U |
| 95-47-6 | Xylene (Ortho) | 520 | U |
| 100-42-5 | Styrene | 520 | U |
| 75-25-2 | Bromoform | 520 | U |
| 98-82-8 | Isopropylbenzene | 520 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-1

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_1 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 9_1.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/20/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 520 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 520 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 520 | U |
| 103-65-1 | n-Propylbenzene | | 520 | U |
| 95-49-8 | 2-Chlorotoluene | | 520 | U |
| 106-43-4 | 4-Chlorotoluene | | 520 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 520 | U |
| 98-06-6 | tert-Butylbenzene | | 520 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 520 | U |
| 135-98-8 | sec-Butylbenzene | | 520 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 520 | U |
| 99-87-6 | 4-Isopropyltoluene | | 520 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 520 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 520 | U |
| 104-51-8 | n-Butylbenzene | | 520 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 520 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 520 | U |
| 87-68-3 | Hexachlorobutadiene | | 520 | U |
| 91-20-3 | Naphthalene | | 520 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 520 | U |
| 1634-04-4 | MTBE | | 520 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 520 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-2

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_2 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 9_2.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/20/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | | 520 | U |
| 74-87-3 | Chloromethane | | 520 | U |
| 75-01-4 | Vinyl chloride | | 520 | U |
| 74-83-9 | Bromomethane | | 520 | U |
| 75-00-3 | Chloroethane | | 520 | U |
| 75-69-4 | Trichlorofluoromethane | | 520 | U |
| 75-35-4 | 1,1-Dichloroethene | | 520 | U |
| 75-09-2 | Methylene chloride | | 520 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 520 | U |
| 75-34-3 | 1,1-Dichloroethane | | 520 | U |
| 594-20-7 | 2,2-Dichloropropane | | 520 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 520 | U |
| 74-97-5 | Bromochloromethane | | 520 | U |
| 67-66-3 | Chloroform | | 520 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 520 | U |
| 56-23-5 | Carbon tetrachloride | | 520 | U |
| 563-58-6 | 1,1-Dichloropropene | | 520 | U |
| 71-43-2 | Benzene | | 520 | U |
| 107-06-2 | 1,2-Dichloroethane | | 520 | U |
| 79-01-6 | Trichloroethene | | 520 | U |
| 78-87-5 | 1,2-Dichloropropane | | 520 | U |
| 74-95-3 | Dibromomethane | | 520 | U |
| 75-27-4 | Bromodichloromethane | | 520 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 520 | U |
| 108-88-3 | Toluene | | 520 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 520 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 520 | U |
| 127-18-4 | Tetrachloroethene | | 4500 | D |
| 142-28-9 | 1,3-Dichloropropane | | 520 | U |
| 124-48-1 | Dibromochloromethane | | 520 | U |
| 106-93-4 | 1,2-Dibromoethane | | 520 | U |
| 108-90-7 | Chlorobenzene | | 520 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 520 | U |
| 100-41-4 | Ethylbenzene | | 520 | U |
| 1330-20-7 | Xylene (para & meta) | | 520 | U |
| 95-47-6 | Xylene (Ortho) | | 520 | U |
| 100-42-5 | Styrene | | 520 | U |
| 75-25-2 | Bromoform | | 520 | U |
| 98-82-8 | Isopropylbenzene | | 520 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-2

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_2 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 9_2.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/20/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 520 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 520 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 520 | U |
| 103-65-1 | n-Propylbenzene | | 520 | U |
| 95-49-8 | 2-Chlorotoluene | | 520 | U |
| 106-43-4 | 4-Chlorotoluene | | 520 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 520 | U |
| 98-06-6 | tert-Butylbenzene | | 520 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 520 | U |
| 135-98-8 | sec-Butylbenzene | | 520 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 520 | U |
| 99-87-6 | 4-Isopropyltoluene | | 520 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 520 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 520 | U |
| 104-51-8 | n-Butylbenzene | | 520 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 520 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 520 | U |
| 87-68-3 | Hexachlorobutadiene | | 520 | U |
| 91-20-3 | Naphthalene | | 520 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 520 | U |
| 1634-04-4 | MTBE | | 520 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 520 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-03

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_03 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_3.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/19/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | | 520 | U |
| 74-87-3 | Chloromethane | | 520 | U |
| 75-01-4 | Vinyl chloride | | 520 | U |
| 74-83-9 | Bromomethane | | 520 | U |
| 75-00-3 | Chloroethane | | 520 | U |
| 75-69-4 | Trichlorofluoromethane | | 520 | U |
| 75-35-4 | 1,1-Dichloroethene | | 520 | U |
| 75-09-2 | Methylene chloride | | 520 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 520 | U |
| 75-34-3 | 1,1-Dichloroethane | | 520 | U |
| 594-20-7 | 2,2-Dichloropropane | | 520 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 520 | U |
| 74-97-5 | Bromochloromethane | | 520 | U |
| 67-66-3 | Chloroform | | 520 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 520 | U |
| 56-23-5 | Carbon tetrachloride | | 520 | U |
| 563-58-6 | 1,1-Dichloropropene | | 520 | U |
| 71-43-2 | Benzene | | 520 | U |
| 107-06-2 | 1,2-Dichloroethane | | 520 | U |
| 79-01-6 | Trichloroethene | | 520 | U |
| 78-87-5 | 1,2-Dichloropropane | | 520 | U |
| 74-95-3 | Dibromomethane | | 520 | U |
| 75-27-4 | Bromodichloromethane | | 520 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 520 | U |
| 108-88-3 | Toluene | | 520 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 520 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 520 | U |
| 127-18-4 | Tetrachloroethene | | 2600 | D |
| 142-28-9 | 1,3-Dichloropropane | | 520 | U |
| 124-48-1 | Dibromochloromethane | | 520 | U |
| 106-93-4 | 1,2-Dibromoethane | | 520 | U |
| 108-90-7 | Chlorobenzene | | 520 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 520 | U |
| 100-41-4 | Ethylbenzene | | 520 | U |
| 1330-20-7 | Xylene (para & meta) | | 520 | U |
| 95-47-6 | Xylene (Ortho) | | 520 | U |
| 100-42-5 | Styrene | | 520 | U |
| 75-25-2 | Bromoform | | 520 | U |
| 98-82-8 | Isopropylbenzene | | 520 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-03

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_03 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_3.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/19/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 520 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 520 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 520 | U |
| 103-65-1 | n-Propylbenzene | | 520 | U |
| 95-49-8 | 2-Chlorotoluene | | 520 | U |
| 106-43-4 | 4-Chlorotoluene | | 520 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 520 | U |
| 98-06-6 | tert-Butylbenzene | | 520 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 520 | U |
| 135-98-8 | sec-Butylbenzene | | 520 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 520 | U |
| 99-87-6 | 4-Isopropyltoluene | | 520 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 520 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 520 | U |
| 104-51-8 | n-Butylbenzene | | 520 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 520 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 520 | U |
| 87-68-3 | Hexachlorobutadiene | | 520 | U |
| 91-20-3 | Naphthalene | | 520 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 520 | U |
| 1634-04-4 | MTBE | | 520 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 520 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-04

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_04 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_4.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/19/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | | 520 | U |
| 74-87-3 | Chloromethane | | 520 | U |
| 75-01-4 | Vinyl chloride | | 520 | U |
| 74-83-9 | Bromomethane | | 520 | U |
| 75-00-3 | Chloroethane | | 520 | U |
| 75-69-4 | Trichlorofluoromethane | | 520 | U |
| 75-35-4 | 1,1-Dichloroethene | | 520 | U |
| 75-09-2 | Methylene chloride | | 520 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 520 | U |
| 75-34-3 | 1,1-Dichloroethane | | 520 | U |
| 594-20-7 | 2,2-Dichloropropane | | 520 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 520 | U |
| 74-97-5 | Bromochloromethane | | 520 | U |
| 67-66-3 | Chloroform | | 520 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 520 | U |
| 56-23-5 | Carbon tetrachloride | | 520 | U |
| 563-58-6 | 1,1-Dichloropropene | | 520 | U |
| 71-43-2 | Benzene | | 520 | U |
| 107-06-2 | 1,2-Dichloroethane | | 520 | U |
| 79-01-6 | Trichloroethene | | 520 | U |
| 78-87-5 | 1,2-Dichloropropane | | 520 | U |
| 74-95-3 | Dibromomethane | | 520 | U |
| 75-27-4 | Bromodichloromethane | | 520 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 520 | U |
| 108-88-3 | Toluene | | 520 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 520 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 520 | U |
| 127-18-4 | Tetrachloroethene | | 4400 | D |
| 142-28-9 | 1,3-Dichloropropane | | 520 | U |
| 124-48-1 | Dibromochloromethane | | 520 | U |
| 106-93-4 | 1,2-Dibromoethane | | 520 | U |
| 108-90-7 | Chlorobenzene | | 520 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 520 | U |
| 100-41-4 | Ethylbenzene | | 520 | U |
| 1330-20-7 | Xylene (para & meta) | | 520 | U |
| 95-47-6 | Xylene (Ortho) | | 520 | U |
| 100-42-5 | Styrene | | 520 | U |
| 75-25-2 | Bromoform | | 520 | U |
| 98-82-8 | Isopropylbenzene | | 520 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-04

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_04 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_4.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/19/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 520 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 520 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 520 | U |
| 103-65-1 | n-Propylbenzene | | 520 | U |
| 95-49-8 | 2-Chlorotoluene | | 520 | U |
| 106-43-4 | 4-Chlorotoluene | | 520 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 520 | U |
| 98-06-6 | tert-Butylbenzene | | 520 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 520 | U |
| 135-98-8 | sec-Butylbenzene | | 520 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 520 | U |
| 99-87-6 | 4-Isopropyltoluene | | 520 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 520 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 520 | U |
| 104-51-8 | n-Butylbenzene | | 520 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 520 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 520 | U |
| 87-68-3 | Hexachlorobutadiene | | 520 | U |
| 91-20-3 | Naphthalene | | 520 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 520 | U |
| 1634-04-4 | MTBE | | 520 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 520 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-05

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_05 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_5.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/19/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | | 520 | U |
| 74-87-3 | Chloromethane | | 520 | U |
| 75-01-4 | Vinyl chloride | | 520 | U |
| 74-83-9 | Bromomethane | | 520 | U |
| 75-00-3 | Chloroethane | | 520 | U |
| 75-69-4 | Trichlorofluoromethane | | 520 | U |
| 75-35-4 | 1,1-Dichloroethene | | 520 | U |
| 75-09-2 | Methylene chloride | | 520 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 520 | U |
| 75-34-3 | 1,1-Dichloroethane | | 520 | U |
| 594-20-7 | 2,2-Dichloropropane | | 520 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 520 | U |
| 74-97-5 | Bromochloromethane | | 520 | U |
| 67-66-3 | Chloroform | | 520 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 520 | U |
| 56-23-5 | Carbon tetrachloride | | 520 | U |
| 563-58-6 | 1,1-Dichloropropene | | 520 | U |
| 71-43-2 | Benzene | | 520 | U |
| 107-06-2 | 1,2-Dichloroethane | | 520 | U |
| 79-01-6 | Trichloroethene | | 520 | U |
| 78-87-5 | 1,2-Dichloropropane | | 520 | U |
| 74-95-3 | Dibromomethane | | 520 | U |
| 75-27-4 | Bromodichloromethane | | 520 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 520 | U |
| 108-88-3 | Toluene | | 520 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 520 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 520 | U |
| 127-18-4 | Tetrachloroethene | | 2800 | D |
| 142-28-9 | 1,3-Dichloropropane | | 520 | U |
| 124-48-1 | Dibromochloromethane | | 520 | U |
| 106-93-4 | 1,2-Dibromoethane | | 520 | U |
| 108-90-7 | Chlorobenzene | | 520 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 520 | U |
| 100-41-4 | Ethylbenzene | | 520 | U |
| 1330-20-7 | Xylene (para & meta) | | 520 | U |
| 95-47-6 | Xylene (Ortho) | | 520 | U |
| 100-42-5 | Styrene | | 520 | U |
| 75-25-2 | Bromoform | | 520 | U |
| 98-82-8 | Isopropylbenzene | | 520 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-05

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_05 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_5.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/19/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 520 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 520 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 520 | U |
| 103-65-1 | n-Propylbenzene | | 520 | U |
| 95-49-8 | 2-Chlorotoluene | | 520 | U |
| 106-43-4 | 4-Chlorotoluene | | 520 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 520 | U |
| 98-06-6 | tert-Butylbenzene | | 520 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 520 | U |
| 135-98-8 | sec-Butylbenzene | | 520 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 520 | U |
| 99-87-6 | 4-Isopropyltoluene | | 520 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 520 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 520 | U |
| 104-51-8 | n-Butylbenzene | | 520 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 520 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 520 | U |
| 87-68-3 | Hexachlorobutadiene | | 520 | U |
| 91-20-3 | Naphthalene | | 520 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 520 | U |
| 1634-04-4 | MTBE | | 520 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 520 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-06

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_06 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_6.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/19/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | | 520 | U |
| 74-87-3 | Chloromethane | | 520 | U |
| 75-01-4 | Vinyl chloride | | 520 | U |
| 74-83-9 | Bromomethane | | 520 | U |
| 75-00-3 | Chloroethane | | 520 | U |
| 75-69-4 | Trichlorofluoromethane | | 520 | U |
| 75-35-4 | 1,1-Dichloroethene | | 520 | U |
| 75-09-2 | Methylene chloride | | 520 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 520 | U |
| 75-34-3 | 1,1-Dichloroethane | | 520 | U |
| 594-20-7 | 2,2-Dichloropropane | | 520 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 520 | U |
| 74-97-5 | Bromochloromethane | | 520 | U |
| 67-66-3 | Chloroform | | 520 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 520 | U |
| 56-23-5 | Carbon tetrachloride | | 520 | U |
| 563-58-6 | 1,1-Dichloropropene | | 520 | U |
| 71-43-2 | Benzene | | 520 | U |
| 107-06-2 | 1,2-Dichloroethane | | 520 | U |
| 79-01-6 | Trichloroethene | | 520 | U |
| 78-87-5 | 1,2-Dichloropropane | | 520 | U |
| 74-95-3 | Dibromomethane | | 520 | U |
| 75-27-4 | Bromodichloromethane | | 520 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 520 | U |
| 108-88-3 | Toluene | | 520 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 520 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 520 | U |
| 127-18-4 | Tetrachloroethene | | 3900 | D |
| 142-28-9 | 1,3-Dichloropropane | | 520 | U |
| 124-48-1 | Dibromochloromethane | | 520 | U |
| 106-93-4 | 1,2-Dibromoethane | | 520 | U |
| 108-90-7 | Chlorobenzene | | 520 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 520 | U |
| 100-41-4 | Ethylbenzene | | 520 | U |
| 1330-20-7 | Xylene (para & meta) | | 520 | U |
| 95-47-6 | Xylene (Ortho) | | 520 | U |
| 100-42-5 | Styrene | | 520 | U |
| 75-25-2 | Bromoform | | 520 | U |
| 98-82-8 | Isopropylbenzene | | 520 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-06

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_06 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_6.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/19/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 520 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 520 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 520 | U |
| 103-65-1 | n-Propylbenzene | | 520 | U |
| 95-49-8 | 2-Chlorotoluene | | 520 | U |
| 106-43-4 | 4-Chlorotoluene | | 520 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 520 | U |
| 98-06-6 | tert-Butylbenzene | | 520 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 520 | U |
| 135-98-8 | sec-Butylbenzene | | 520 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 520 | U |
| 99-87-6 | 4-Isopropyltoluene | | 520 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 520 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 520 | U |
| 104-51-8 | n-Butylbenzene | | 520 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 520 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 520 | U |
| 87-68-3 | Hexachlorobutadiene | | 520 | U |
| 91-20-3 | Naphthalene | | 520 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 520 | U |
| 1634-04-4 | MTBE | | 520 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 520 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-07

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_07 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_7.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/19/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | | 520 | U |
| 74-87-3 | Chloromethane | | 520 | U |
| 75-01-4 | Vinyl chloride | | 520 | U |
| 74-83-9 | Bromomethane | | 520 | U |
| 75-00-3 | Chloroethane | | 520 | U |
| 75-69-4 | Trichlorofluoromethane | | 520 | U |
| 75-35-4 | 1,1-Dichloroethene | | 520 | U |
| 75-09-2 | Methylene chloride | | 520 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 520 | U |
| 75-34-3 | 1,1-Dichloroethane | | 520 | U |
| 594-20-7 | 2,2-Dichloropropane | | 520 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 520 | U |
| 74-97-5 | Bromochloromethane | | 520 | U |
| 67-66-3 | Chloroform | | 520 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 520 | U |
| 56-23-5 | Carbon tetrachloride | | 520 | U |
| 563-58-6 | 1,1-Dichloropropene | | 520 | U |
| 71-43-2 | Benzene | | 520 | U |
| 107-06-2 | 1,2-Dichloroethane | | 520 | U |
| 79-01-6 | Trichloroethene | | 520 | U |
| 78-87-5 | 1,2-Dichloropropane | | 520 | U |
| 74-95-3 | Dibromomethane | | 520 | U |
| 75-27-4 | Bromodichloromethane | | 520 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 520 | U |
| 108-88-3 | Toluene | | 520 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 520 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 520 | U |
| 127-18-4 | Tetrachloroethene | | 2800 | D |
| 142-28-9 | 1,3-Dichloropropane | | 520 | U |
| 124-48-1 | Dibromochloromethane | | 520 | U |
| 106-93-4 | 1,2-Dibromoethane | | 520 | U |
| 108-90-7 | Chlorobenzene | | 520 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 520 | U |
| 100-41-4 | Ethylbenzene | | 520 | U |
| 1330-20-7 | Xylene (para & meta) | | 520 | U |
| 95-47-6 | Xylene (Ortho) | | 520 | U |
| 100-42-5 | Styrene | | 520 | U |
| 75-25-2 | Bromoform | | 520 | U |
| 98-82-8 | Isopropylbenzene | | 520 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-07

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_07 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_7.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/19/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 520 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 520 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 520 | U |
| 103-65-1 | n-Propylbenzene | | 520 | U |
| 95-49-8 | 2-Chlorotoluene | | 520 | U |
| 106-43-4 | 4-Chlorotoluene | | 520 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 520 | U |
| 98-06-6 | tert-Butylbenzene | | 520 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 520 | U |
| 135-98-8 | sec-Butylbenzene | | 520 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 520 | U |
| 99-87-6 | 4-Isopropyltoluene | | 520 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 520 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 520 | U |
| 104-51-8 | n-Butylbenzene | | 520 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 520 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 520 | U |
| 87-68-3 | Hexachlorobutadiene | | 520 | U |
| 91-20-3 | Naphthalene | | 520 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 520 | U |
| 1634-04-4 | MTBE | | 520 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 520 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-08

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_08 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_8.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/19/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | | 520 | U |
| 74-87-3 | Chloromethane | | 520 | U |
| 75-01-4 | Vinyl chloride | | 520 | U |
| 74-83-9 | Bromomethane | | 520 | U |
| 75-00-3 | Chloroethane | | 520 | U |
| 75-69-4 | Trichlorofluoromethane | | 520 | U |
| 75-35-4 | 1,1-Dichloroethene | | 520 | U |
| 75-09-2 | Methylene chloride | | 520 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 520 | U |
| 75-34-3 | 1,1-Dichloroethane | | 520 | U |
| 594-20-7 | 2,2-Dichloropropane | | 520 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 520 | U |
| 74-97-5 | Bromochloromethane | | 520 | U |
| 67-66-3 | Chloroform | | 520 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 520 | U |
| 56-23-5 | Carbon tetrachloride | | 520 | U |
| 563-58-6 | 1,1-Dichloropropene | | 520 | U |
| 71-43-2 | Benzene | | 520 | U |
| 107-06-2 | 1,2-Dichloroethane | | 520 | U |
| 79-01-6 | Trichloroethene | | 520 | U |
| 78-87-5 | 1,2-Dichloropropane | | 520 | U |
| 74-95-3 | Dibromomethane | | 520 | U |
| 75-27-4 | Bromodichloromethane | | 520 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 520 | U |
| 108-88-3 | Toluene | | 520 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 520 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 520 | U |
| 127-18-4 | Tetrachloroethene | | 5800 | D |
| 142-28-9 | 1,3-Dichloropropane | | 520 | U |
| 124-48-1 | Dibromochloromethane | | 520 | U |
| 106-93-4 | 1,2-Dibromoethane | | 520 | U |
| 108-90-7 | Chlorobenzene | | 520 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 520 | U |
| 100-41-4 | Ethylbenzene | | 520 | U |
| 1330-20-7 | Xylene (para & meta) | | 520 | U |
| 95-47-6 | Xylene (Ortho) | | 520 | U |
| 100-42-5 | Styrene | | 520 | U |
| 75-25-2 | Bromoform | | 520 | U |
| 98-82-8 | Isopropylbenzene | | 520 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-08

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_08 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_8.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/19/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 520 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 520 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 520 | U |
| 103-65-1 | n-Propylbenzene | | 520 | U |
| 95-49-8 | 2-Chlorotoluene | | 520 | U |
| 106-43-4 | 4-Chlorotoluene | | 520 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 520 | U |
| 98-06-6 | tert-Butylbenzene | | 520 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 520 | U |
| 135-98-8 | sec-Butylbenzene | | 520 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 520 | U |
| 99-87-6 | 4-Isopropyltoluene | | 520 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 520 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 520 | U |
| 104-51-8 | n-Butylbenzene | | 520 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 520 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 520 | U |
| 87-68-3 | Hexachlorobutadiene | | 520 | U |
| 91-20-3 | Naphthalene | | 520 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 520 | U |
| 1634-04-4 | MTBE | | 520 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 520 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-09

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_09 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_9.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/19/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | | 520 | U |
| 74-87-3 | Chloromethane | | 520 | U |
| 75-01-4 | Vinyl chloride | | 520 | U |
| 74-83-9 | Bromomethane | | 520 | U |
| 75-00-3 | Chloroethane | | 520 | U |
| 75-69-4 | Trichlorofluoromethane | | 520 | U |
| 75-35-4 | 1,1-Dichloroethene | | 520 | U |
| 75-09-2 | Methylene chloride | | 520 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 520 | U |
| 75-34-3 | 1,1-Dichloroethane | | 520 | U |
| 594-20-7 | 2,2-Dichloropropane | | 520 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 520 | U |
| 74-97-5 | Bromochloromethane | | 520 | U |
| 67-66-3 | Chloroform | | 520 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 520 | U |
| 56-23-5 | Carbon tetrachloride | | 520 | U |
| 563-58-6 | 1,1-Dichloropropene | | 520 | U |
| 71-43-2 | Benzene | | 520 | U |
| 107-06-2 | 1,2-Dichloroethane | | 520 | U |
| 79-01-6 | Trichloroethene | | 520 | U |
| 78-87-5 | 1,2-Dichloropropane | | 520 | U |
| 74-95-3 | Dibromomethane | | 520 | U |
| 75-27-4 | Bromodichloromethane | | 520 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 520 | U |
| 108-88-3 | Toluene | | 520 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 520 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 520 | U |
| 127-18-4 | Tetrachloroethene | | 3100 | D |
| 142-28-9 | 1,3-Dichloropropane | | 520 | U |
| 124-48-1 | Dibromochloromethane | | 520 | U |
| 106-93-4 | 1,2-Dibromoethane | | 520 | U |
| 108-90-7 | Chlorobenzene | | 520 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 520 | U |
| 100-41-4 | Ethylbenzene | | 520 | U |
| 1330-20-7 | Xylene (para & meta) | | 520 | U |
| 95-47-6 | Xylene (Ortho) | | 520 | U |
| 100-42-5 | Styrene | | 520 | U |
| 75-25-2 | Bromoform | | 520 | U |
| 98-82-8 | Isopropylbenzene | | 520 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-09

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_09 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_9.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/19/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 520 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 520 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 520 | U |
| 103-65-1 | n-Propylbenzene | | 520 | U |
| 95-49-8 | 2-Chlorotoluene | | 520 | U |
| 106-43-4 | 4-Chlorotoluene | | 520 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 520 | U |
| 98-06-6 | tert-Butylbenzene | | 520 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 520 | U |
| 135-98-8 | sec-Butylbenzene | | 520 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 520 | U |
| 99-87-6 | 4-Isopropyltoluene | | 520 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 520 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 520 | U |
| 104-51-8 | n-Butylbenzene | | 520 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 520 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 520 | U |
| 87-68-3 | Hexachlorobutadiene | | 520 | U |
| 91-20-3 | Naphthalene | | 520 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 520 | U |
| 1634-04-4 | MTBE | | 520 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 520 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-10

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_10 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_10.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/20/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | | 520 | U |
| 74-87-3 | Chloromethane | | 520 | U |
| 75-01-4 | Vinyl chloride | | 520 | U |
| 74-83-9 | Bromomethane | | 520 | U |
| 75-00-3 | Chloroethane | | 520 | U |
| 75-69-4 | Trichlorofluoromethane | | 520 | U |
| 75-35-4 | 1,1-Dichloroethene | | 520 | U |
| 75-09-2 | Methylene chloride | | 520 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 520 | U |
| 75-34-3 | 1,1-Dichloroethane | | 520 | U |
| 594-20-7 | 2,2-Dichloropropane | | 520 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 520 | U |
| 74-97-5 | Bromochloromethane | | 520 | U |
| 67-66-3 | Chloroform | | 520 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 520 | U |
| 56-23-5 | Carbon tetrachloride | | 520 | U |
| 563-58-6 | 1,1-Dichloropropene | | 520 | U |
| 71-43-2 | Benzene | | 520 | U |
| 107-06-2 | 1,2-Dichloroethane | | 520 | U |
| 79-01-6 | Trichloroethene | | 520 | U |
| 78-87-5 | 1,2-Dichloropropane | | 520 | U |
| 74-95-3 | Dibromomethane | | 520 | U |
| 75-27-4 | Bromodichloromethane | | 520 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 520 | U |
| 108-88-3 | Toluene | | 520 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 520 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 520 | U |
| 127-18-4 | Tetrachloroethene | | 3600 | D |
| 142-28-9 | 1,3-Dichloropropane | | 520 | U |
| 124-48-1 | Dibromochloromethane | | 520 | U |
| 106-93-4 | 1,2-Dibromoethane | | 520 | U |
| 108-90-7 | Chlorobenzene | | 520 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 520 | U |
| 100-41-4 | Ethylbenzene | | 520 | U |
| 1330-20-7 | Xylene (para & meta) | | 520 | U |
| 95-47-6 | Xylene (Ortho) | | 520 | U |
| 100-42-5 | Styrene | | 520 | U |
| 75-25-2 | Bromoform | | 520 | U |
| 98-82-8 | Isopropylbenzene | | 520 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-10

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_10 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_10.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/20/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 520 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 520 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 520 | U |
| 103-65-1 | n-Propylbenzene | | 520 | U |
| 95-49-8 | 2-Chlorotoluene | | 520 | U |
| 106-43-4 | 4-Chlorotoluene | | 520 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 520 | U |
| 98-06-6 | tert-Butylbenzene | | 520 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 520 | U |
| 135-98-8 | sec-Butylbenzene | | 520 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 520 | U |
| 99-87-6 | 4-Isopropyltoluene | | 520 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 520 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 520 | U |
| 104-51-8 | n-Butylbenzene | | 520 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 520 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 520 | U |
| 87-68-3 | Hexachlorobutadiene | | 520 | U |
| 91-20-3 | Naphthalene | | 520 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 520 | U |
| 1634-04-4 | MTBE | | 520 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 520 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-11

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_11 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_11.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/20/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|----|
| 75-71-8 | Dichlorodifluoromethane | | 520 | U |
| 74-87-3 | Chloromethane | | 520 | U |
| 75-01-4 | Vinyl chloride | | 520 | U |
| 74-83-9 | Bromomethane | | 520 | U |
| 75-00-3 | Chloroethane | | 520 | U |
| 75-69-4 | Trichlorofluoromethane | | 520 | U |
| 75-35-4 | 1,1-Dichloroethene | | 520 | U |
| 75-09-2 | Methylene chloride | | 520 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 520 | U |
| 75-34-3 | 1,1-Dichloroethane | | 520 | U |
| 594-20-7 | 2,2-Dichloropropane | | 520 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 520 | U |
| 74-97-5 | Bromochloromethane | | 520 | U |
| 67-66-3 | Chloroform | | 520 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 520 | U |
| 56-23-5 | Carbon tetrachloride | | 520 | U |
| 563-58-6 | 1,1-Dichloropropene | | 520 | U |
| 71-43-2 | Benzene | | 520 | U |
| 107-06-2 | 1,2-Dichloroethane | | 520 | U |
| 79-01-6 | Trichloroethene | | 520 | U |
| 78-87-5 | 1,2-Dichloropropane | | 520 | U |
| 74-95-3 | Dibromomethane | | 520 | U |
| 75-27-4 | Bromodichloromethane | | 520 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 520 | U |
| 108-88-3 | Toluene | | 520 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 520 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 520 | U |
| 127-18-4 | Tetrachloroethene | | 19000 | ED |
| 142-28-9 | 1,3-Dichloropropane | | 520 | U |
| 124-48-1 | Dibromochloromethane | | 520 | U |
| 106-93-4 | 1,2-Dibromoethane | | 520 | U |
| 108-90-7 | Chlorobenzene | | 520 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 520 | U |
| 100-41-4 | Ethylbenzene | | 520 | U |
| 1330-20-7 | Xylene (para & meta) | | 520 | U |
| 95-47-6 | Xylene (Ortho) | | 520 | U |
| 100-42-5 | Styrene | | 520 | U |
| 75-25-2 | Bromoform | | 520 | U |
| 98-82-8 | Isopropylbenzene | | 520 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-11

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_11 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_11.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/20/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 520 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 520 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 520 | U |
| 103-65-1 | n-Propylbenzene | | 520 | U |
| 95-49-8 | 2-Chlorotoluene | | 520 | U |
| 106-43-4 | 4-Chlorotoluene | | 520 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 520 | U |
| 98-06-6 | tert-Butylbenzene | | 520 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 520 | U |
| 135-98-8 | sec-Butylbenzene | | 520 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 520 | U |
| 99-87-6 | 4-Isopropyltoluene | | 520 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 520 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 520 | U |
| 104-51-8 | n-Butylbenzene | | 520 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 520 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 520 | U |
| 87-68-3 | Hexachlorobutadiene | | 520 | U |
| 91-20-3 | Naphthalene | | 520 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 520 | U |
| 1634-04-4 | MTBE | | 520 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 520 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-11

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_11 0.20 ml
 Sample wt/vol: 0.2 (g/ml) ML Lab File ID: 7489_11.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/20/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 210.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 1000 | U |
| 74-87-3 | Chloromethane | | 1000 | U |
| 75-01-4 | Vinyl chloride | | 1000 | U |
| 74-83-9 | Bromomethane | | 1000 | U |
| 75-00-3 | Chloroethane | | 1000 | U |
| 75-69-4 | Trichlorofluoromethane | | 1000 | U |
| 75-35-4 | 1,1-Dichloroethene | | 1000 | U |
| 75-09-2 | Methylene chloride | | 1000 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 1000 | U |
| 75-34-3 | 1,1-Dichloroethane | | 1000 | U |
| 594-20-7 | 2,2-Dichloropropane | | 1000 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 1000 | U |
| 74-97-5 | Bromochloromethane | | 1000 | U |
| 67-66-3 | Chloroform | | 1000 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 1000 | U |
| 56-23-5 | Carbon tetrachloride | | 1000 | U |
| 563-58-6 | 1,1-Dichloropropene | | 1000 | U |
| 71-43-2 | Benzene | | 1000 | U |
| 107-06-2 | 1,2-Dichloroethane | | 1000 | U |
| 79-01-6 | Trichloroethene | | 1000 | U |
| 78-87-5 | 1,2-Dichloropropane | | 1000 | U |
| 74-95-3 | Dibromomethane | | 1000 | U |
| 75-27-4 | Bromodichloromethane | | 1000 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 1000 | U |
| 108-88-3 | Toluene | | 1000 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 1000 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 1000 | U |
| 127-18-4 | Tetrachloroethene | | 18000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 1000 | U |
| 124-48-1 | Dibromochloromethane | | 1000 | U |
| 106-93-4 | 1,2-Dibromoethane | | 1000 | U |
| 108-90-7 | Chlorobenzene | | 1000 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 1000 | U |
| 100-41-4 | Ethylbenzene | | 1000 | U |
| 1330-20-7 | Xylene (para & meta) | | 1000 | U |
| 95-47-6 | Xylene (Ortho) | | 1000 | U |
| 100-42-5 | Styrene | | 1000 | U |
| 75-25-2 | Bromoform | | 1000 | U |
| 98-82-8 | Isopropylbenzene | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-11

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_11 0.20 ml
 Sample wt/vol: 0.2 (g/ml) ML Lab File ID: 7489_11.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/20/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 210.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 1000 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 1000 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 1000 | U |
| 103-65-1 | n-Propylbenzene | | 1000 | U |
| 95-49-8 | 2-Chlorotoluene | | 1000 | U |
| 106-43-4 | 4-Chlorotoluene | | 1000 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 1000 | U |
| 98-06-6 | tert-Butylbenzene | | 1000 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 1000 | U |
| 135-98-8 | sec-Butylbenzene | | 1000 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 1000 | U |
| 99-87-6 | 4-Isopropyltoluene | | 1000 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 1000 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 1000 | U |
| 104-51-8 | n-Butylbenzene | | 1000 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 1000 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 1000 | U |
| 87-68-3 | Hexachlorobutadiene | | 1000 | U |
| 91-20-3 | Naphthalene | | 1000 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 1000 | U |
| 1634-04-4 | MTBE | | 1000 | U |
| 67-64-1 | Acetone | | 2100 | U |
| 75-15-0 | Carbon disulfide | | 1000 | U |
| 78-93-3 | 2-Butanone | | 2100 | U |
| 109-99-9 | Tetrahydrofuran | | 2100 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 2100 | U |
| 591-78-6 | 2-Hexanone | | 2100 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 2100 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-12

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_12 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_12.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/20/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | | 520 | U |
| 74-87-3 | Chloromethane | | 520 | U |
| 75-01-4 | Vinyl chloride | | 520 | U |
| 74-83-9 | Bromomethane | | 520 | U |
| 75-00-3 | Chloroethane | | 520 | U |
| 75-69-4 | Trichlorofluoromethane | | 520 | U |
| 75-35-4 | 1,1-Dichloroethene | | 520 | U |
| 75-09-2 | Methylene chloride | | 520 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 520 | U |
| 75-34-3 | 1,1-Dichloroethane | | 520 | U |
| 594-20-7 | 2,2-Dichloropropane | | 520 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 520 | U |
| 74-97-5 | Bromochloromethane | | 520 | U |
| 67-66-3 | Chloroform | | 520 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 520 | U |
| 56-23-5 | Carbon tetrachloride | | 520 | U |
| 563-58-6 | 1,1-Dichloropropene | | 520 | U |
| 71-43-2 | Benzene | | 520 | U |
| 107-06-2 | 1,2-Dichloroethane | | 520 | U |
| 79-01-6 | Trichloroethene | | 520 | U |
| 78-87-5 | 1,2-Dichloropropane | | 520 | U |
| 74-95-3 | Dibromomethane | | 520 | U |
| 75-27-4 | Bromodichloromethane | | 520 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 520 | U |
| 108-88-3 | Toluene | | 520 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 520 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 520 | U |
| 127-18-4 | Tetrachloroethene | | 2500 | D |
| 142-28-9 | 1,3-Dichloropropane | | 520 | U |
| 124-48-1 | Dibromochloromethane | | 520 | U |
| 106-93-4 | 1,2-Dibromoethane | | 520 | U |
| 108-90-7 | Chlorobenzene | | 520 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 520 | U |
| 100-41-4 | Ethylbenzene | | 520 | U |
| 1330-20-7 | Xylene (para & meta) | | 520 | U |
| 95-47-6 | Xylene (Ortho) | | 520 | U |
| 100-42-5 | Styrene | | 520 | U |
| 75-25-2 | Bromoform | | 520 | U |
| 98-82-8 | Isopropylbenzene | | 520 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7489-12

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7489_12 0.40 ml
 Sample wt/vol: 0.4 (g/ml) ML Lab File ID: 89_12.D
 Level: (low/med) LOW Date Received: 12/19/06
 % Moisture: not dec. _____ Date Analyzed: 12/20/06
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 105.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 520 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 520 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 520 | U |
| 103-65-1 | n-Propylbenzene | | 520 | U |
| 95-49-8 | 2-Chlorotoluene | | 520 | U |
| 106-43-4 | 4-Chlorotoluene | | 520 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 520 | U |
| 98-06-6 | tert-Butylbenzene | | 520 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 520 | U |
| 135-98-8 | sec-Butylbenzene | | 520 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 520 | U |
| 99-87-6 | 4-Isopropyltoluene | | 520 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 520 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 520 | U |
| 104-51-8 | n-Butylbenzene | | 520 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 520 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 520 | U |
| 87-68-3 | Hexachlorobutadiene | | 520 | U |
| 91-20-3 | Naphthalene | | 520 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 520 | U |
| 1634-04-4 | MTBE | | 520 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 520 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

3.0 Raw Data and Chromatograms

Raw data and Chromatograms are attached.

Acq On : 20 Dec 2006 1:50 am

Sample : 7489_1 0.40 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Dec 20 9:57 2006

Vial: 12

Operator: TS

Inst : GC/MS #2

Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.06 | 96 | 200667 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.91 | 117 | 113851 | 25.00 | ug/L | -0.04 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 45713 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|--------|-------|------|--------------------|
| 16) Dibromofluoromethane | 8.37 | 113 | 51077 | 26.33 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 105.32% |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 41721 | 24.85 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 99.40% |
| 28) Toluene-d8 | 13.13 | 98 | 158156 | 23.52 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 94.08% |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 44510 | 21.10 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 84.40% |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. | d |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#)=qualifier out of range (m)=manual integration

Data File : C:\HPCHEM\1\DATA\12_19_06\9_1.D

Acq On : 20 Dec 2006 1:50 am

Sample : 7489_1 0.40 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Dec 20 9:57 2006

Vial: 12

Operator: TS

Inst : GC/MS #2

Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.48 | 166 | 101936 | 62.93 | ug/L | 96 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

9_1.D 71NV16_6.M

Wed Jan 03 16:25:27 2007

GC#2

Page 2

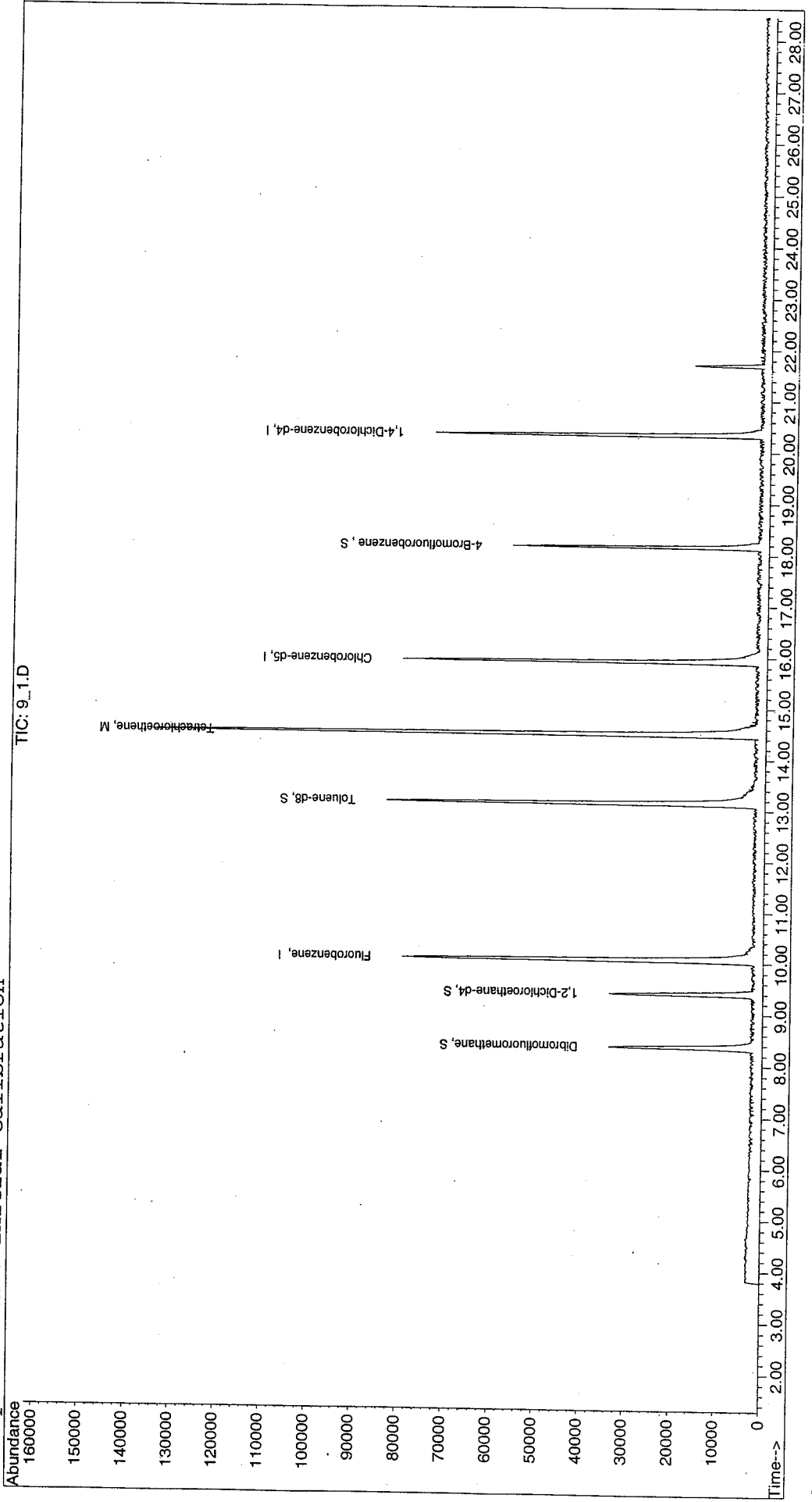
Quantitation Report

Data File : C:\HPCHEM\1\DATA\12_19_06\9_1.D
Acq On : 20 Dec 2006 1:50 am
Sample : 7489_1 0.40 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Dec 20 9:57 2006

Vial: 12
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Nov 27 09:29:43 2006
Response via : Initial Calibration



Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.06 | 96 | 200089 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.91 | 117 | 115201 | 25.00 | ug/L | -0.04 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 48672 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|-------|
| 16) Dibromofluoromethane | 8.37 | 113 | 50898 | 26.31 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 105.24% | |
| 19) 1,2-Dichloroethane-d4 | 9.41 | 65 | 40704 | 24.31 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 97.24% | |
| 28) Toluene-d8 | 13.13 | 98 | 153920 | 22.95 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 91.80% | |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 46044 | 21.57 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 86.28% | |

Target Compounds

| | | | | | Qvalue |
|-------------------------------|------|-----|---|------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\12_19_06\9_2.D
 Acq On : 20 Dec 2006 2:24 am
 Sample : 7489_2 dup 0.40 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Dec 20 9:57 2006

Vial: 13
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

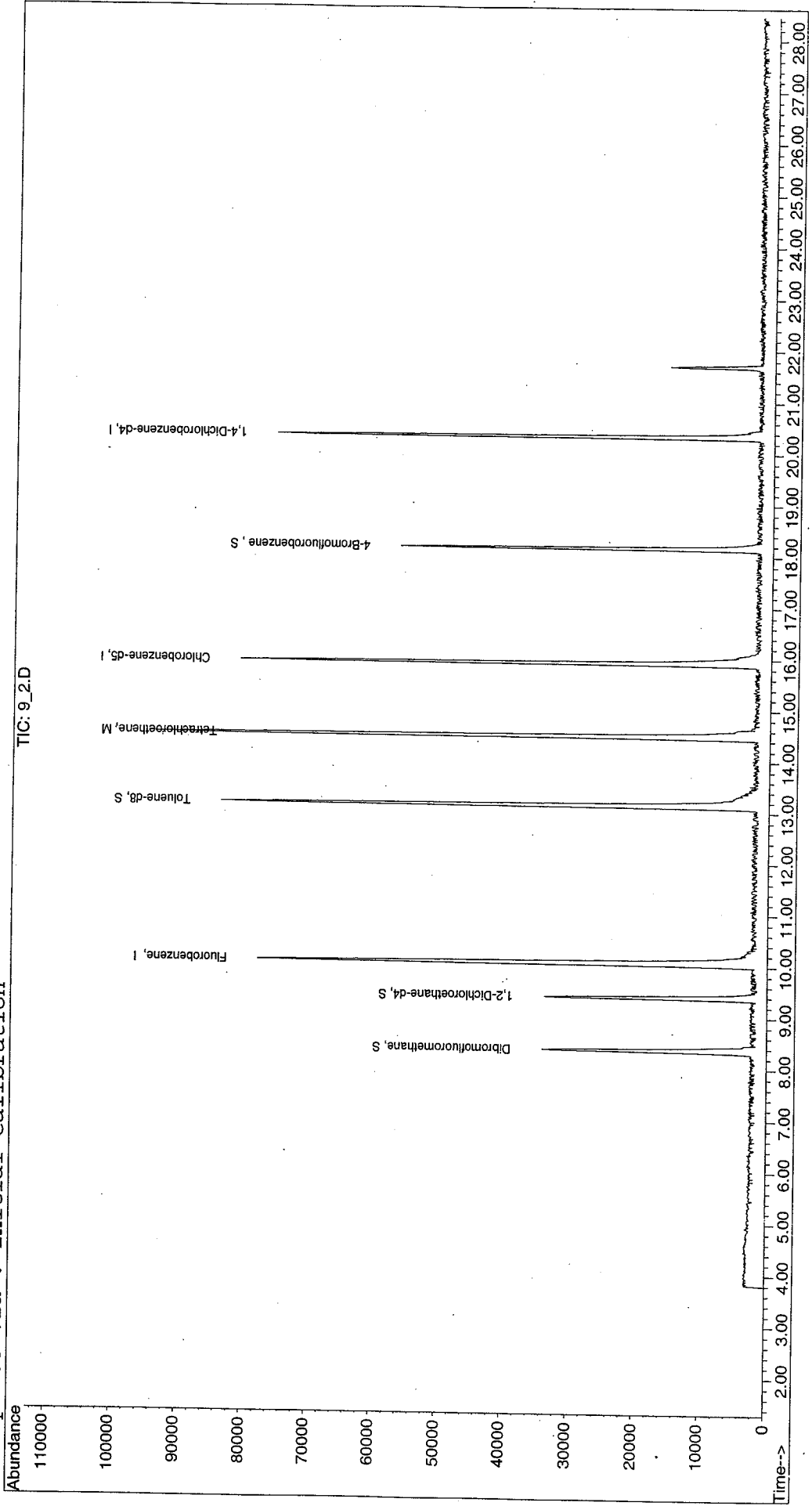
Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Thu Nov 16 17:01:55 2006
 Response via : Initial Calibration
 DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.48 | 166 | 69511 | 43.04 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\12_19_06\9_2.D
Acq On : 20 Dec 2006 2:24 am
Sample : 7489_2 dup 0.40 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Dec 20 9:57 2006
Quant Results File: 71NV16_6.RES
Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Nov 27 09:29:43 2006
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\12_19_06\89_3.D

Vial: 2

Acq On : 19 Dec 2006 8:04 pm

Operator: TS

Sample : 7489_03 0.40 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Dec 20 9:51 2006

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------|--------|------|----------|-------|---------|----------|
| 1) Fluorobenzene | 10.07 | 96 | 237315 | 25.00 | ug/L | -0.03 |
| 36) Chlorobenzene-d5 | 15.92 | 117 | 144261 | 25.00 | ug/L | -0.03 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 56095 | 25.00 | ug/L | -0.03 |
| System Monitoring Compounds | | | | | | |
| 16) Dibromofluoromethane | 8.38 | 113 | 61948 | 27.00 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 108.00% | |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 49279 | 24.82 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 99.28% | |
| 28) Toluene-d8 | 13.14 | 98 | 191268 | 24.05 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 96.20% | |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 57248 | 21.41 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 85.64% | |

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|--------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. d | |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\12_19_06\89_3.D
Acq On : 19 Dec 2006 8:04 pm
Sample : 7489_03 0.40 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Dec 20 9:51 2006

Vial: 2
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Thu Nov 16 17:01:55 2006
Response via : Initial Calibration
DataAcq Meth : RUN1

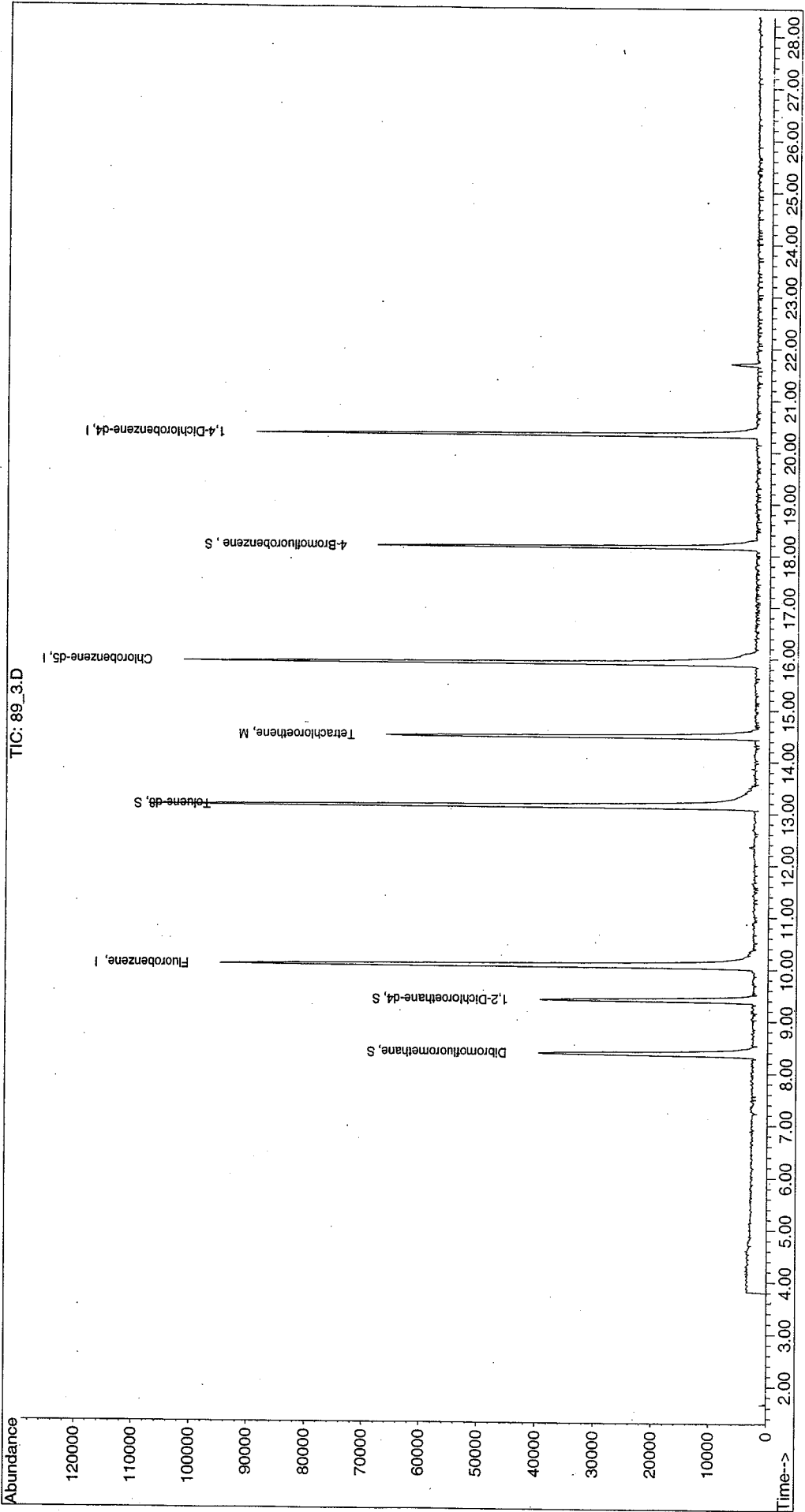
| Compound | R.T. | QIon | Response | Conc Unit | Qvalue |
|---------------------------------|-------|------|----------|------------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 48032 | 25.07 ug/L | 97 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. d | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | |

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\12_19_06\89_3.D
Acq On : 19 Dec 2006 8:04 pm
Sample : 7489_03 0.40 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Dec 20 9:51 2006
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Nov 27 09:29:43 2006
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\12_19_06\89_4.D

Acq On : 19 Dec 2006 8:39 pm

Sample : 7489_04 0.40 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Dec 20 9:52 2006

Vial: 3

Operator: TS

Inst : GC/MS #2

Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.07 | 96 | 229964 | 25.00 | ug/L | -0.03 |
| 36) Chlorobenzene-d5 | 15.92 | 117 | 144284 | 25.00 | ug/L | -0.03 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 55982 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|-------|
| 16) Dibromofluoromethane | 8.37 | 113 | 59293 | 26.67 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 106.68% | |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 47740 | 24.81 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 99.24% | |
| 28) Toluene-d8 | 13.14 | 98 | 189539 | 24.59 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 98.36% | |
| 45) 4-Bromofluorobenzene | 18.17 | 95 | 56184 | 21.01 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 84.04% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. | d |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\12_19_06\89_4.D

Acq On : 19 Dec 2006 8:39 pm

Sample : 7489_04 0.40 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Dec 20 9:52 2006

Vial: 3

Operator: TS

Inst : GC/MS #2

Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 77537 | 41.77 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | d | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

89_4.D 71NV16_6.M

Wed Jan 03 16:25:08 2007

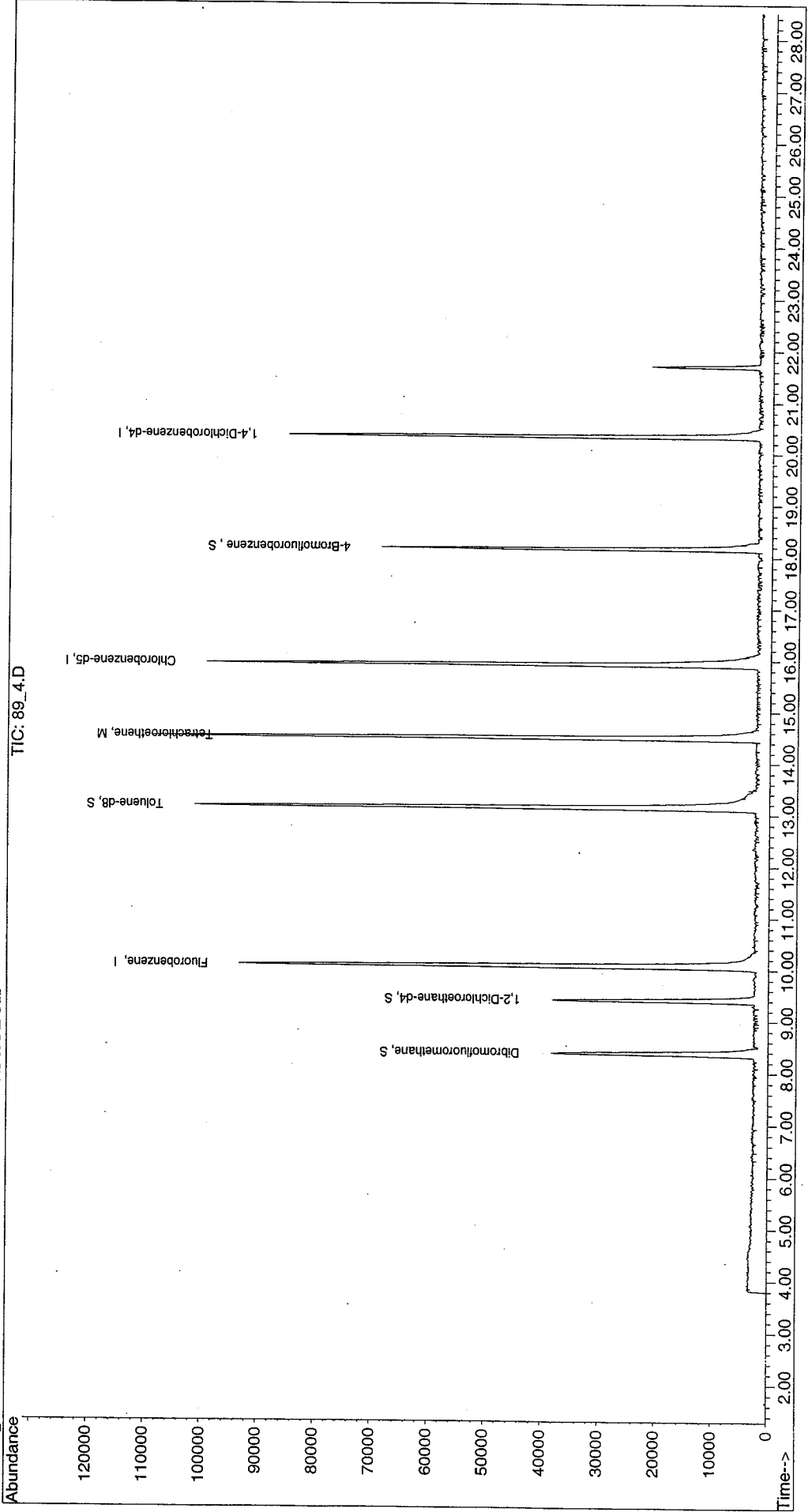
GC#2

Page 2

Quantitation Report

Data File : C:\HPCHEM\1\DATA\12_19_06\89_4.D
Acq On : 19 Dec 2006 8:39 pm
Sample : 7489_04 0.40 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Dec 20 9:52 2006
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Nov 27 09:29:43 2006
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\12_19_06\89_5.D
 Acq On : 19 Dec 2006 9:13 pm
 Sample : 7489_05 0.40 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Dec 20 9:53 2006

Vial: 4
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Thu Nov 16 17:01:55 2006
 Response via : Initial Calibration
 DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.07 | 96 | 226549 | 25.00 | ug/L | -0.03 |
| 36) Chlorobenzene-d5 | 15.92 | 117 | 134788 | 25.00 | ug/L | -0.03 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 53671 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|-------|
| 16) Dibromofluoromethane | 8.37 | 113 | 59388 | 27.12 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 108.48% | |
| 19) 1,2-Dichloroethane-d4 | 9.41 | 65 | 48243 | 25.45 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 101.80% | |
| 28) Toluene-d8 | 13.14 | 98 | 185291 | 24.40 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 97.60% | |
| 45) 4-Bromofluorobenzene | 18.17 | 95 | 54821 | 21.95 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 87.80% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. | d |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\12_19_06\89_5.D

Acq On : 19 Dec 2006 9:13 pm

Sample : 7489_05 0.40 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Dec 20 9:53 2006

Vial: 4

Operator: TS

Inst : GC/MS #2

Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 48174 | 26.34 | ug/L | 99 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | d | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

89_5.D 71NV16_6.M

Wed Jan 03 16:25:11 2007

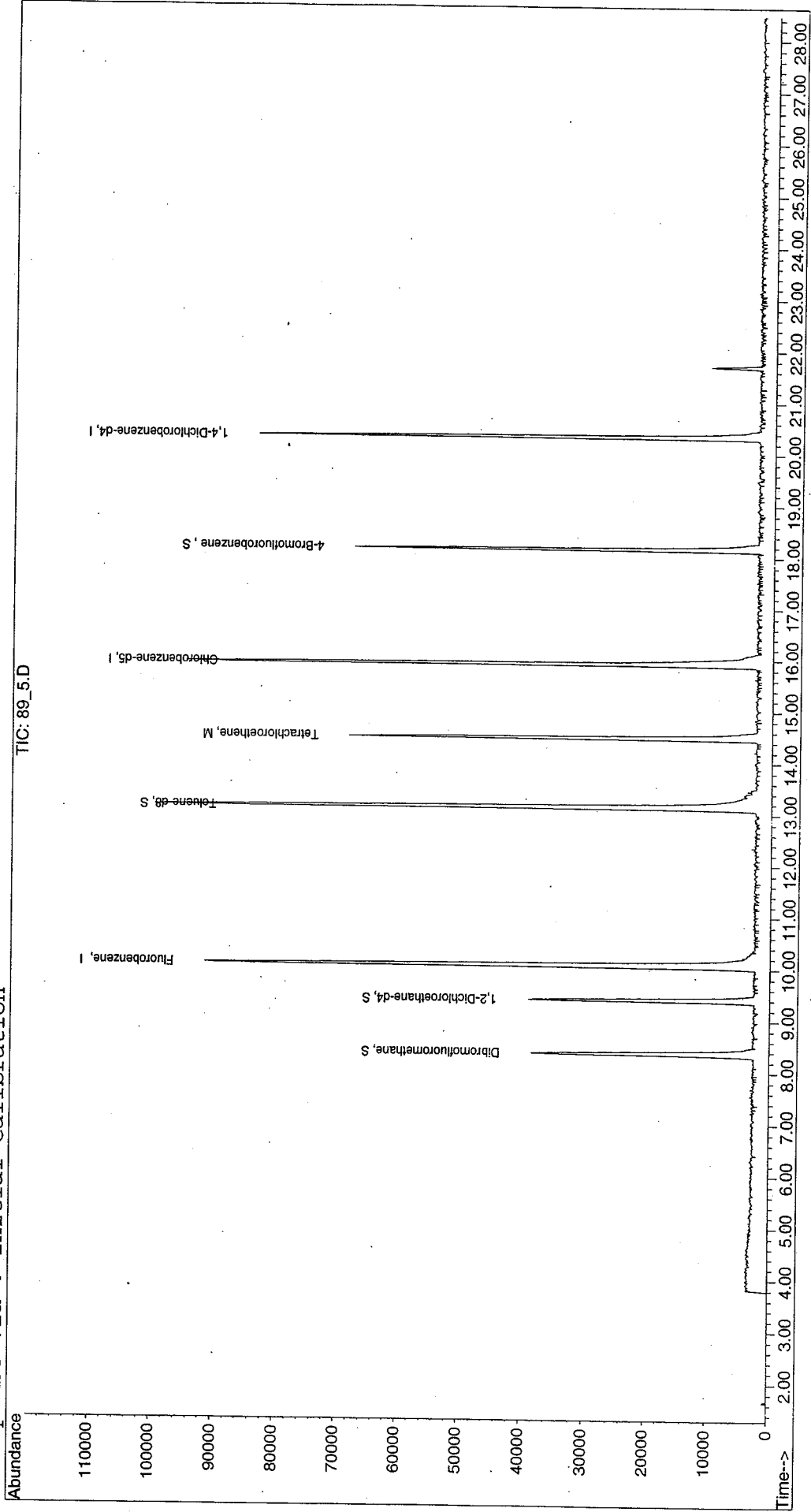
GC#2

Page 2

Quantitation Report

Data File : C:\HPCHEM\1\DATA\12_19_06\89_5.D
Acq On : 19 Dec 2006 9:13 pm
Sample : 7489_05 0.40 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Dec 20 9:53 2006
Vial: 4
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Nov 27 09:29:43 2006
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\12_19_06\89_6.D

Acq On : 19 Dec 2006 9:47 pm

Sample : 7489_06 0.40 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Dec 20 9:54 2006

Vial: 5

Operator: TS

Inst : GC/MS #2

Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.06 | 96 | 228951 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.92 | 117 | 140153 | 25.00 | ug/L | -0.03 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 53721 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|--------|-------|------|--------------------|
| 16) Dibromofluoromethane | 8.37 | 113 | 58491 | 26.43 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 105.72% |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 47138 | 24.61 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 98.44% |
| 28) Toluene-d8 | 13.13 | 98 | 186234 | 24.27 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 97.08% |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 55248 | 21.27 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 85.08% |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Acq On : 19 Dec 2006 9:47 pm

Operator: TS

Sample : 7489_06 0.40 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Dec 20 9:54 2006

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

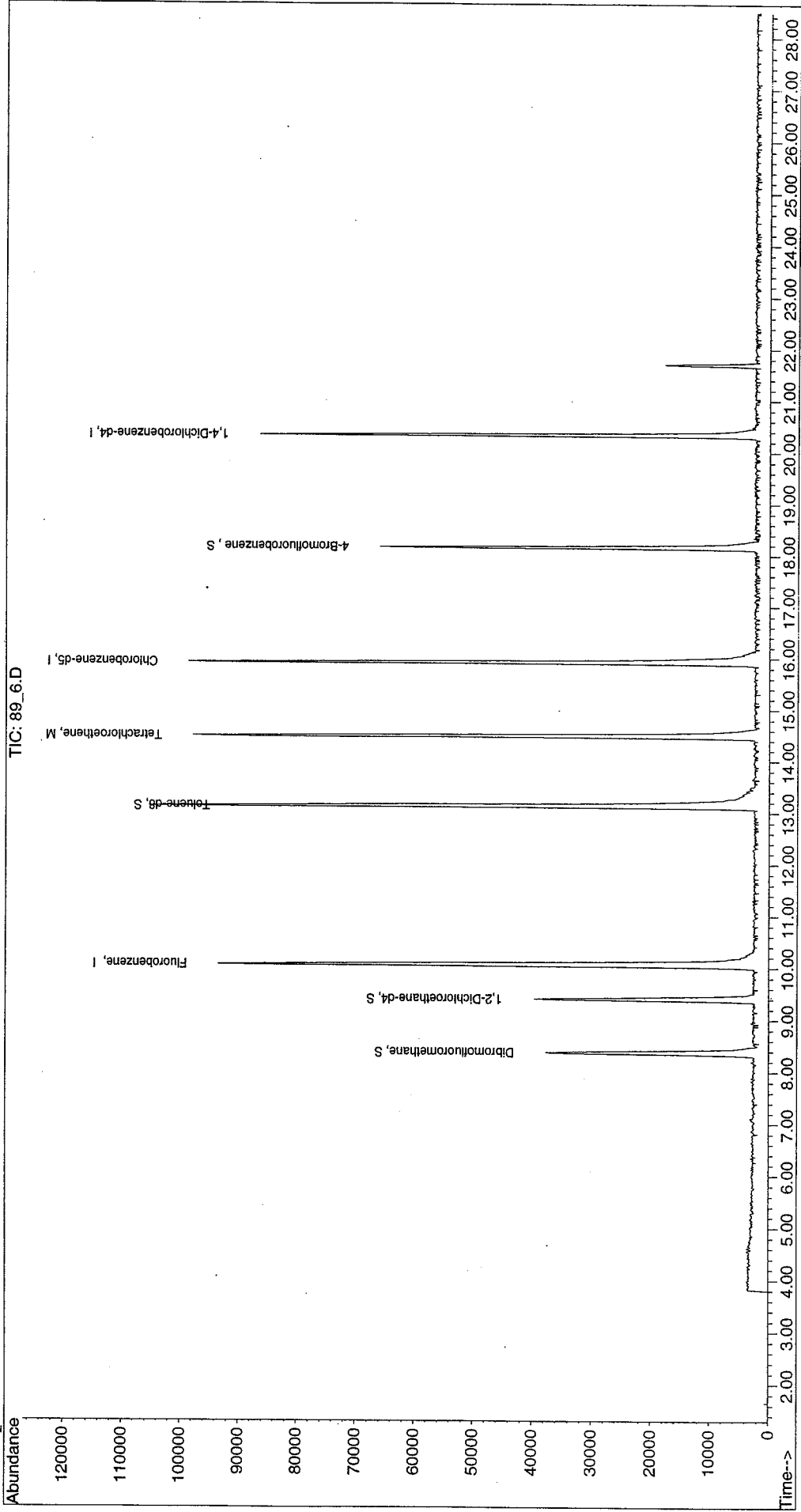
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 69211 | 37.45 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | d | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\12_19_06\89_6.D Vial: 5
Acq On : 19 Dec 2006 9:47 pm Operator: TS
Sample : 7489_06 0.40 ml Inst : GC/MS #2
Misc : Multiplr: 1.00
MS Integration Params: ODD.P
Quant Time: Dec 20 9:54 2006 Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Nov 27 09:29:43 2006
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\12_19_06\89_7.D
 Acq On : 19 Dec 2006 10:22 pm
 Sample : 7489_07 0.40 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Dec 20 9:54 2006

Vial: 6
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Thu Nov 16 17:01:55 2006
 Response via : Initial Calibration
 DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.06 | 96 | 224331 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.92 | 117 | 139403 | 25.00 | ug/L | -0.03 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 53025 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|-------|
| 16) Dibromofluoromethane | 8.37 | 113 | 59260 | 27.33 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 109.32% | |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 46557 | 24.80 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 99.20% | |
| 28) Toluene-d8 | 13.14 | 98 | 181464 | 24.14 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 96.56% | |
| 45) 4-Bromofluorobenzene | 18.17 | 95 | 54607 | 21.14 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 84.56% | |

Target Compounds

| | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\12_19_06\89_7.D

Acq On : 19 Dec 2006 10:22 pm

Sample : 7489_07 0.40 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Dec 20 9:54 2006

Vial: 6

Operator: TS

Inst : GC/MS #2

Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 48636 | 26.86 | ug/L | 97 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

89_7.D 71NV16_6.M

Wed Jan 03 16:25:18 2007

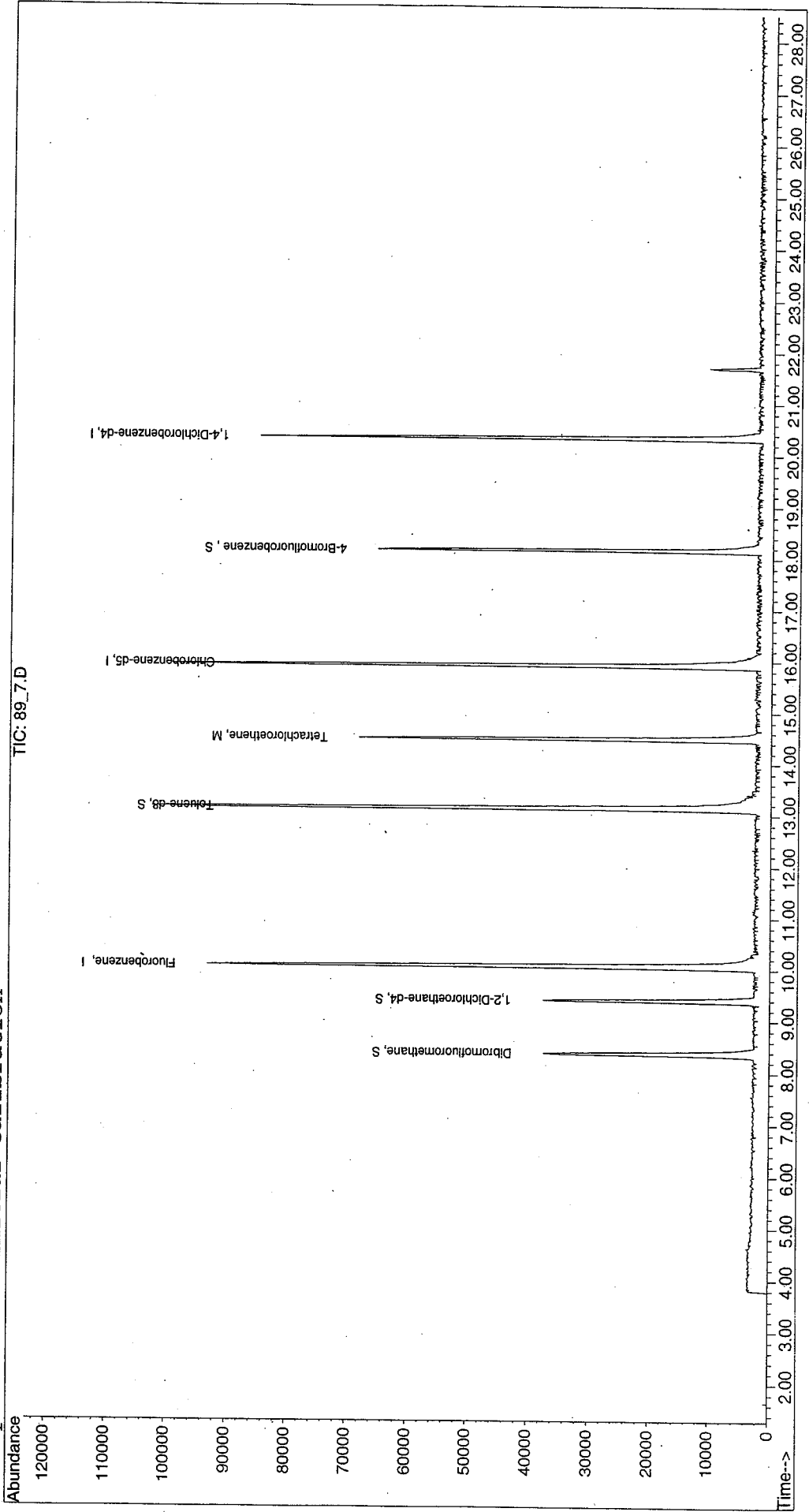
GC#2

Page 2

Quantitation report

Data File : C:\HPCHEM\1\DATA\12_19_06\89_7.D Vial: 6
Acq On : 19 Dec 2006 10:22 pm Operator: TS
Sample : 7489_07 0.40 ml Inst : GC/MS #2
Misc : Multiplr: 1.00
MS Integration Params: ODD.P
Quant Time: Dec 20 9:54 2006 Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Nov 27 09:29:43 2006
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\12_19_06\89_8.D
 Acq On : 19 Dec 2006 10:57 pm
 Sample : 7489_08 0.40 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Dec 20 9:55 2006

Vial: 7
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Thu Nov 16 17:01:55 2006
 Response via : Initial Calibration
 DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------|--------|------|----------|-------|---------|----------|
| 1) Fluorobenzene | 10.06 | 96 | 217961 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.92 | 117 | 138613 | 25.00 | ug/L | -0.03 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 52214 | 25.00 | ug/L | -0.03 |
| System Monitoring Compounds | | | | | | |
| 16) Dibromofluoromethane | 8.38 | 113 | 56720 | 26.92 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 107.68% | |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 46401 | 25.44 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 101.76% | |
| 28) Toluene-d8 | 13.13 | 98 | 180618 | 24.73 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 98.92% | |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 53895 | 20.98 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 83.92% | |

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Acq On : 19 Dec 2006 10:57 pm

Operator: TS

Sample : 7489_08 0.40 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Dec 20 9:55 2006

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

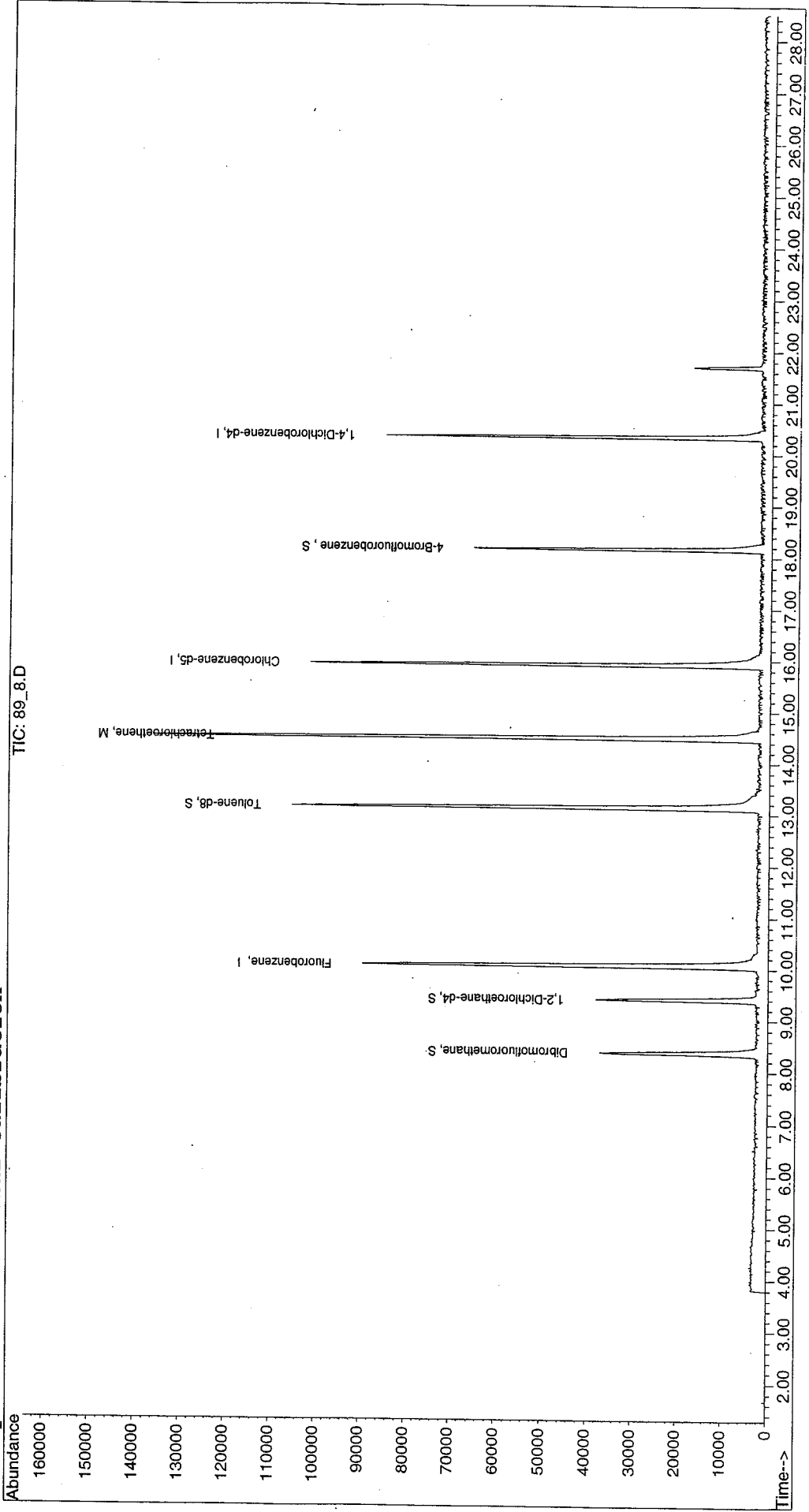
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 97432 | 55.38 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Quantitation report

Data File : C:\HPCHEM\1\DATA\12_19_06\89_8.D
Acq On : 19 Dec 2006 10:57 pm
Sample : 7489_08 0.40 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Dec 20 9:55 2006
Vial: 7
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Nov 27 09:29:43 2006
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\12_19_06\89_9.D
 Acq On : 19 Dec 2006 11:31 pm
 Sample : 7489_09 0.40 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Dec 20 9:56 2006

Vial: 8
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Thu Nov 16 17:01:55 2006
 Response via : Initial Calibration
 DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.06 | 96 | 214869 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.91 | 117 | 130309 | 25.00 | ug/L | -0.04 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 51410 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|-------|
| 16) Dibromofluoromethane | 8.38 | 113 | 54236 | 26.11 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 104.44% | |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 43536 | 24.22 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 96.88% | |
| 28) Toluene-d8 | 13.14 | 98 | 173524 | 24.10 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 96.40% | |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 50573 | 20.94 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 83.76% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. | d |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\12_19_06\89_9.D
 Acq On : 19 Dec 2006 11:31 pm
 Sample : 7489_09 0.40 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Dec 20 9:56 2006

Vial: 8
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Thu Nov 16 17:01:55 2006
 Response via : Initial Calibration
 DataAcq Meth : RUN1

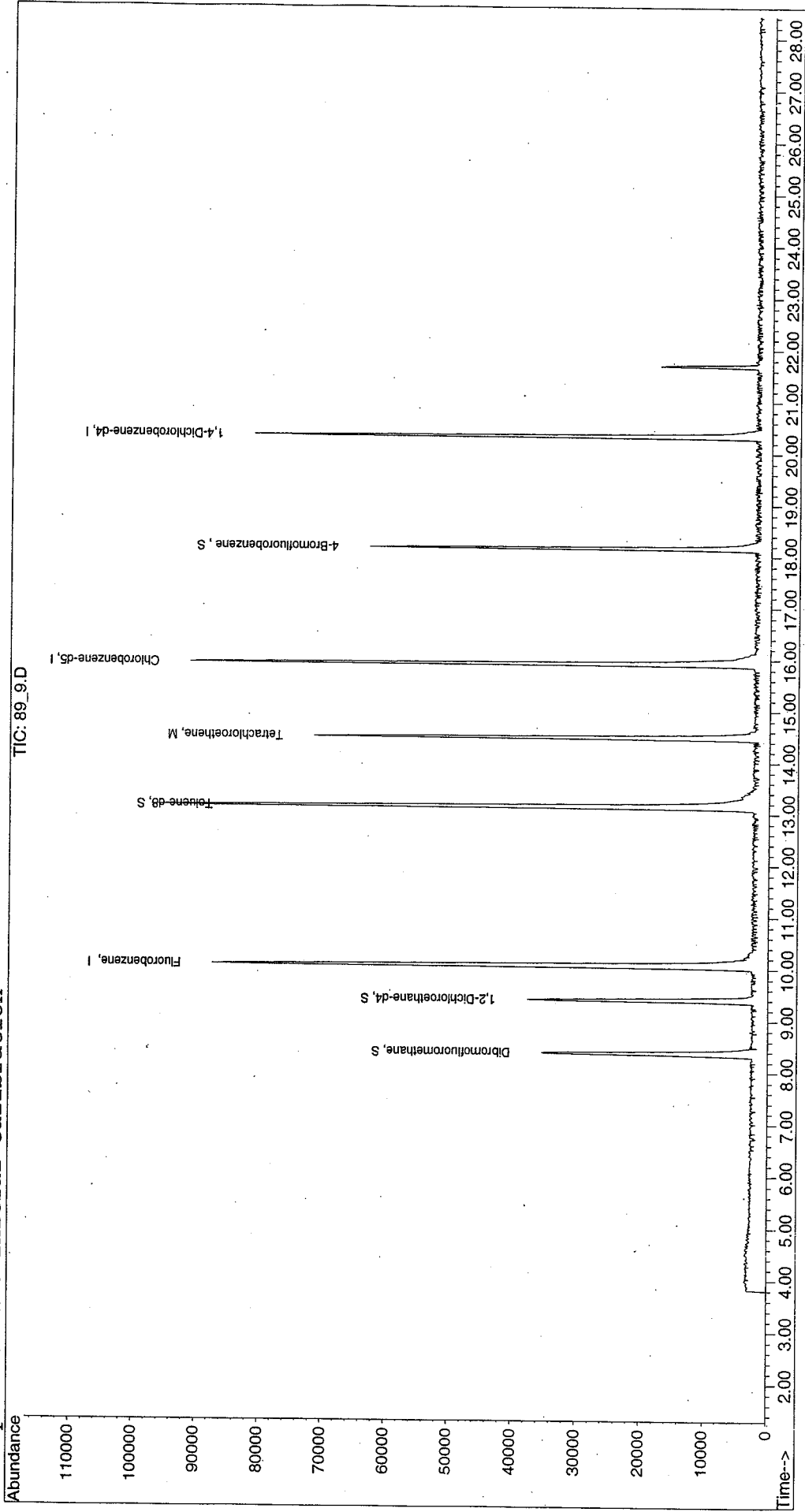
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|--------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 50521 | 29.13 | ug/L | 95 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | | N.D. | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | | N.D. d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | | N.D. | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | | N.D. | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | | N.D. | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | | N.D. | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | | N.D. | |
| 42) Styrene | 0.00 | 104 | 0 | | N.D. | |
| 43) Bromoform | 0.00 | 173 | 0 | | N.D. | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 46) Bromobenzene | 0.00 | 156 | 0 | | N.D. | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | | N.D. | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | | N.D. | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | | N.D. | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | | N.D. | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | | N.D. | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | | N.D. | |
| 65) Naphthalene | 0.00 | 128 | 0 | | N.D. | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | | N.D. | |
| 68) Acetone | 0.00 | 43 | 0 | | N.D. | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | | N.D. | |
| 70) 2-Butanone | 0.00 | 43 | 0 | | N.D. | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | | N.D. | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | | N.D. | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | | N.D. | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\12_19_06\89_9.D
Acq On : 19 Dec 2006 11:31 pm
Sample : 7489_09 0.40 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Dec 20 9:56 2006
Vial: 8
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Nov 27 09:29:43 2006
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\12_19_06\89_10.D
 Acq On : 20 Dec 2006 12:08 am
 Sample : 7489_10 0.40 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Dec 20 9:45 2006

Vial: 9
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Thu Nov 16 17:01:55 2006
 Response via : Initial Calibration
 DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.07 | 96 | 210874 | 25.00 | ug/L | -0.03 |
| 36) Chlorobenzene-d5 | 15.92 | 117 | 121461 | 25.00 | ug/L | -0.03 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 48640 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|-------|
| 16) Dibromofluoromethane | 8.37 | 113 | 54252 | 26.61 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 106.44% | |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 43586 | 24.70 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 98.80% | |
| 28) Toluene-d8 | 13.14 | 98 | 163613 | 23.15 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 92.60% | |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 46247 | 20.55 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 82.20% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. | d |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\12_19_06\89_10.D
Acq On : 20 Dec 2006 12:08 am
Sample : 7489_10 0.40 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Dec 20 9:45 2006

Vial: 9
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Thu Nov 16 17:01:55 2006
Response via : Initial Calibration
DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.48 | 166 | 58771 | 34.53 | ug/L | 94 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

89_10.D 71NV16_6.M

Wed Jan 03 16:24:50 2007

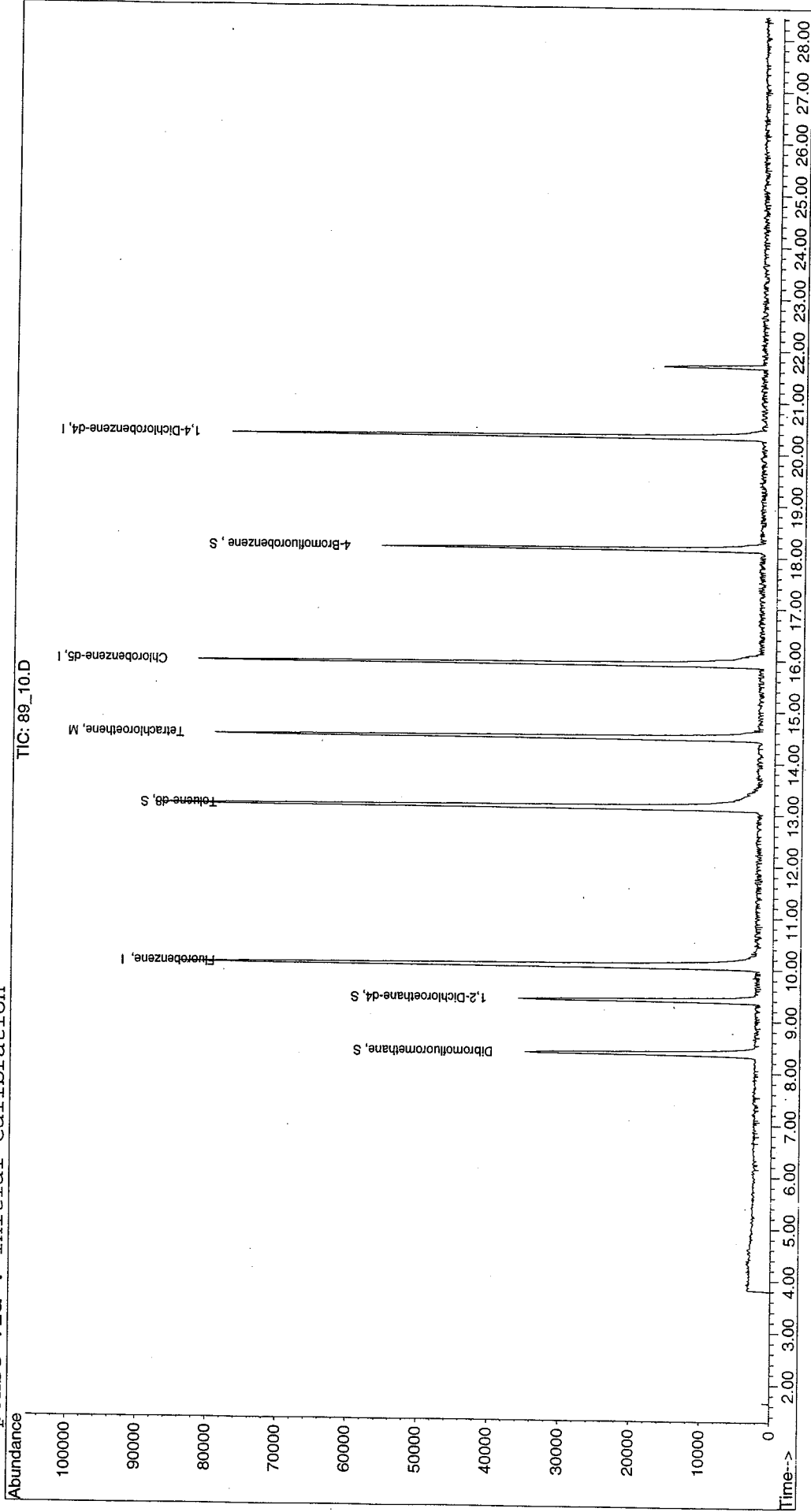
GC#2

Page 2

Quantitation Report

Data File : C:\HPCHEM\1\DATA\12_19_06\89_10.D
Acq On : 20 Dec 2006 12:08 am
Sample : 7489_10 0.40 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Dec 20 9:45 2006
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Nov 27 09:29:43 2006
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\12_19_06\89_11.D
 Acq On : 20 Dec 2006 12:42 am
 Sample : 7489_11 0.40 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Dec 20 9:46 2006

Vial: 10
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Thu Nov 16 17:01:55 2006
 Response via : Initial Calibration
 DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.06 | 96 | 206389 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.91 | 117 | 124677 | 25.00 | ug/L | -0.04 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 47886 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|--------|-------|------|--------------------|
| 16) Dibromofluoromethane | 8.38 | 113 | 52814 | 26.47 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 105.88% |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 43598 | 25.25 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 101.00% |
| 28) Toluene-d8 | 13.13 | 98 | 164892 | 23.84 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 95.36% |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 48726 | 21.09 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 84.36% |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | d | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | d | |

(#) = qualifier out of range (m) = manual integration
 89_11.D 71NV16_6.M Wed Jan 03 16:24:52 2007

Data File : C:\HPCHEM\1\DATA\12_19_06\89_11.D
 Acq On : 20 Dec 2006 12:42 am
 Sample : 7489_11 0.40 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Dec 20 9:46 2006

Vial: 10
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Thu Nov 16 17:01:55 2006
 Response via : Initial Calibration
 DataAcq Meth : RUN1

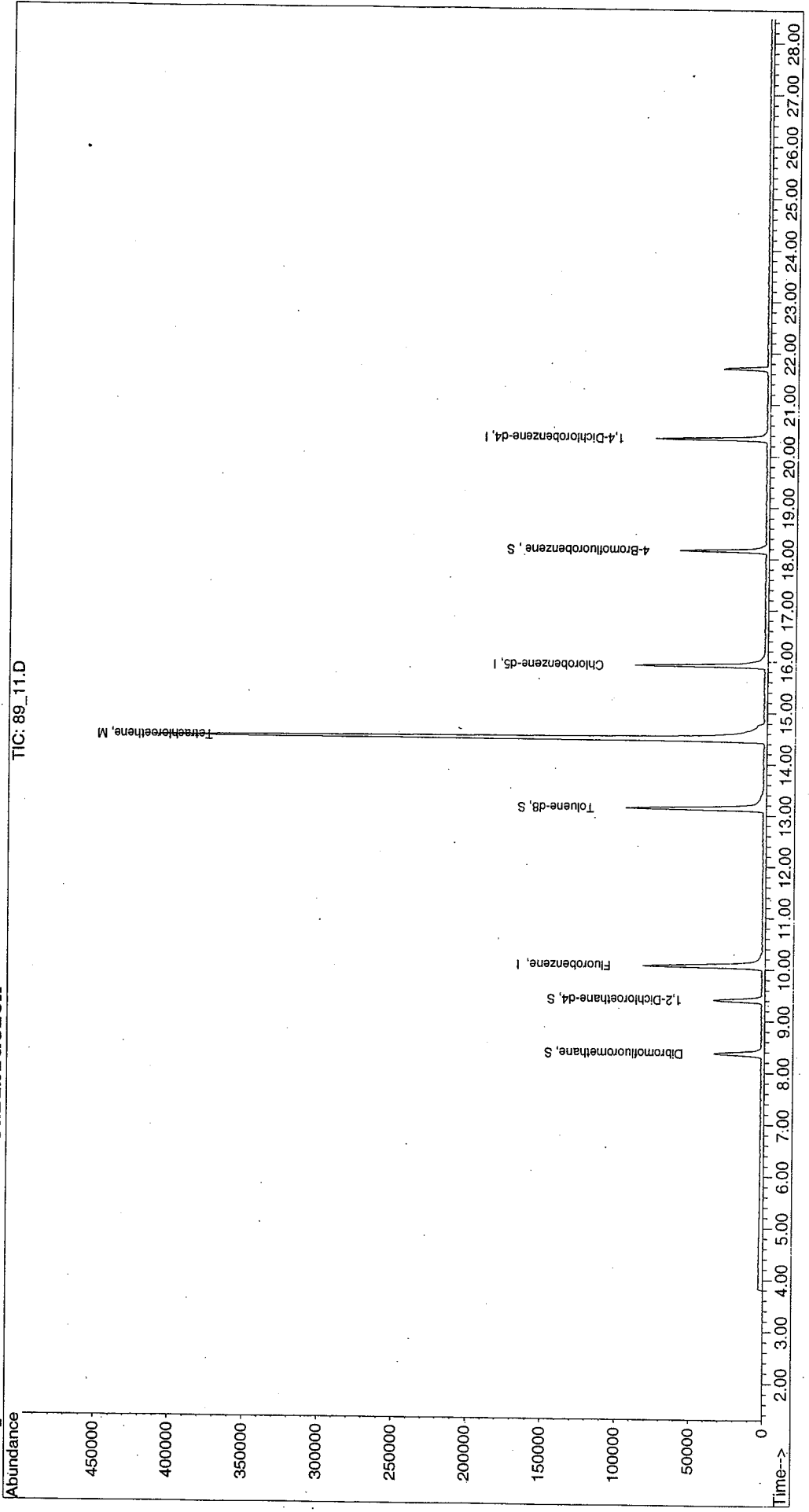
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|--------|------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 305794 | 183.55 | ug/L | 99 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Quantitation report

Data File : C:\HPCHEM\1\DATA\12_19_06\89_11.D Vial: 10
Acq On : 20 Dec 2006 12:42 am Operator: TS
Sample : 7489_11 0.40 ml Inst : GC/MS #2
Misc : Multiplr: 1.00
MS Integration Params: ODD.P
Quant Time: Dec 20 9:46 2006 Quant Results File: 71NV16_6.RE5

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Nov 27 09:29:43 2006
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\12_20_06\7489_11.D
 Acq On : 20 Dec 2006 5:15 pm
 Sample : 7489_11 0.20 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Dec 26 12:34 2006

Vial: 3
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Thu Nov 16 17:01:55 2006
 Response via : Initial Calibration
 DataAcq Meth : RUN1

Internal Standards R.T. QIon Response Conc Units Dev(Min)

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.06 | 96 | 191779 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.92 | 117 | 122777 | 25.00 | ug/L | -0.03 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 45496 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|-------|
| 16) Dibromofluoromethane | 8.37 | 113 | 49792 | 26.86 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 107.44% | |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 39754 | 24.77 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 99.08% | |
| 28) Toluene-d8 | 13.14 | 98 | 159867 | 24.87 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 99.48% | |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 48597 | 21.36 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 85.44% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\12_20_06\7489_11.D
 Acq On : 20 Dec 2006 5:15 pm
 Sample : 7489_11 0.20 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Dec 26 12:34 2006

Vial: 3
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Thu Nov 16 17:01:55 2006
 Response via : Initial Calibration
 DataAcq Meth : RUN1

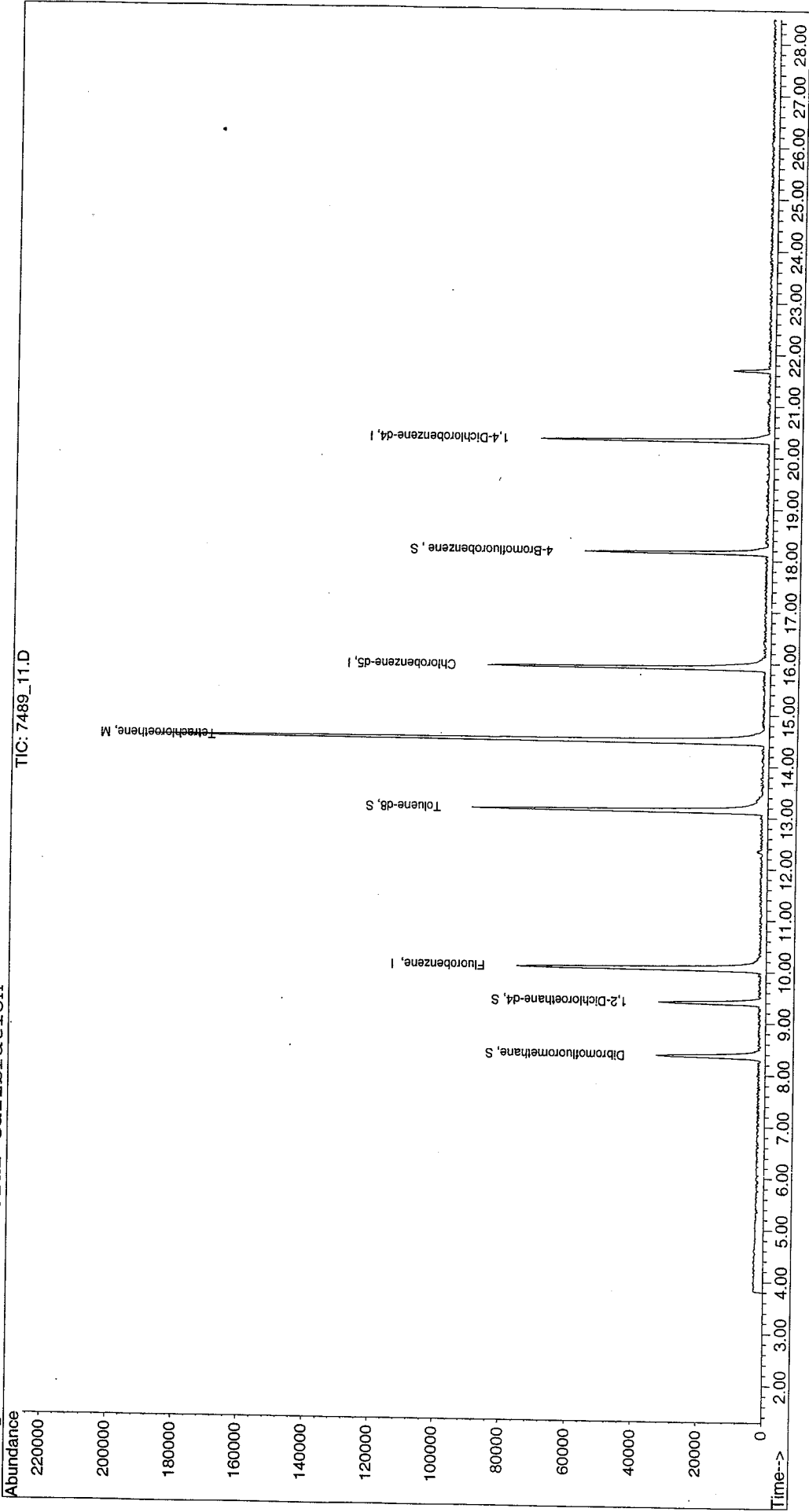
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 134291 | 86.75 | ug/L | 97 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Quantitation report

Data File : C:\HPCHEM\1\DATA\12_20_06\7489_11.D Vial: 3
Acq On : 20 Dec 2006 5:15 pm Operator: TS
Sample : 7489_11 0.20 ml Inst : GC/MS #2
Misc : Multiplr: 1.00
MS Integration Params: ODD.P
Quant Time: Dec 26 12:34 2006 Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Nov 27 09:29:43 2006
Response via : Initial Calibration



Acq On : 20 Dec 2006 1:16 am

Operator: TS

Sample : 7489_12 0.40 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Dec 20 9:47 2006

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|--------------------|------|------|----------|------|-------|----------|
|--------------------|------|------|----------|------|-------|----------|

| | | | | | | |
|----------------------------|-------|-----|--------|-------|------|-------|
| 1) Fluorobenzene | 10.06 | 96 | 199406 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.91 | 117 | 122946 | 25.00 | ug/L | -0.04 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 47826 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|--------|-------|------|--------------------|
| 16) Dibromofluoromethane | 8.37 | 113 | 51479 | 26.70 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 106.80% |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 41876 | 25.10 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 100.40% |
| 28) Toluene-d8 | 13.13 | 98 | 161919 | 24.23 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 96.92% |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 46205 | 20.28 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 81.12% |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\12_19_06\89_12.D
Acq On : 20 Dec 2006 1:16 am
Sample : 7489_12 0.40 ml
Misc :

Vial: 11
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00

MS Integration Params: ODD.P
Quant Time: Dec 20 9:47 2006

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Thu Nov 16 17:01:55 2006
Response via : Initial Calibration
DataAcq Meth : RUN1

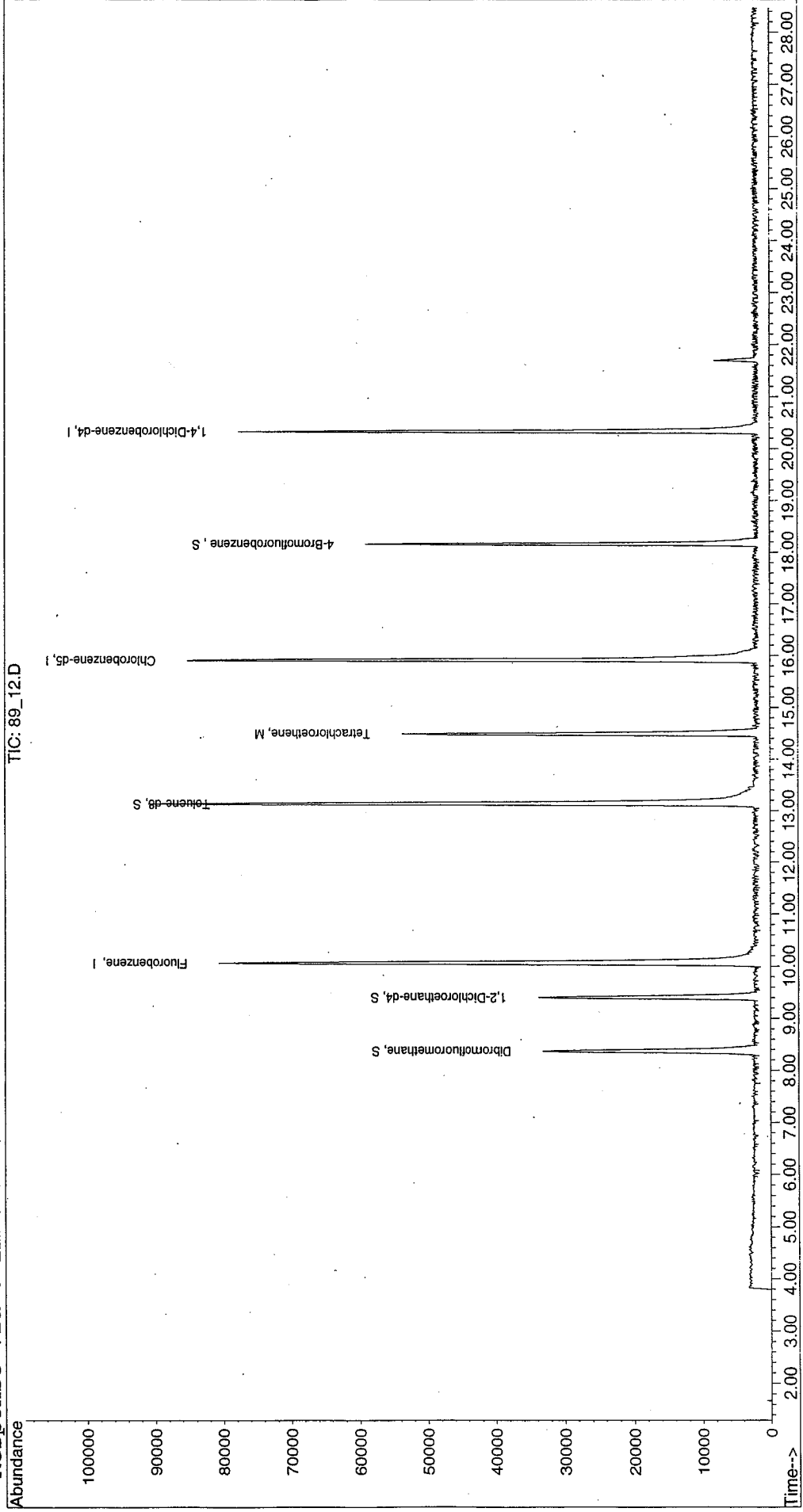
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|--------|--------|
| 32) Tetrachloroethene | 14.48 | 166 | 37810 | 23.49 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | | N.D. | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | | N.D. d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | | N.D. | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | | N.D. | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | | N.D. | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | | N.D. | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | | N.D. | |
| 42) Styrene | 0.00 | 104 | 0 | | N.D. | |
| 43) Bromoform | 0.00 | 173 | 0 | | N.D. | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 46) Bromobenzene | 0.00 | 156 | 0 | | N.D. | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | | N.D. | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | | N.D. | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | | N.D. | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | | N.D. | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | | N.D. | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | | N.D. | |
| 65) Naphthalene | 0.00 | 128 | 0 | | N.D. | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | | N.D. | |
| 68) Acetone | 0.00 | 43 | 0 | | N.D. | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | | N.D. | |
| 70) 2-Butanone | 0.00 | 43 | 0 | | N.D. | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | | N.D. | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | | N.D. | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | | N.D. | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\12_19_06\89_12.D
Acq On : 20 Dec 2006 1:16 am
Sample : 7489_12 0.40 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Dec 20 9:47 2006
Vial: 11
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Nov 27 09:29:43 2006
Response via : Initial Calibration





17 Princess Rd
Lawrenceville, New Jersey 08648
Tel: 609/895-5370
Fax: 609/895-1858

Volatile Organic Compound Data Summary Package

Prepared for
Gorham Textron

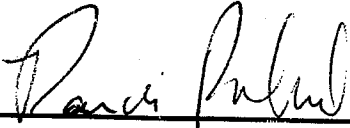
Lab ID
7502

Project Number 101960 02000000

Samples Received
27-Jan-06

Reported
23-Jan-07

NJDEP Certified Lab 11001


Randi K Rothmel, PhD

1/23/07
Date

Laboratory Director

1.0 Chain of Custody

CHAIN OF CUSTODY

Ref. Document #

Project Number/Cost code: 101960 / 02

Project Name / Location: Garham Textron

Send bill to:

Project Contact: Charles Schaefer, Chuck Condee
(Name & phone #)

Send Report To: Chuck Condee
Phone/Fax Number:
Address:
City/State:

Shipment date:
Lab Destination:
Lab Contact Name / ph. #:

Sampler's Name(s): Sheryl Stregger

| Lab No. | Sample ID Number | Sample Description | Collection Information | | Matrix | # of containers | Container type | Preservative | | | | VOC | MEE | Turn Around Time Requested | |
|---------|------------------|--------------------|------------------------|------|--------|-----------------|----------------|--------------|-----|------|------------------|-----|-----|----------------------------|--------------------------------|
| | | | Date | Time | | | | G/C | HCL | NaOH | HNO ₃ | | | | H ₂ SO ₄ |
| 1 | Killed Control A | | 12/27/06 | 7:00 | g | Aq | 1 | 10mL | X | | | | | X | Std |
| 2 | Killed Control B | | 12/27/06 | 7:00 | g | Aq | 1 | 5mL | | | | | X | | Std |
| 3 | Killed Control C | | 12/27/06 | 7:00 | g | Aq | 1 | 10mL | X | | | | | X | Std |
| 4 | Live Control A | | 12/27/06 | 7:00 | g | Aq | 1 | 5mL | | | | | X | | Std |
| 5 | Live Control B | | 12/27/06 | 7:00 | g | Aq | 1 | 10mL | X | | | | | X | Std |
| 6 | Live Control C | | 12/27/06 | 7:00 | g | Aq | 1 | 5mL | | | | | X | | Std |
| 7 | Bloaug 1 L A | | 12/27/06 | 7:00 | g | Aq | 1 | 10mL | X | | | | | X | Std |
| 8 | Bloaug 1 L B | | 12/27/06 | 7:00 | g | Aq | 1 | 5mL | | | | | X | | Std |
| 9 | Bloaug 1 L C | | 12/27/06 | 7:00 | g | Aq | 1 | 10mL | X | | | | | X | Std |
| 10 | Bloaug 2 E A | | 12/27/06 | 7:00 | g | Aq | 1 | 5mL | | | | | X | | Std |
| 11 | Bloaug 2 E B | | 12/27/06 | 7:00 | g | Aq | 1 | 10mL | X | | | | | X | Std |

* Note: PCE ~ 30 PPM!

Special Instructions:
RCRA PCB/dioxin PAH/oll
QC/Data Package Level Required: I II III IV NJ EDD GIS EDD Preliminary data
Known Waste Stream Circle: RAD Corrosive Flammable Reactive
C = Composite G = Grab

Relinquished By: *Sheryl Stregger* Date: 12-27-06 Time: 1027
Relinquished By: *[Signature]* Date: 12/27/06 Time: 1230
Date: Date: Time: Time:

Level II = data summary + basic QC
Level III = New Jersey QC reduced deliverable
Level IV = Full deliverable CLP package
Cooler temperature upon arrival at Lab:



COC Continuation Page

COC Ref. Document #

Page 2 of 2

762

Project Number/ cost code: 101960 / 02 Shipment Date:

Project Name / Location: Gorham, Textron

| Lab No. | Sample ID Number | Sample Description | Collection Information | | # of containers | Container Type | Preservative | | | | | VOC | MEE | Anions | VFA | Turn Around Time Requested | |
|---------|-------------------|--------------------|------------------------|------|-----------------|----------------|--------------|------|------|------------------|--------------------------------|-----|-----|--------|-----|----------------------------|------|
| | | | Date | Time | | | G/C | HCL | NaOH | HNO ₃ | H ₂ SO ₄ | | | | | | None |
| 12 | Bioaug 2 E C " | | 12/27/06 | 7:00 | g | Aq | 1 | 10mL | X | | | | | | | | |
| 13 | Killed Control D | | 12/27/06 | 7:00 | g | Aq | 1 | 5mL | | X | X | | | | | | |
| 14 | Live Control D | | 12/27/06 | 7:00 | g | Aq | 1 | 15mL | | X | X | | | X | X | | |
| 15 | Bioaug 1 L D | | 12/27/06 | 7:00 | g | Aq | 1 | 15mL | | X | X | | | X | X | | |
| 16 | Bioaug 2 E D | | 12/27/06 | 7:00 | g | Aq | 1 | 15mL | | X | X | | | X | X | | |
| | | | | | | | | | | X | X | | | X | X | | |

Analyses Requested

2.0. Sample Summaries

Sample summaries are enclosed

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-01

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-01 0.050 ml
 Sample wt/vol: 0.1 ^{0.05} (g/ml) ML Lab File ID: 7502_01.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 18000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-01

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-01 0.050 ml
 Sample wt/vol: 0.1 A? 0.05 (g/ml) ML Lab File ID: 7502_01.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-02

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-02 0.050 ml
 Sample wt/vol: 0.170.05 (g/ml) ML Lab File ID: 7502_02.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 16000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-02

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-02 0.050 ml
 Sample wt/vol: 0.1¹⁰⁰ 0.05 (g/ml) ML Lab File ID: 7502_02.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | <u>UG/L</u> | Q |
|-----------|-----------------------------|-----------------|-------------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-03

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-03 0.050 ml
 Sample wt/vol: 0.05 (g/ml) ML Lab File ID: 7502_03.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 16000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-03

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-03 0.050 ml
 Sample wt/vol: 0.05 (g/ml) ML Lab File ID: 7502_03.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-04

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-04 0.050 ml
 Sample wt/vol: ~~0.1~~ 0.05 (g/ml) ML Lab File ID: 7502_04.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 7300 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-04

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-04 0.050 ml
 Sample wt/vol: 0.170.03 (g/ml) ML Lab File ID: 7502_04.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-05

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-05 0.050 ml
 Sample wt/vol: 0.2705 (g/ml) ML Lab File ID: 7502_05.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 14000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-05

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-05 0.050 ml
 Sample wt/vol: 0.1 * 0.05 (g/ml) ML Lab File ID: 7502_05.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-06

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-06 0.050 ml
 Sample wt/vol: 0.105 (g/ml) ML Lab File ID: 7502_06.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 14000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-06

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-06 0.050 ml
 Sample wt/vol: 0.1^g 0.05^g (g/ml) ML Lab File ID: 7502_06.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-07

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-07 0.050 ml
 Sample wt/vol: 0.1870.05 (g/ml) ML Lab File ID: 7502_07.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 16000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-07

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-07 0.050 ml
 Sample wt/vol: 0.1 0.05 (g/ml) ML Lab File ID: 7502_07.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-08

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-08 0.050 ml
 Sample wt/vol: 0.170 (g/ml) ML Lab File ID: 7502_08.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 16000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-08

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-08 0.050 ml
 Sample wt/vol: 0.050 (g/ml) ML Lab File ID: 7502_08.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-9

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-09 0.050 ml
 Sample wt/vol: 0.1205 (g/ml) ML Lab File ID: 7502_9.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 9800 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-9

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-09 0.050 ml
 Sample wt/vol: 0.127 0.050 (g/ml) ML Lab File ID: 7502_9.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/02/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7502-10

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-10 0.10 ml
 Sample wt/vol: 0.1 (g/ml) ML Lab File ID: 02_10.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/17/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 100.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | | 500 | U |
| 74-87-3 | Chloromethane | | 500 | U |
| 75-01-4 | Vinyl chloride | | 500 | U |
| 74-83-9 | Bromomethane | | 500 | U |
| 75-00-3 | Chloroethane | | 500 | U |
| 75-69-4 | Trichlorofluoromethane | | 500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 500 | U |
| 75-09-2 | Methylene chloride | | 500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 500 | U |
| 74-97-5 | Bromochloromethane | | 500 | U |
| 67-66-3 | Chloroform | | 500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 500 | U |
| 56-23-5 | Carbon tetrachloride | | 500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 500 | U |
| 71-43-2 | Benzene | | 500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 500 | U |
| 79-01-6 | Trichloroethene | | 500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 500 | U |
| 74-95-3 | Dibromomethane | | 500 | U |
| 75-27-4 | Bromodichloromethane | | 500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 500 | U |
| 108-88-3 | Toluene | | 500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 500 | U |
| 127-18-4 | Tetrachloroethene | | 2500 | D |
| 142-28-9 | 1,3-Dichloropropane | | 500 | U |
| 124-48-1 | Dibromochloromethane | | 500 | U |
| 108-90-7 | Chlorobenzene | | 500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 500 | U |
| 100-41-4 | Ethylbenzene | | 500 | U |
| 1330-20-7 | Xylene (para & meta) | | 500 | U |
| 95-47-6 | Xylene (Ortho) | | 500 | U |
| 100-42-5 | Styrene | | 500 | U |
| 75-25-2 | Bromoform | | 500 | U |
| 98-82-8 | Isopropylbenzene | | 500 | U |
| 108-86-1 | Bromobenzene | | 500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7502-10

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-10 0.10 ml
 Sample wt/vol: 0.1 (g/ml) ML Lab File ID: 02_10.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/17/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 100.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 500 | U |
| 103-65-1 | n-Propylbenzene | | 500 | U |
| 95-49-8 | 2-Chlorotoluene | | 500 | U |
| 106-43-4 | 4-Chlorotoluene | | 500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 500 | U |
| 98-06-6 | tert-Butylbenzene | | 500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 500 | U |
| 135-98-8 | sec-Butylbenzene | | 500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 500 | U |
| 104-51-8 | n-Butylbenzene | | 500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 500 | U |
| 87-68-3 | Hexachlorobutadiene | | 500 | U |
| 91-20-3 | Naphthalene | | 500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 500 | U |
| 1634-04-4 | MTBE | | 500 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 500 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-11

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-11 0.050 ml
 Sample wt/vol: 0.95 (g/ml) ML Lab File ID: 7502_11.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/03/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 14000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-11

Lab Name: ATL Analyst: Tony Soto

NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA

Matrix: (soil/water) WATER Lab Sample ID: 7502-11 0.050 ml

Sample wt/vol: 0.050 (g/ml) ML Lab File ID: 7502_11.D

Level: (low/med) LOW Date Received: 12/27/06

% Moisture: not dec. _____ Date Analyzed: 01/03/07

GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0

Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-12

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-12 0.050 ml
 Sample wt/vol: 0.120.050 (g/ml) ML Lab File ID: 7502_12.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. Date Analyzed: 01/03/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND. | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 5100 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7502-12

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7502-12 0.050 ml
 Sample wt/vol: 0.050 (g/ml) ML Lab File ID: 7502_12.D
 Level: (low/med) LOW Date Received: 12/27/06
 % Moisture: not dec. _____ Date Analyzed: 01/03/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

3.0 Raw Data and Chromatograms

Raw data and Chromatograms are attached.

Acq On : 2 Jan 2007 6:30 pm

Sample : 7502-01 0.050 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Jan 3 8:32 2007

Vial: 3

Operator: TS

Inst : GC/MS #2

Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.07 | 96 | 218524 | 25.00 | ug/L | -0.03 |
| 36) Chlorobenzene-d5 | 15.92 | 117 | 146184 | 25.00 | ug/L | -0.03 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 60713 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|-------|
| 16) Dibromofluoromethane | 8.38 | 113 | 57357 | 27.15 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 108.60% | |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 46174 | 25.25 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 101.00% | |
| 28) Toluene-d8 | 13.13 | 98 | 175412 | 23.95 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 95.80% | |
| 45) 4-Bromofluorobenzene | 18.17 | 95 | 58933 | 21.75 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 87.00% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#)= qualifier out of range (m) = manual integration

7502_01.D 71NV16_6.M

Wed Jan 17 11:13:30 2007

GC#2

Page 1

Data File : C:\HPCHEM\1\DATA\010207\7502_01.D

Vial: 3

Acq On : 2 Jan 2007 6:30 pm

Operator: TS

Sample : 7502-01 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 3 8:32 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 37194 | 21.09 | ug/L | 99 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | d | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7502_01.D 71NV16_6.M

Wed Jan 17 11:13:30 2007

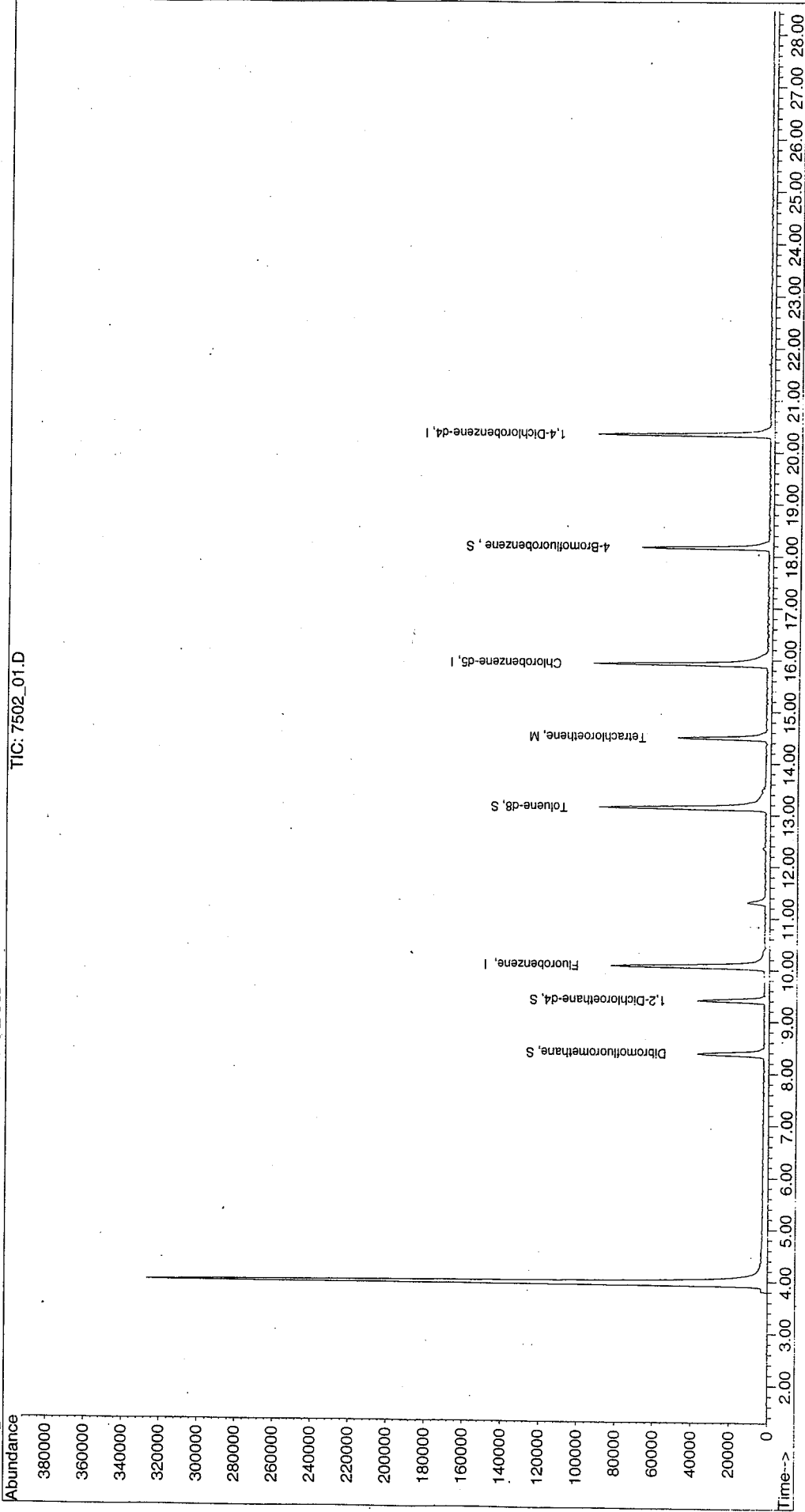
GC#2

Page 2

Quantitation Report

Data File : C:\HPCHEM\1\DATA\010207\7502_01.D
Acq On : 2 Jan 2007 6:30 pm
Sample : 7502-01 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 3 8:32 2007
Vial: 3
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00
Quant Results File: 71NV16_6.REB

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Acq On : 2 Jan 2007 7:05 pm

Operator: TS

Sample : 7502-02 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 3 8:33 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.06 | 96 | 209059 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.92 | 117 | 139439 | 25.00 | ug/L | -0.03 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 59133 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|--------|-------|------|--------------------|
| 16) Dibromofluoromethane | 8.37 | 113 | 59228 | 29.31 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 117.24% |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 46944 | 26.84 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 107.36% |
| 28) Toluene-d8 | 13.13 | 98 | 172809 | 24.66 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 98.64% |
| 45) 4-Bromofluorobenzene | 18.17 | 95 | 58042 | 22.46 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 89.84% |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. | d |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#)= qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\010207\7502_02.D

Acq On : 2 Jan 2007 7:05 pm

Sample : 7502-02 0.050 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Jan 3 8:33 2007

Vial: 4

Operator: TS

Inst : GC/MS #2

Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|--------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 32592 | 19.31 | ug/L | 96 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | | N.D. | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | | N.D. d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | | N.D. | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | | N.D. | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | | N.D. | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | | N.D. | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | | N.D. | |
| 42) Styrene | 0.00 | 104 | 0 | | N.D. | |
| 43) Bromoform | 0.00 | 173 | 0 | | N.D. | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 46) Bromobenzene | 0.00 | 156 | 0 | | N.D. | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | | N.D. | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | | N.D. | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | | N.D. | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | | N.D. | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | | N.D. | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | | N.D. | |
| 65) Naphthalene | 0.00 | 128 | 0 | | N.D. | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | | N.D. | |
| 68) Acetone | 0.00 | 43 | 0 | | N.D. | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | | N.D. d | |
| 70) 2-Butanone | 0.00 | 43 | 0 | | N.D. | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | | N.D. | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | | N.D. | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | | N.D. | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

7502_02.D 71NV16_6.M

Wed Jan 17 11:13:33 2007

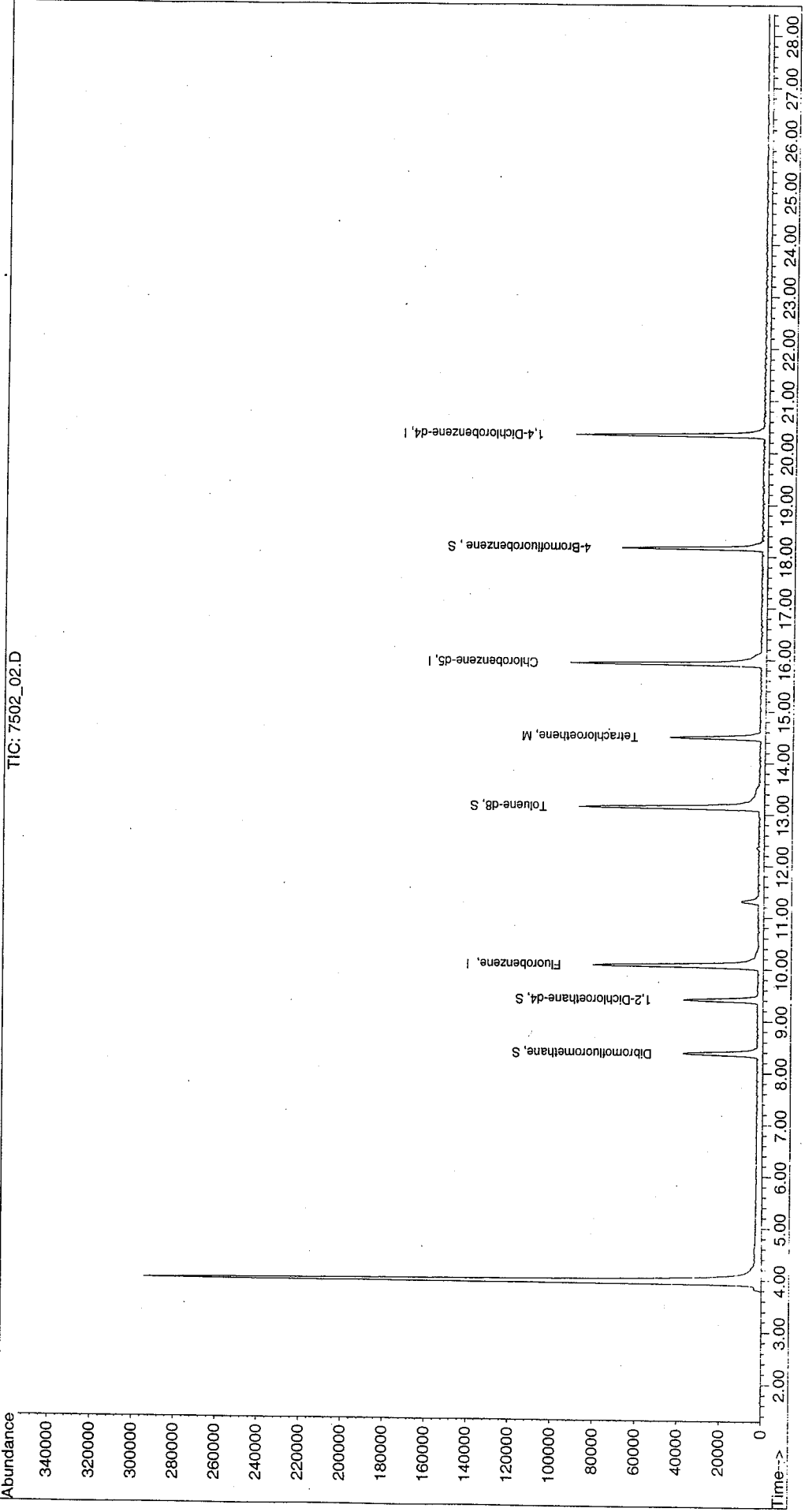
GC#2

Page 2

Quantitation Report

Data File : C:\HPCHEM\1\DATA\010207\7502_02.D
Acq On : 2 Jan 2007 7:05 pm
Sample : 7502-02 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 3 8:33 2007
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\010207\7502_03.D

Acq On : 2 Jan 2007 7:40 pm

Sample : 7502-03 0.050 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Jan 3 8:34 2007

Vial: 5

Operator: TS

Inst : GC/MS #2

Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.06 | 96 | 209415 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.92 | 117 | 148010 | 25.00 | ug/L | -0.03 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 58941 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|-------|
| 16) Dibromofluoromethane | 8.38 | 113 | 57631 | 28.47 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | | | | |
| | | | Recovery | = | 113.88% | |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 46754 | 26.68 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | Recovery | = | 106.72% | |
| 28) Toluene-d8 | 13.14 | 98 | 183248 | 26.11 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | Recovery | = | 104.44% | |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 58320 | 21.26 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | Recovery | = | 85.04% | |

Target Compounds

| | | | | | Qvalue |
|-------------------------------|------|-----|---|------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | |

(#) = qualifier out of range (m) = manual integration

7502_03.D 71NV16_6.M

Wed Jan 17 11:13:36 2007

GC#2

Page 1

Data File : C:\HPCHEM\1\DATA\010207\7502_03.D

Vial: 5

Acq On : 2 Jan 2007 7:40 pm

Operator: TS

Sample : 7502-03 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 3 8:34 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 33169 | 19.62 | ug/L | 95 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | d | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7502_03.D 71NV16_6.M

Wed Jan 17 11:13:36 2007

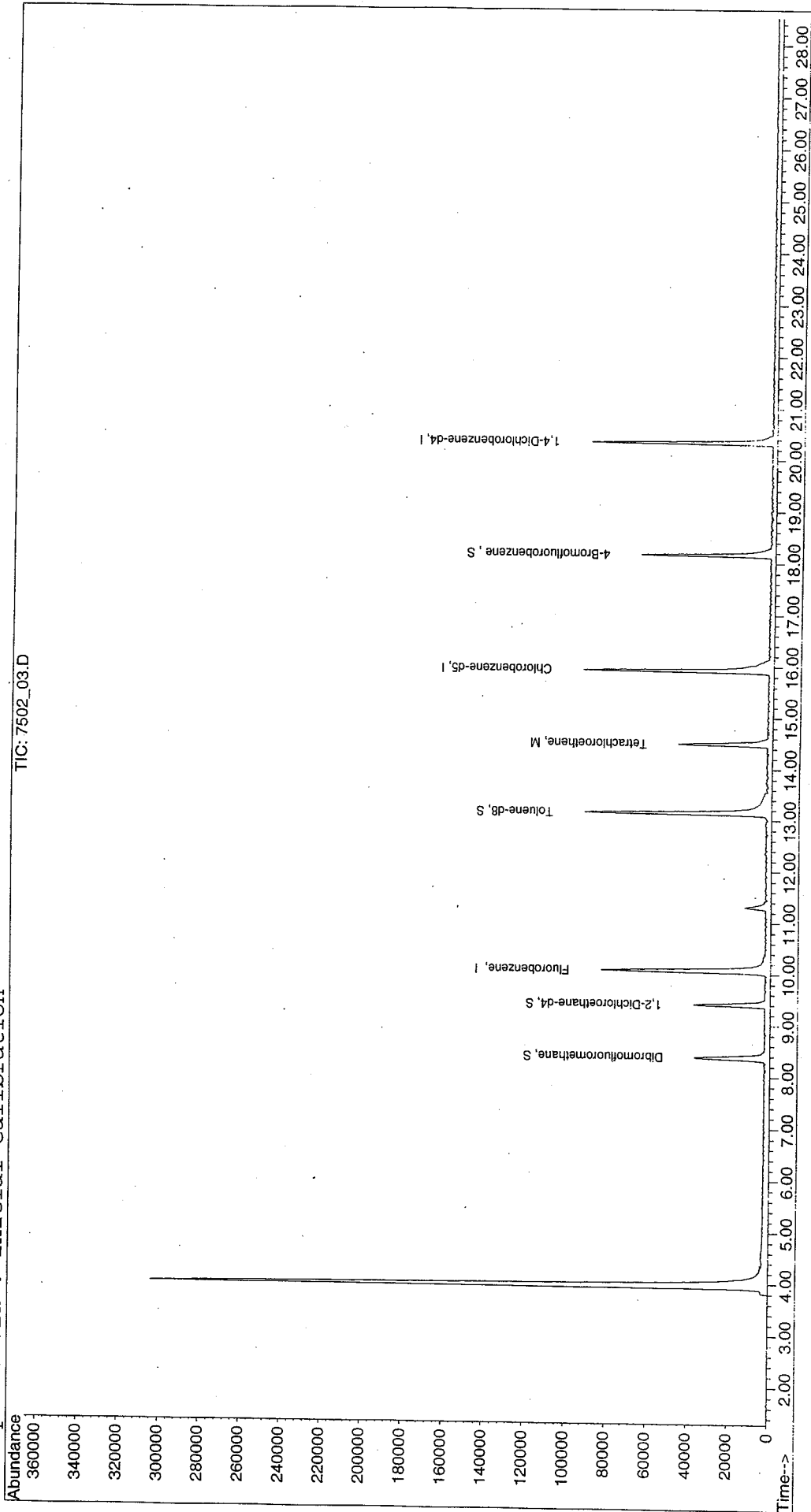
GC#2

Page 2

Quantitation Report

Data File : C:\HPCHEM\1\DATA\010207\7502_03.D
Acq On : 2 Jan 2007 7:40 pm Vial: 5
Sample : 7502-03 0.050 ml Operator: TS
Misc : Inst : GC/MS #2
MS Integration Params: ODD.P Multiplr: 1.00
Quant Time: Jan 3 8:34 2007 Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\010207\7502_04.D

Acq On : 2 Jan 2007 8:14 pm

Sample : 7502-04 0.050 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Jan 3 8:34 2007

Vial: 6

Operator: TS

Inst : GC/MS #2

Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev (Min) |
|-----------------------------|--------|------|----------|-------|---------|-----------|
| 1) Fluorobenzene | 10.06 | 96 | 208932 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.92 | 117 | 139335 | 25.00 | ug/L | -0.03 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 55894 | 25.00 | ug/L | -0.03 |
| System Monitoring Compounds | | | | | | |
| 16) Dibromofluoromethane | 8.37 | 113 | 57454 | 28.45 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 113.80% | |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 45801 | 26.20 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 104.80% | |
| 28) Toluene-d8 | 13.13 | 98 | 182086 | 26.00 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 104.00% | |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 57022 | 22.08 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 88.32% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\010207\7502_04.D

Vial: 6

Acq On : 2 Jan 2007 8:14 pm

Operator: TS

Sample : 7502-04 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 3 8:34 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|------|------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 14735 | 8.74 | ug/L | 99 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7502_04.D 71NV16_6.M

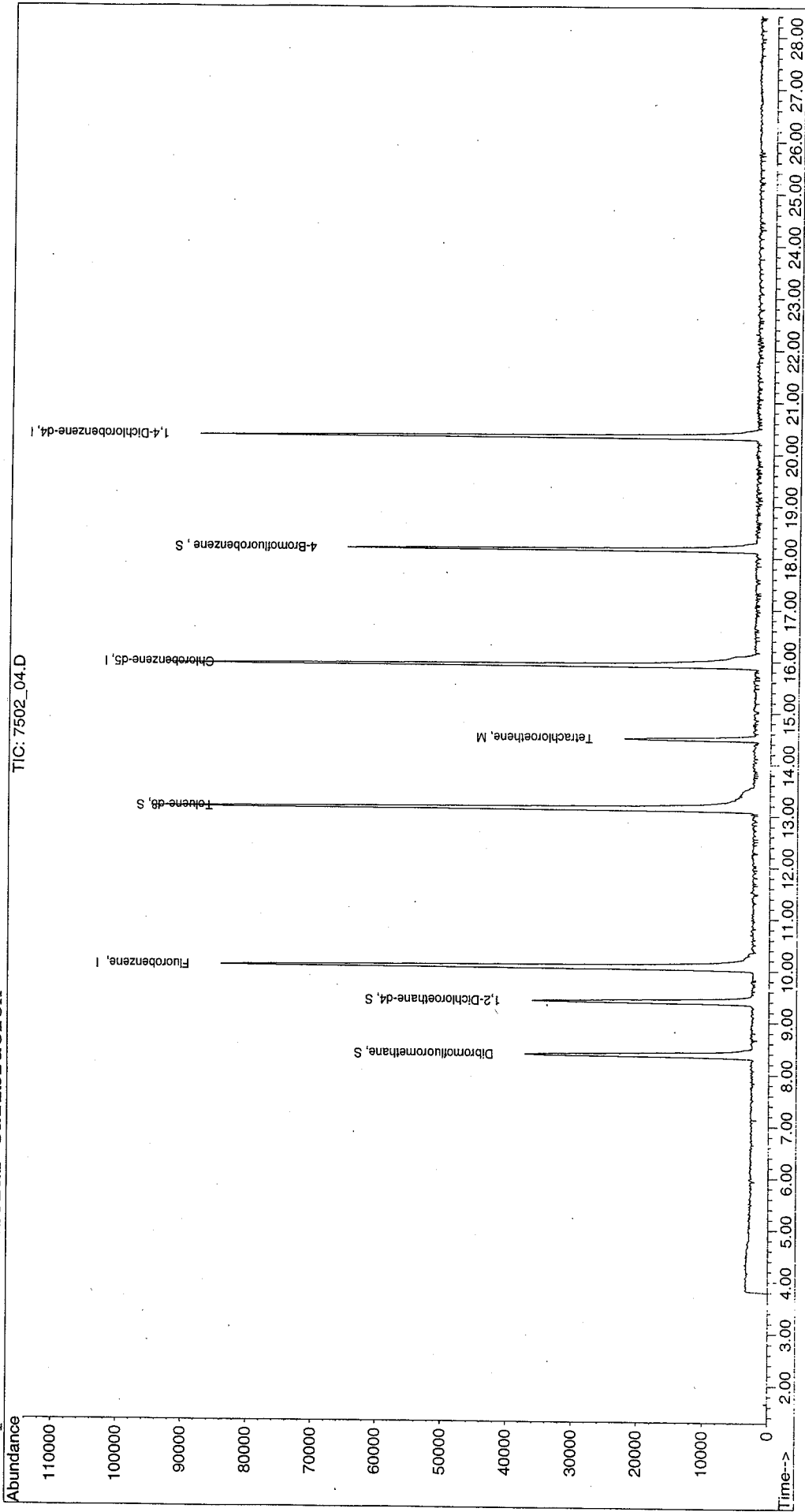
Wed Jan 17 11:13:39 2007

GC#2

Page 2

Quantitation Report

Data File : C:\HPCHEM\1\DATA\010207\7502_04.D
Acq On : 2 Jan 2007 8:14 pm Vial: 6
Sample : 7502-04 0.050 ml Operator: TS
Misc : Inst : GC/MS #2
MS Integration Params: ODD.P Multiplr: 1.00
Quant Time: Jan 3 8:34 2007 Quant Results File: 71NV16_6.RES
Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\010207\7502_05.D

Acq On : 2 Jan 2007 8:49 pm

Sample : 7502-05 0.050 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Jan 3 8:35 2007

Vial: 7

Operator: TS

Inst : GC/MS #2

Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.06 | 96 | 205610 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.91 | 117 | 135760 | 25.00 | ug/L | -0.04 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 57608 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|-------|
| 16) Dibromofluoromethane | 8.37 | 113 | 57296 | 28.83 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | Recovery | = | 115.32% | |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 46198 | 26.85 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | Recovery | = | 107.40% | |
| 28) Toluene-d8 | 13.14 | 98 | 162533 | 23.59 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | Recovery | = | 94.36% | |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 56388 | 22.41 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | Recovery | = | 89.64% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. | d |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

7502_05.D 71NV16_6.M

Wed Jan 17 11:13:42 2007

GC#2

Page 1

Data File : C:\HPCHEM\1\DATA\010207\7502_05.D

Vial: 7

Acq On : 2 Jan 2007 8:49 pm

Operator: TS

Sample : 7502-05 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 3 8:35 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 27698 | 16.69 | ug/L | 95 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7502_05.D 71NV16_6.M

Wed Jan 17 11:13:42 2007

GC#2

Page 2

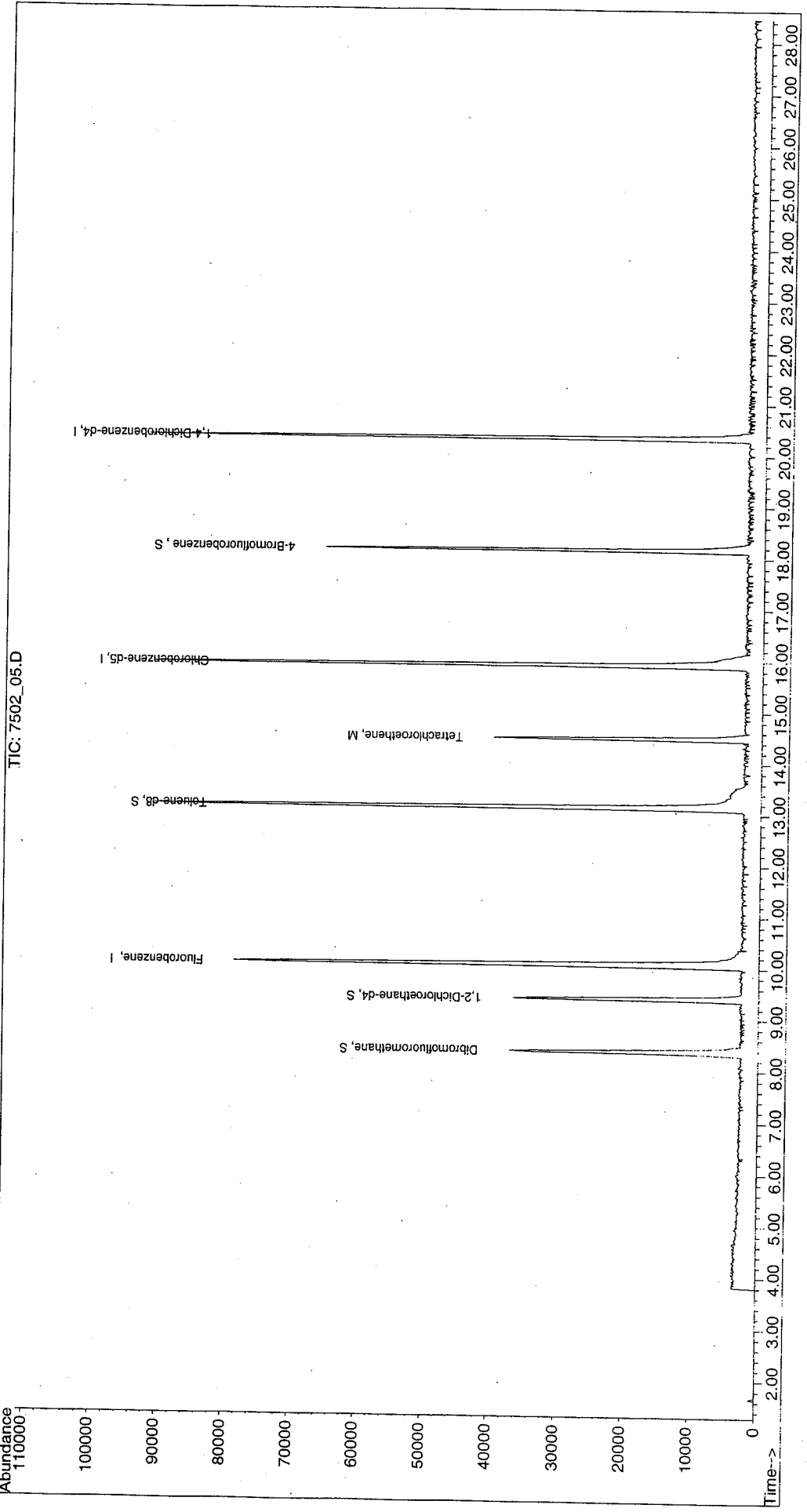
Quantitation Report

Data File : C:\HPCHEM\1\DATA\010207\7502_05.D
Acq On : 2 Jan 2007 8:49 pm
Sample : 7502-05 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 3 8:35 2007

Vial: 7
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Acq On : 2 Jan 2007 9:24 pm

Operator: TS

Sample : 7502-06 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 3 8:35 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.06 | 96 | 208220 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.92 | 117 | 134786 | 25.00 | ug/L | -0.03 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 56015 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|-------|
| 16) Dibromofluoromethane | 8.37 | 113 | 55492 | 27.57 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 110.28% | |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 44457 | 25.52 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 102.08% | |
| 28) Toluene-d8 | 13.13 | 98 | 164898 | 23.63 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 94.52% | |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 54881 | 21.97 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 87.88% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#)= qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\010207\7502_06.D

Vial: 8

Acq On : 2 Jan 2007 9:24 pm

Operator: TS

Sample : 7502-06 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 3 8:35 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|--------|--------|
| 32) Tetrachloroethene | 14.48 | 166 | 28391 | 16.89 | ug/L | 95 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | | N.D. | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | | N.D. d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | | N.D. | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | | N.D. | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | | N.D. | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | | N.D. | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | | N.D. | |
| 42) Styrene | 0.00 | 104 | 0 | | N.D. | |
| 43) Bromoform | 0.00 | 173 | 0 | | N.D. | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 46) Bromobenzene | 0.00 | 156 | 0 | | N.D. | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | | N.D. | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | | N.D. | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | | N.D. | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | | N.D. | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | | N.D. | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | | N.D. | |
| 65) Naphthalene | 0.00 | 128 | 0 | | N.D. | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | | N.D. | |
| 68) Acetone | 0.00 | 43 | 0 | | N.D. | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | | N.D. | |
| 70) 2-Butanone | 0.00 | 43 | 0 | | N.D. | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | | N.D. | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | | N.D. | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | | N.D. | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

7502_06.D 71NV16_6.M

Wed Jan 17 11:13:45 2007

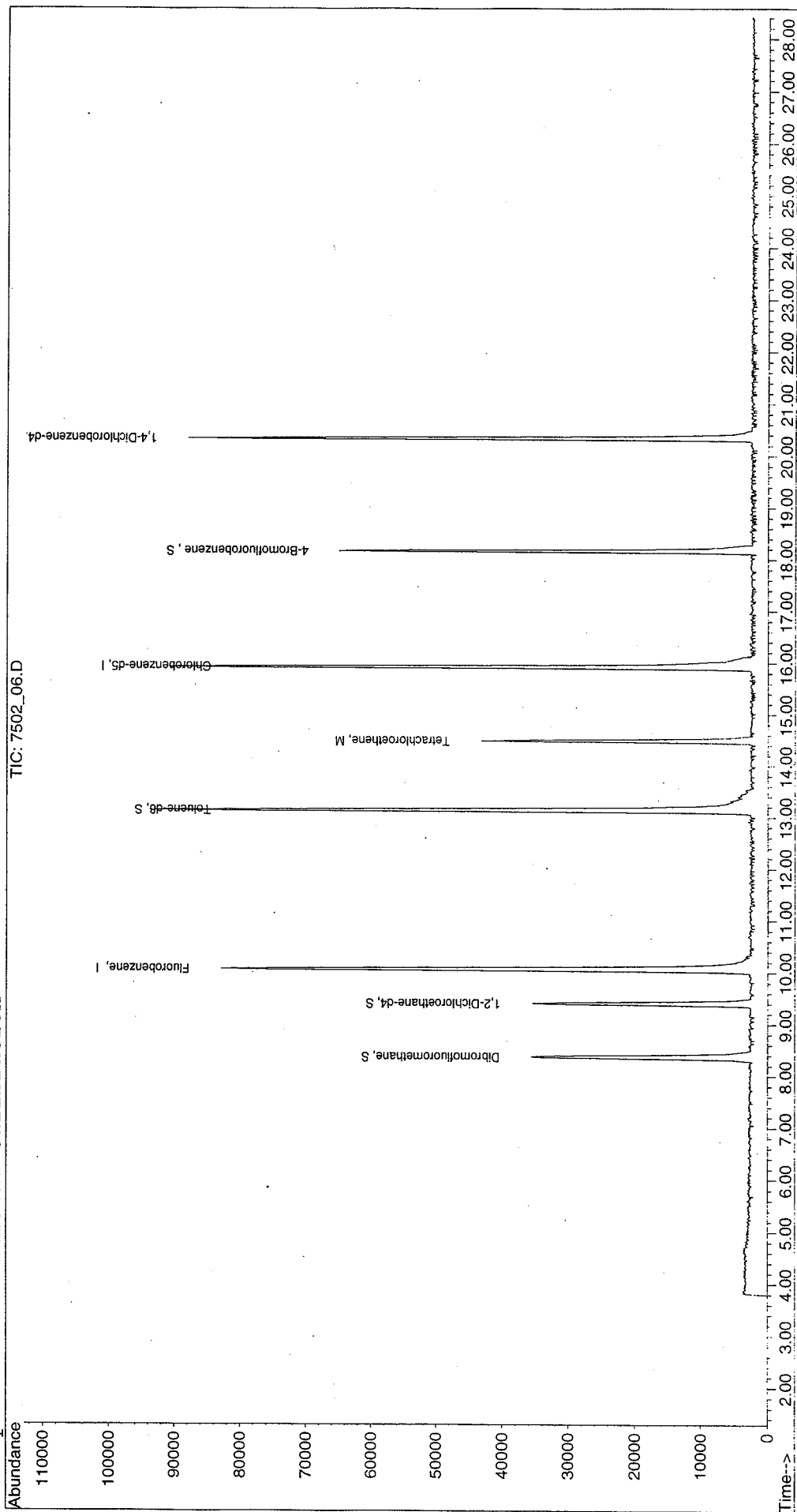
GC#2

Page 2

Quantitation Report

Data File : C:\HPCHEM\1\DATA\010207\7502_06.D
Acq On : 2 Jan 2007 9:24 pm
Sample : 7502-06 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 3 8:35 2007
Vial: 8
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\010207\7502_07.D

Vial: 9

Acq On : 2 Jan 2007 9:58 pm

Operator: TS

Sample : 7502-07 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 3 8:36 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.06 | 96 | 199536 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.92 | 117 | 130099 | 25.00 | ug/L | -0.03 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 57259 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|-------|
| 16) Dibromofluoromethane | 8.37 | 113 | 54549 | 28.28 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 113.12% | |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 43342 | 25.96 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 103.84% | |
| 28) Toluene-d8 | 13.14 | 98 | 158908 | 23.76 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 95.04% | |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 55236 | 22.91 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 91.64% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7502_07.D 71NV16_6.M

Wed Jan 17 11:13:48 2007

GC#2

Page 1

Acq On : 2 Jan 2007 9:58 pm

Operator: TS

Sample : 7502-07 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 3 8:36 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

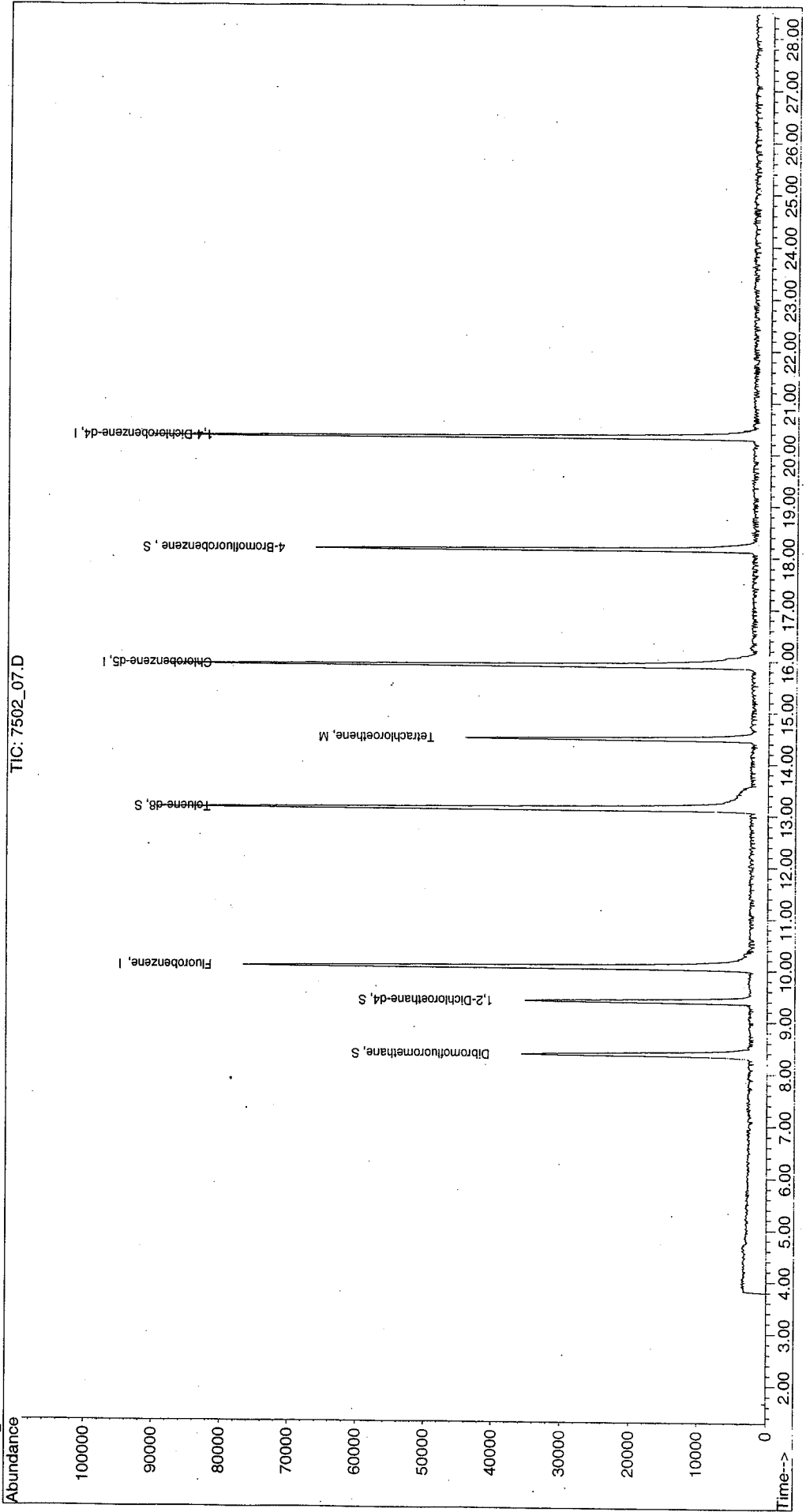
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|--------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 31328 | 19.45 | ug/L | 95 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | | N.D. | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | | N.D. d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | | N.D. | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | | N.D. | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | | N.D. | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | | N.D. | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | | N.D. | |
| 42) Styrene | 0.00 | 104 | 0 | | N.D. | |
| 43) Bromoform | 0.00 | 173 | 0 | | N.D. | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 46) Bromobenzene | 0.00 | 156 | 0 | | N.D. | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | | N.D. | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | | N.D. | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | | N.D. | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | | N.D. | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | | N.D. | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | | N.D. | |
| 65) Naphthalene | 0.00 | 128 | 0 | | N.D. | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | | N.D. | |
| 68) Acetone | 0.00 | 43 | 0 | | N.D. | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | | N.D. | |
| 70) 2-Butanone | 0.00 | 43 | 0 | | N.D. | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | | N.D. | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | | N.D. | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | | N.D. | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | | N.D. | |

 (#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\010207\7502_07.D
Acq On : 2 Jan 2007 9:58 pm
Sample : 7502-07 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 3 8:36 2007
Vial: 9
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\010207\7502_08.D

Vial: 10

Acq On : 2 Jan 2007 10:33 pm

Operator: TS

Sample : 7502-08 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 3 8:37 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.06 | 96 | 185826 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.91 | 117 | 119626 | 25.00 | ug/L | -0.04 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 54276 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|-------|
| 16) Dibromofluoromethane | 8.37 | 113 | 54583 | 30.38 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 121.52% | |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 43387 | 27.90 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 111.60% | |
| 28) Toluene-d8 | 13.13 | 98 | 146260 | 23.48 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 93.92% | |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 49225 | 22.21 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 88.84% | |

Target Compounds

Qvalue

| | | | | | |
|-------------------------------|------|-----|---|------|---|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | |

(#) = qualifier out of range (m) = manual integration

7502_08.D 71NV16_6.M

Wed Jan 17 11:13:51 2007

GC#2

Page 1

Data File : C:\HPCHEM\1\DATA\010207\7502_08.D

Acq On : 2 Jan 2007 10:33 pm

Sample : 7502-08 0.050 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Jan 3 8:37 2007

Vial: 10

Operator: TS

Inst : GC/MS #2

Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 29395 | 19.60 | ug/L | 96 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7502_08.D 71NV16_6.M

Wed Jan 17 11:13:51 2007

GC#2

Page 2

Quantitation Report

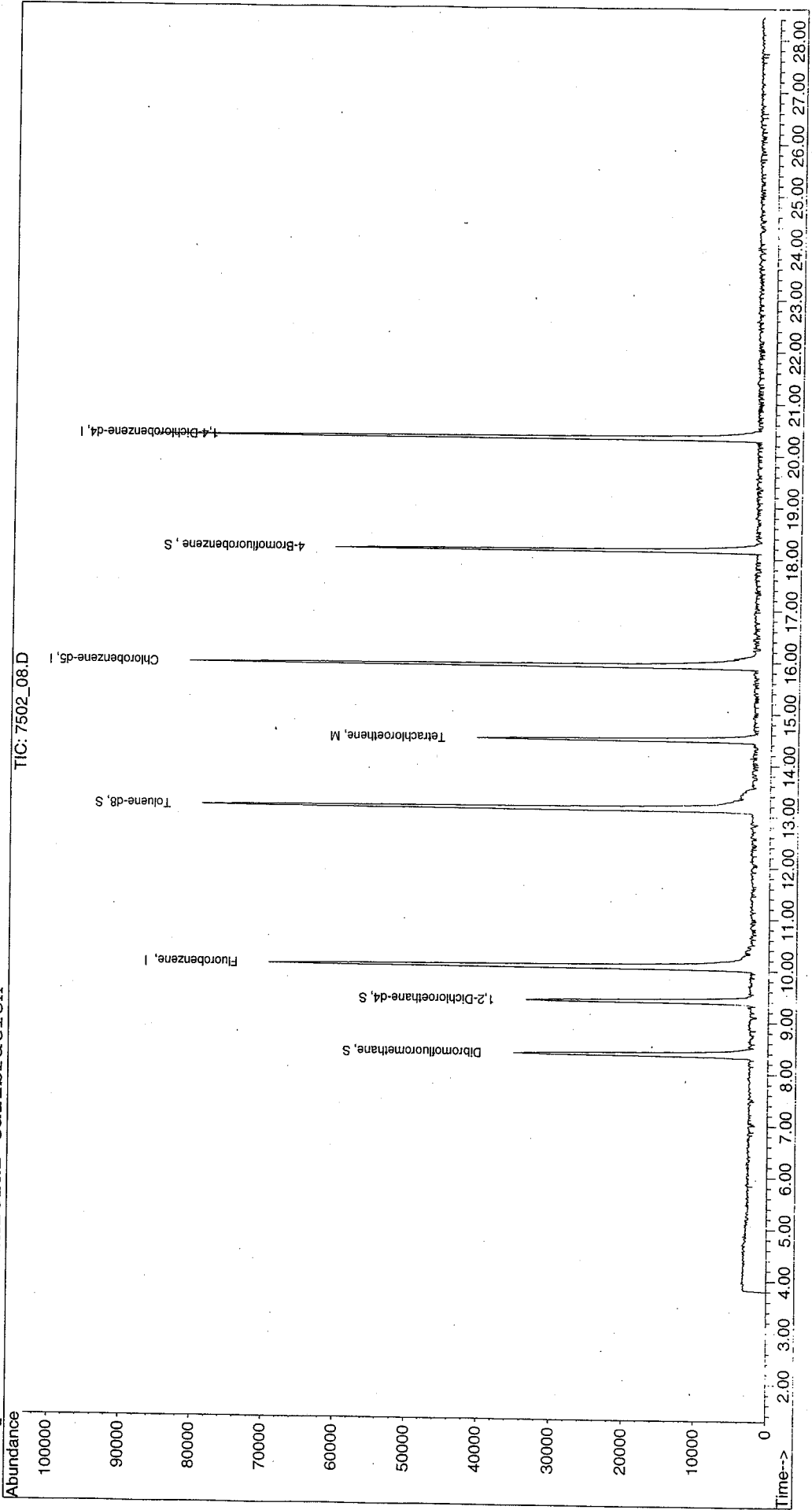
Data File : C:\HPCHEM\1\DATA\010207\7502_08.D
Acq On : 2 Jan 2007 10:33 pm
Sample : 7502-08 0.050 ml
Misc :

Vial: 10
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00

MS Integration Params: ODD.P
Quant Time: Jan 3 8:37 2007

Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\010207\7502_9.D

Acq On : 2 Jan 2007 11:08 pm

Sample : 7502-09 0.050 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Jan 3 8:43 2007

Vial: 11

Operator: TS

Inst : GC/MS #2

Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.06 | 96 | 196763 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.91 | 117 | 129394 | 25.00 | ug/L | -0.04 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 54222 | 25.00 | ug/L | -0.03 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|-------|
| 16) Dibromofluoromethane | 8.37 | 113 | 53206 | 27.97 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 111.88% | |
| 19) 1,2-Dichloroethane-d4 | 9.40 | 65 | 42493 | 25.81 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 103.24% | |
| 28) Toluene-d8 | 13.14 | 98 | 161752 | 24.53 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 98.12% | |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 54551 | 22.75 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 91.00% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7502_9.D 71NV16_6.M

Wed Jan 17 11:14:03 2007

GC#2

Page 1

Data File : C:\HPCHEM\1\DATA\O10207\7502_9.D

Acq On : 2 Jan 2007 11:08 pm

Sample : 7502-09 0.050 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Jan 3 8:43 2007

Vial: 11

Operator: TS

Inst : GC/MS #2

Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|--------|--------|
| 32) Tetrachloroethene | 14.48 | 166 | 18469 | 11.63 | ug/L | 95 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | | N.D. | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | | N.D. d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | | N.D. | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | | N.D. | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | | N.D. | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | | N.D. | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | | N.D. | |
| 42) Styrene | 0.00 | 104 | 0 | | N.D. | |
| 43) Bromoform | 0.00 | 173 | 0 | | N.D. | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 46) Bromobenzene | 0.00 | 156 | 0 | | N.D. | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | | N.D. | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | | N.D. | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | | N.D. | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | | N.D. | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | | N.D. | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | | N.D. | |
| 65) Naphthalene | 0.00 | 128 | 0 | | N.D. | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | | N.D. | |
| 68) Acetone | 0.00 | 43 | 0 | | N.D. | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | | N.D. | |
| 70) 2-Butanone | 0.00 | 43 | 0 | | N.D. | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | | N.D. | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | | N.D. | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | | N.D. | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

7502_9.D 71NV16_6.M

Wed Jan 17 11:14:03 2007

GC#2

Page 2

Quantitation Report

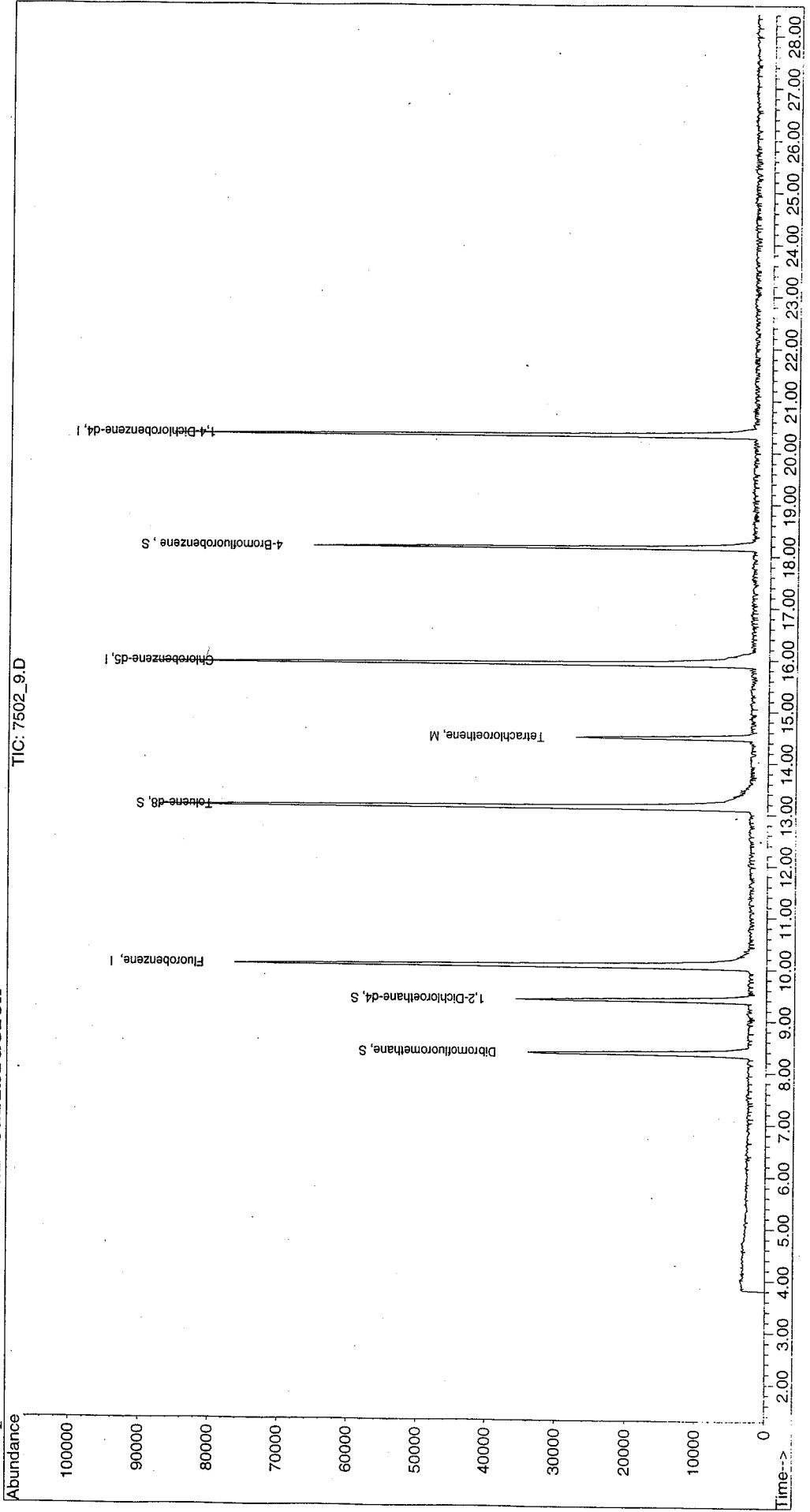
Data File : C:\HPCHEM\1\DATA\010207\7502_9.D
Acq On : 2 Jan 2007 11:08 pm
Sample : 7502-09 0.050 ml
Misc :

Vial: 11
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00

MS Integration Params: ODD.P
Quant Time: Jan 3 8:43 2007

Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Acq On : 17 Jan 2007 5:01 pm

Operator: RR/AS

Sample : 7502-10 0.20 ml

Inst : GC/MS Ins

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 22 11:03 2007

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:54:07 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.73 | 96 | 275913 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 20.11 | 117 | 240158 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.35 | 152 | 136416 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|--------|-------|------------|---------|
| 16) Dibromofluoromethane | 13.29 | 113 | 84635 | 28.26 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | Recovery = | 113.04% |
| 19) 1,2-Dichloroethane-d4 | 14.24 | 65 | 83122 | 27.40 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | Recovery = | 109.60% |
| 28) Toluene-d8 | 17.41 | 98 | 311004 | 24.96 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | Recovery = | 99.84% |
| 45) 4-Bromofluorobenzene | 22.27 | 95 | 87963 | 21.45 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | Recovery = | 85.80% |

Target Compounds

Qvalue

| | | | | |
|-------------------------------|------|-----|---|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. |
| 21) Benzene | 0.00 | 78 | 0 | N.D. |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. |
| 29) Toluene | 0.00 | 92 | 0 | N.D. |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. |

(#)= qualifier out of range (m) = manual integration

7502_10.D 01_12_07.M Tue Jan 23 08:53:27 2007

Acq On : 17 Jan 2007 5:01 pm

Operator: RR/AS

Sample : 7502-10 0.20 ml

Inst : GC/MS Ins

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 22 11:03 2007

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:54:07 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 18.18 | 166 | 300564 | 55.76 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | d | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

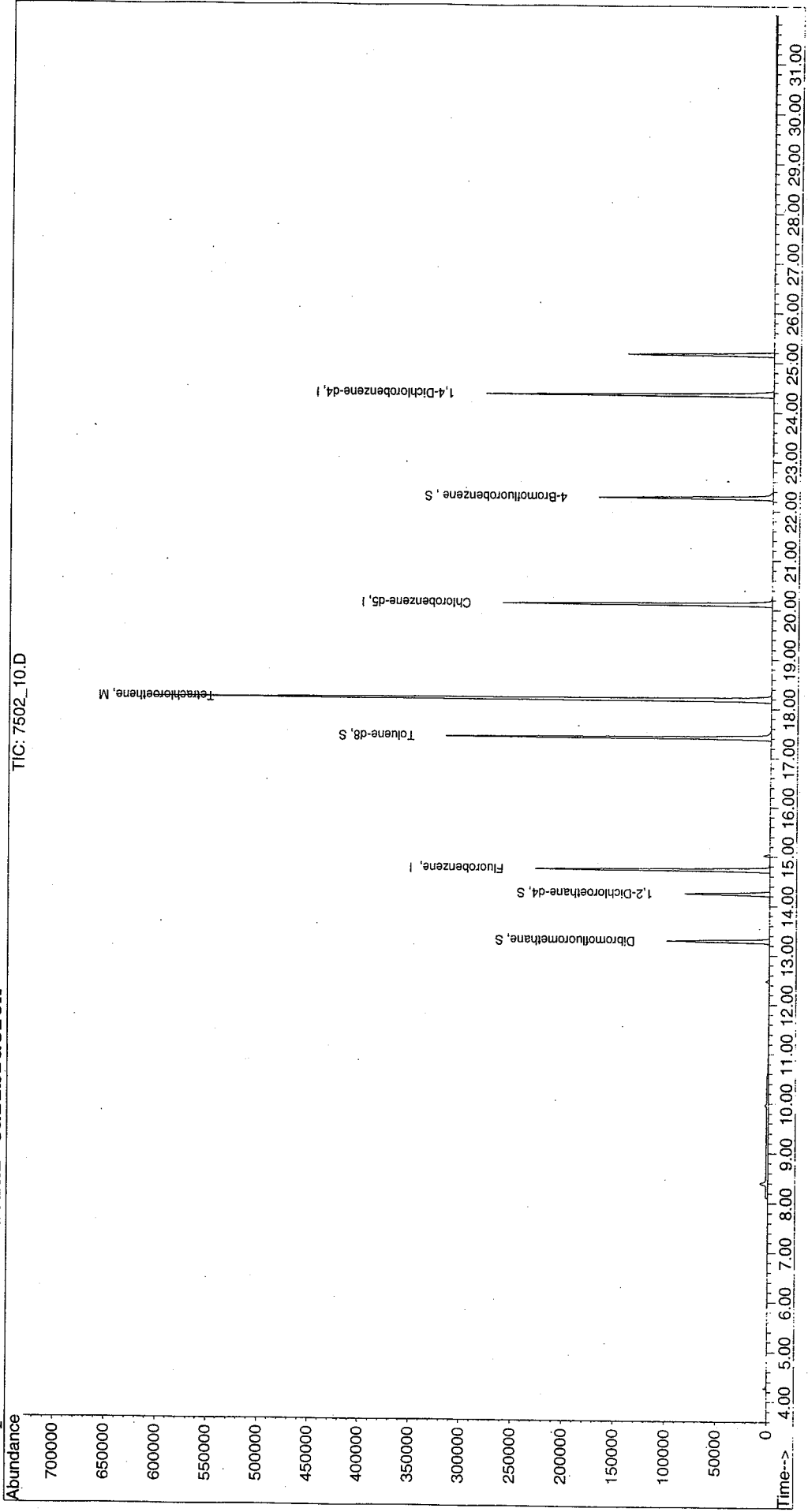
(#) = qualifier out of range (m) = manual integration

7502_10.D 01_12_07.M Tue Jan 23 08:53:29 2007

Quantitation Report

Data File : C:\HPCHEM\1\DATA\0101707\7502_10.D
Acq On : 17 Jan 2007 5:01 pm
Sample : 7502-10 0.20 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 22 11:03 2007
Vial: 1
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: 01_12_07.RE5

Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:54:07 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\010207\7502_11.D
 Acq On : 3 Jan 2007 12:18 am
 Sample : 7502-11 0.050 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 3 8:41 2007

Vial: 13
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Thu Nov 16 17:01:55 2006
 Response via : Initial Calibration
 DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.07 | 96 | 186571 | 25.00 | ug/L | -0.03 |
| 36) Chlorobenzene-d5 | 15.92 | 117 | 122872 | 25.00 | ug/L | -0.03 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 50083 | 25.00 | ug/L | -0.03 |

| System Monitoring Compounds | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------|--------|------|----------|-------|---------|----------|
| 16) Dibromofluoromethane | 8.37 | 113 | 50870 | 28.20 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 112.80% | |
| 19) 1,2-Dichloroethane-d4 | 9.41 | 65 | 40603 | 26.01 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 104.04% | |
| 28) Toluene-d8 | 13.14 | 98 | 154547 | 24.72 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 98.88% | |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 49381 | 21.69 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 86.76% | |

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\010207\7502_11.D
Acq On : 3 Jan 2007 12:18 am
Sample : 7502-11 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 3 8:41 2007

Vial: 13
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00

Quant Results File: 71NV16_6.RES

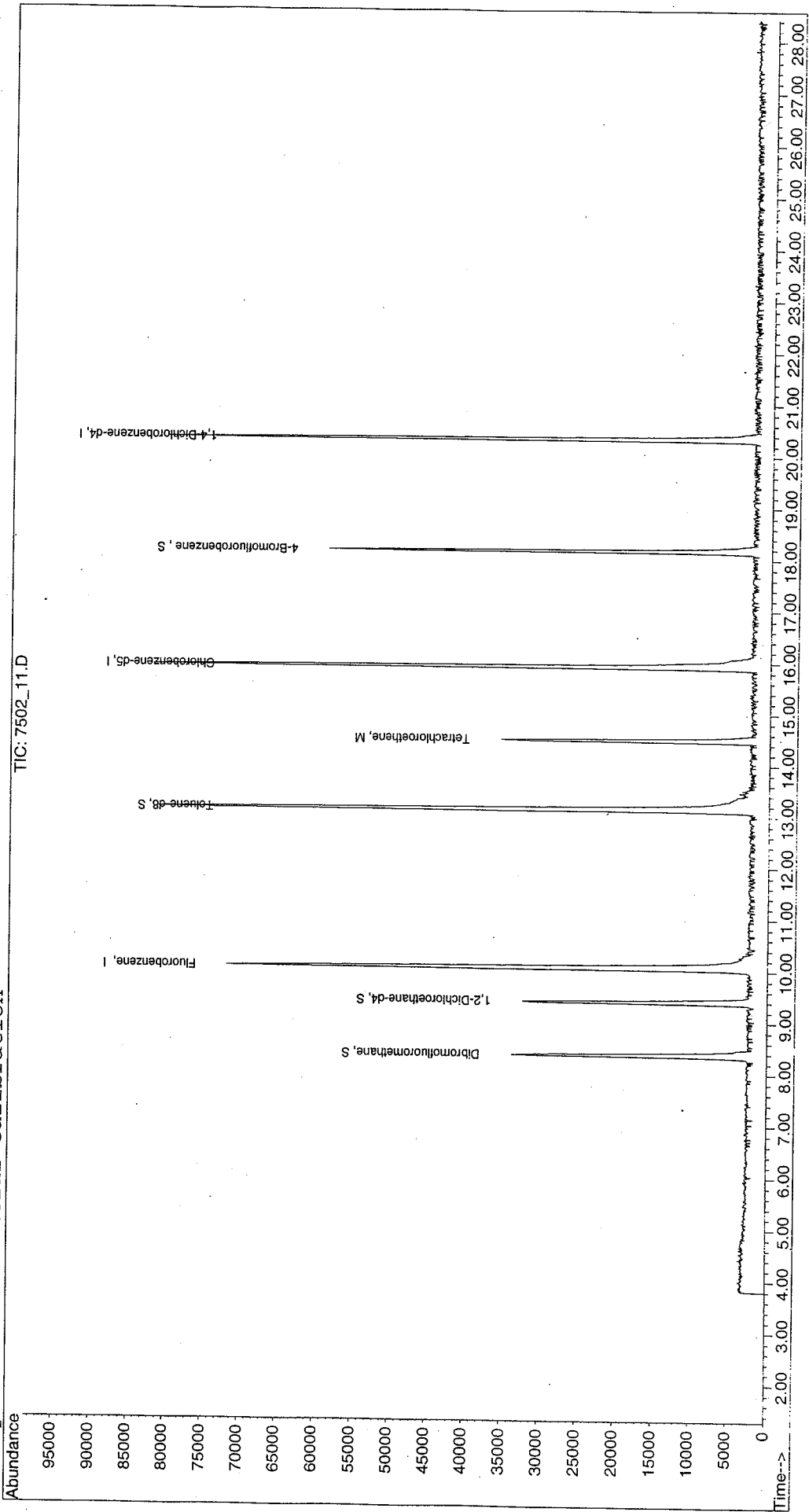
Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Thu Nov 16 17:01:55 2006
Response via : Initial Calibration
DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 25998 | 17.26 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\010207\7502_11.D
Acq On : 3 Jan 2007 12:18 am
Sample : 7502-11 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 3 8:41 2007
Quant Results File: 71NV16_6.REIS
Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Acq On : 3 Jan 2007 12:52 am

Operator: TS

Sample : 7502-12 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 3 8:42 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|--------|------|----------|-------|-------|--------------------|
| 1) Fluorobenzene | 10.06 | 96 | 188298 | 25.00 | ug/L | -0.04 |
| 36) Chlorobenzene-d5 | 15.92 | 117 | 123640 | 25.00 | ug/L | -0.03 |
| 58) 1,4-Dichlorobenzene-d4 | 20.33 | 152 | 51664 | 25.00 | ug/L | -0.03 |
| System Monitoring Compounds | | | | | | |
| 16) Dibromofluoromethane | 8.37 | 113 | 50100 | 27.52 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 110.08% |
| 19) 1,2-Dichloroethane-d4 | 9.41 | 65 | 40213 | 25.52 | ug/L | -0.02 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 102.08% |
| 28) Toluene-d8 | 13.14 | 98 | 152117 | 24.10 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 96.40% |
| 45) 4-Bromofluorobenzene | 18.16 | 95 | 51683 | 22.56 | ug/L | -0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 90.24% |
| Target Compounds | | | | | | |
| | | | | | | Qvalue |
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. | d |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

Acq On : 3 Jan 2007 12:52 am

Operator: TS

Sample : 7502-12 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 3 8:42 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Thu Nov 16 17:01:55 2006

Response via : Initial Calibration

DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|------|------|--------|
| 32) Tetrachloroethene | 14.49 | 166 | 9236 | 6.08 | ug/L | 96 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7502_12.D 71NV16_6.M

Wed Jan 17 11:14:00 2007

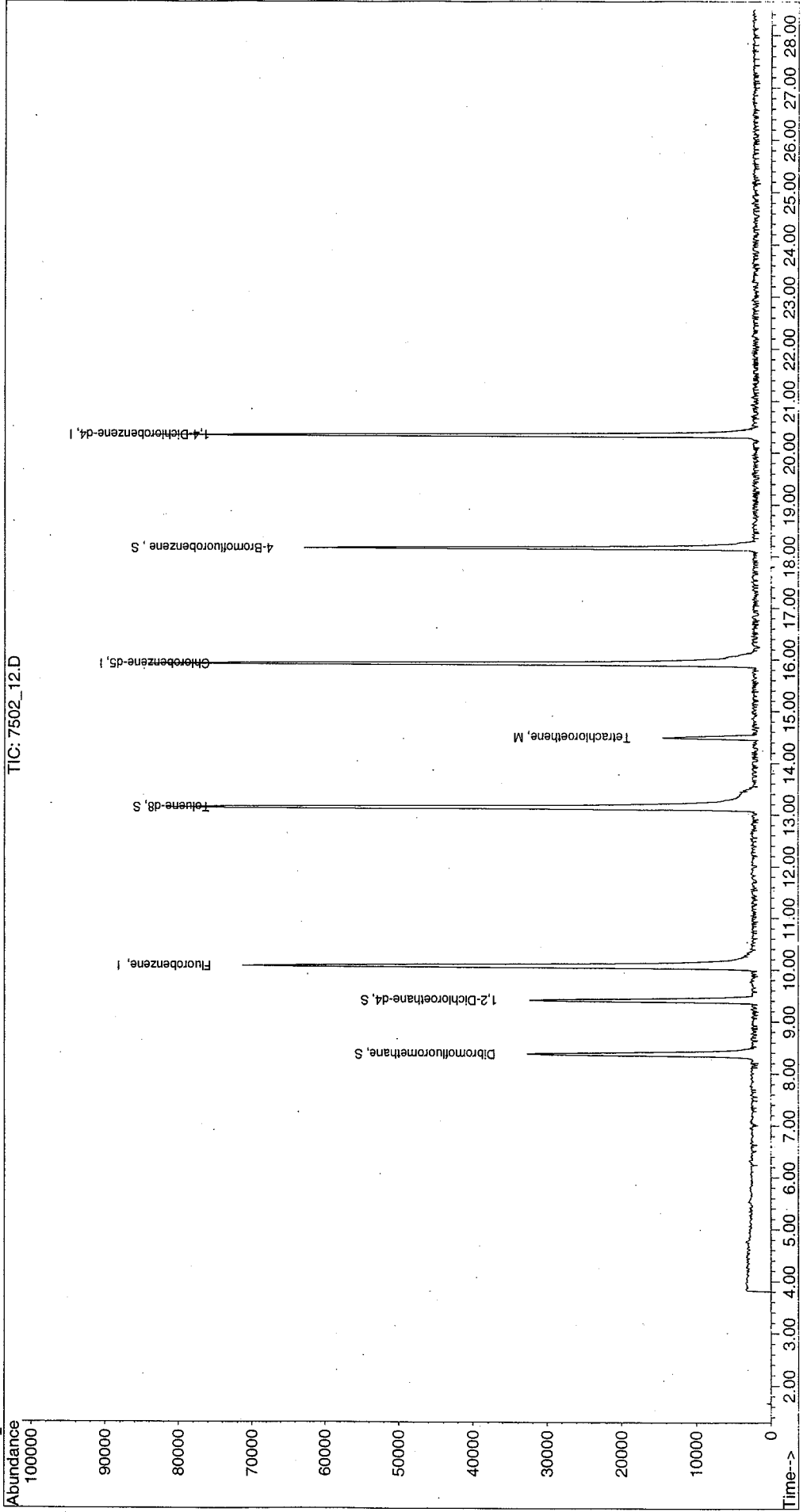
GC#2

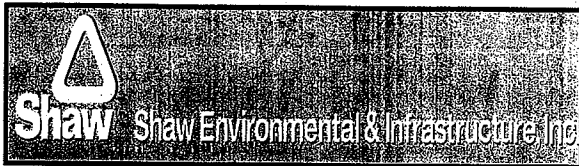
Page 2

Quantitation Report

Data File : C:\HPCHEM\1\DATA\010207\7502_12.D Vial: 14
Acq On : 3 Jan 2007 12:52 am Operator: TS
Sample : 7502-12 0.050 ml Inst : GC/MS #2
Misc : Multiplr: 1.00
MS Integration Params: ODD.P
Quant Time: Jan 3 8:42 2007 Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration





17 Princess Rd
Lawrenceville, New Jersey 08648
Tel: 609/895-5370
Fax: 609/895-1858

Volatile Organic Compound Data Summary Package

Prepared for
Gorham Textron

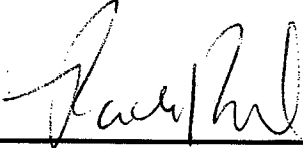
Lab ID
7504

Project Number 101960 02000000

Samples Received
3-Jan-07

Reported
18-Jan-07

NJDEP Certified Lab 11001


Randi K Rothmel, PhD
Laboratory Director


Date

1.0 Chain of Custody

CHAIN OF CUSTODY

Ref. Document #

Page 1 of 1

Project Number/Coast code: 101960 / 02

Project Name / Location: Gorham Textron

Send bill to:

Project Contact: Charles Schaefer, Chuck Condee
(Name & phone #)

Send Report To: Chuck Condee

Phone/Fax Number:

Address:

City/State:

Shipment date:

Lab Destination:

Lab Contact Name / ph. #:

Sampler's Name(s): Sheryl Stieger

Collection Information

| Lab No. | Sample ID Number | Sample Description | Collection Information | | Matrix | # of containers | Container type | Preservative | | | | MeOH | Std | |
|---------|------------------|--------------------|------------------------|------|--------|-----------------|----------------|--------------|------|------------------|--------------------------------|------|-----|------|
| | | | Date | Time | | | | HCL | NaOH | HNO ₃ | H ₂ SO ₄ | | | None |
| 1 | Killed Control A | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | X | Std |
| 2 | " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | X | Std |
| 3 | Killed Control B | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | X | Std |
| 4 | " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | X | Std |
| 5 | Killed Control C | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | X | Std |
| 6 | " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | X | Std |
| 7 | Live Control A | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | X | Std |
| 8 | " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | X | Std |
| 9 | Live Control B | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | X | Std |
| 10 | " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | X | Std |
| 11 | Live Control C | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | X | Std |
| 12 | " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | X | Std |
| 13 | Bloaug 1 LA | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | X | Std |
| 14 | " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | X | Std |
| 15 | Bloaug 1 LB | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | X | Std |
| 16 | " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | X | Std |
| 17 | Bloaug 1 LC | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | X | Std |
| 18 | " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | X | Std |
| 19 | Bloaug 2 EA | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | X | Std |
| 20 | " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | X | Std |
| 21 | Bloaug 2 EB | | 01/03/07 | 9:35 | g | Aq | 1 | 10mL | X | | | | X | Std |
| 22 | " | | 01/03/07 | 9:35 | g | Aq | 1 | 5mL | | | | | X | Std |

Special Instructions:

Killed controls contain 1.5% formaldehyde.

Known Waste Stream Circle:
RCRA PCB/dioxin PAH/oli
QC/Data Package Level Required: I II III IV

Flammable

Corrosive

GIS EDD

Preliminary data

G = Grab

C = Composite

Reactive

NJ EDD

QC Package Codes

Analyses Requested

VOC

MTH

Turn Around Time Requested

Level II = data summary + basic QC
Level III = New Jersey QC reduced deliverable
Level IV = Full deliverable CLP package

Cooler temperature upon arrival at Lab:

Received By:

Date: 1-3-07

Time: 1008

Received By:

Date: 01/03/07

Time: 10:30

Received By:

Date:

Time:

Received By:

Date:

Time:



COC Continuation Page

COC Ref. Document #

7504

Project Number/ cost code: 101960

02

Shipment Date:

Project Name / Location: Gorham Textron

| Lab No. | Sample ID Number | Sample Description | Collection Information | | # of containers | Container Type | HCL | NaOH | HNO ₃ | H ₂ SO ₄ | Preservative | | | VOC | MEE | Anions | VFA | Turn Around Time Requested |
|---------|------------------|--------------------|------------------------|------|-----------------|----------------|-----|------|------------------|--------------------------------|--------------|----------|------|-----|-----|--------|-----|----------------------------|
| | | | Date | Time | | | | | | | G/C | Filtered | None | | | | | |
| 12 | Bloaug 2 E C | | 01/03/07 | 9:35 | g | 10mL | X | | | | X | | | X | | | | |
| 13 | Killed Control D | | 01/03/07 | 9:35 | g | 5mL | | | | | X | | | X | | | | |
| 14 | Live Control D | | 01/03/07 | 9:35 | g | 15mL | | | | | X | | | X | | | | |
| 15 | Bloaug 1 L D | | 01/03/07 | 9:35 | g | 15mL | | | | | X | | | X | | | | |
| 16 | Bloaug 2 E D | | 01/03/07 | 9:35 | g | 15mL | | | | | X | | | X | | | | |
| | | | | | | | | | | | | | | | | | | |

Analyses Requested

Lab ID 7504-16

Client:

Date Received:

Shaw E&I Analytical and Treatability Laboratories Internal Chain of Custody

Pg _____ of _____

| Sample ID | Parameter | Bottle Type | Preservative | Date/Time Removed | Relinquishing Custodian Initials | Receiving Analyst Initials | Date/Time Returned | Receiving Custodian Initials | Relinquishing Analyst Initials |
|-----------|-----------|-------------|--------------|-------------------|----------------------------------|----------------------------|--------------------|------------------------------|--------------------------------|
| 7504-1 | YDC's | 1-10-1 | HCl | 01-06-07 9:30 | AS | AS | 01-06-07 174 | | AS |
| 2 | | | | | | | | | |
| 3 | | | | | | | | | |
| 4 | | | | | | | | | |
| 5 | | | | | | | | | |
| 6 | | | | | | | | | |
| 7 | | | | | | | | | |
| 8 | | | | | | | | | |
| 9 | | | | | | | | | |
| 10 | | | | | | | | | |
| 11 | | | | | | | | | |
| 7504-12 | YDC's | 1-10-1 | HCl | 01-06-07 9:30 | AS | AS | 01-06-07 174 | | AS |
| 7504-1 | MSEP | 1-10-1 | None | 11/5/07 8:00 | PM | PM | 11/5/07 15:00 | | PM |
| 2 | | | | | | | | | |
| 3 | | | | | | | | | |
| 4 | | | | | | | | | |
| 5 | | | | | | | | | |
| 6 | | | | | | | | | |
| 7 | | | | | | | | | |
| 8 | | | | | | | | | |
| 9 | | | | | | | | | |
| 10 | | | | | | | | | |
| 11 | | | | | | | | | |
| 7504-12 | MSEP | 1-10-1 | None | | | | | | |
| 7504-13 | VFA's | 10-1 | None | 1/4/07 12:00 | PM | PM | 1/4/07 12:00 | | PM |
| 14 | | 10-1 | | | | | | | |
| 15 | | 10-1 | | | | | | | |
| 7504-16 | VFA's | 10-1 | None | | | | | | |

2.0. Sample Summaries

Sample summaries are enclosed

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

Lab Name: ATL Analyst: Tony Soto 7504-01
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_01 0.050 ml
 Sample wt/vol: ~~0.10050~~ (g/ml) ML Lab File ID: 7504-01.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | 4200 | | U |
| 74-87-3 | Chloromethane | 4200 | | U |
| 75-01-4 | Vinyl chloride | 4200 | | U |
| 74-83-9 | Bromomethane | 4200 | | U |
| 75-00-3 | Chloroethane | 4200 | | U |
| 75-69-4 | Trichlorofluoromethane | 4200 | | U |
| 75-35-4 | 1,1-Dichloroethene | 4200 | | U |
| 75-09-2 | Methylene chloride | 4200 | | U |
| 156-60-5 | trans-1,2-Dichloroethene | 4200 | | U |
| 75-34-3 | 1,1-Dichloroethane | 4200 | | U |
| 594-20-7 | 2,2-Dichloropropane | 4200 | | U |
| 156-59-2 | cis-1,2-Dichloroethene | 4200 | | U |
| 74-97-5 | Bromochloromethane | 4200 | | U |
| 67-66-3 | Chloroform | 4200 | | U |
| 71-55-6 | 1,1,1-Trichloroethane | 4200 | | U |
| 56-23-5 | Carbon tetrachloride | 4200 | | U |
| 563-58-6 | 1,1-Dichloropropene | 4200 | | U |
| 71-43-2 | Benzene | 4200 | | U |
| 107-06-2 | 1,2-Dichloroethane | 4200 | | U |
| 79-01-6 | Trichloroethene | 4200 | | U |
| 78-87-5 | 1,2-Dichloropropane | 4200 | | U |
| 74-95-3 | Dibromomethane | 4200 | | U |
| 75-27-4 | Bromodichloromethane | 4200 | | U |
| 10061-02-5 | cis-1,3-Dichloropropene | 4200 | | U |
| 108-88-3 | Toluene | 4200 | | U |
| 10061-01-6 | trans-1,3-Dichloropropene | 4200 | | U |
| 79-00-5 | 1,1,2-Trichloroethane | 4200 | | U |
| 127-18-4 | Tetrachloroethene | 4200 | | U |
| 142-28-9 | 1,3-Dichloropropane | 20000 | | D |
| 124-48-1 | Dibromochloromethane | 4200 | | U |
| 106-93-4 | 1,2-Dibromoethane | 4200 | | U |
| 108-90-7 | Chlorobenzene | 4200 | | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 4200 | | U |
| 100-41-4 | Ethylbenzene | 4200 | | U |
| 1330-20-7 | Xylene (para & meta) | 4200 | | U |
| 95-47-6 | Xylene (Ortho) | 4200 | | U |
| 100-42-5 | Styrene | 4200 | | U |
| 75-25-2 | Bromoform | 4200 | | U |
| 98-82-8 | Isopropylbenzene | 4200 | | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-01

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_01 0.050 ml
 Sample wt/vol: 0.050 (g/ml) ML Lab File ID: 7504-01.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-02

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_02 0.050 ml
 Sample wt/vol: 0.1^{AD} 0.050 (g/ml) ML Lab File ID: 7504-02.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 21000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-02

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_02 0.050 ml
 Sample wt/vol: 0.1^A 0.2^{SO} (g/ml) ML Lab File ID: 7504-02.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | CONCENTRATION UNITS: | |
|-----------|-----------------------------|----------------------|------|
| | | (ug/L or ug/Kg) | UG/L |
| 108-86-1 | Bromobenzene | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | 4200 | U |
| 103-65-1 | n-Propylbenzene | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 4200 | U |
| 98-06-6 | tert-Butylbenzene | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 4200 | U |
| 135-98-8 | sec-Butylbenzene | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | 4200 | U |
| 104-51-8 | n-Butylbenzene | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | 4200 | U |
| 91-20-3 | Naphthalene | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 4200 | U |
| 1634-04-4 | MTBE | 4200 | U |
| 67-64-1 | Acetone | 8400 | U |
| 75-15-0 | Carbon disulfide | 4200 | U |
| 78-93-3 | 2-Butanone | 8400 | U |
| 109-99-9 | Tetrahydrofuran | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 8400 | U |
| 591-78-6 | 2-Hexanone | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-03

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_03 0.050 ml
 Sample wt/vol: 0.170.050 (g/ml) ML Lab File ID: 7504-03.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 19000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-03

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_03 0.050 ml
 Sample wt/vol: 0.1^{AS} 0.050 (g/ml) ML Lab File ID: 7504-03.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-04

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_04 0.050 ml
 Sample wt/vol: 0.1305 (g/ml) ML Lab File ID: 7504-04.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 8800 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-04

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_04 0.050 ml
 Sample wt/vol: 0.186.05 (g/ml) ML Lab File ID: 7504-04.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7404-05

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_05 0.050 ml
 Sample wt/vol: 0.113^{0.050} (g/ml) ML Lab File ID: 7504-05.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 17000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7404-05

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_05 0.050 ml
 Sample wt/vol: 0.1^A 0.05⁰ (g/ml) ML Lab File ID: 7504-05.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-06

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_06 0.050 ml
 Sample wt/vol: 0.1^g 0.05^g (g/ml) ML Lab File ID: 7504-06.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 18000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-06

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_06 0.050 ml
 Sample wt/vol: 0.1170.050 (g/ml) ML Lab File ID: 7504-06.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-07

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_07 0.050 ml
 Sample wt/vol: 0.1¹ 0.05⁰ (g/ml) ML Lab File ID: 7504-07.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 19000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-07

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_07 0.050 ml
 Sample wt/vol: 0.125 (g/ml) ML Lab File ID: 7504-07.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-08

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_08 0.050 ml
 Sample wt/vol: 0.105 (g/ml) ML Lab File ID: 7504-08.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 26000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-08

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_08 0.050 ml
 Sample wt/vol: 0.117 0.050 (g/ml) ML Lab File ID: 7504-08.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-09

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_09 0.050 ml
 Sample wt/vol: 0.17050 (g/ml) ML Lab File ID: 7504-09.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 14000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-09

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_09 0.050 ml
 Sample wt/vol: 0.120 (g/ml) ML Lab File ID: 7504-09.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-10

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_10 0.050 ml
 Sample wt/vol: 0.1 (g/ml) ML Lab File ID: 7504-10.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|----|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 4200 | U |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U* |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-10

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_10 0.050 ml
 Sample wt/vol: 0.1 (g/ml) ML Lab File ID: 7504-10.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7404-11

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_11 0.050 ml
 Sample wt/vol: 0.14 → 0.05 (g/ml) ML Lab File ID: 7504-11.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 19000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7404-11

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_11 0.050 ml
 Sample wt/vol: 0.1A, 0.050 (g/ml) ML Lab File ID: 7504-11.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-12

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_12 0.050 ml
 Sample wt/vol: 0.125 (g/ml) ML Lab File ID: 7504-12.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 4200 | U |
| 74-87-3 | Chloromethane | | 4200 | U |
| 75-01-4 | Vinyl chloride | | 4200 | U |
| 74-83-9 | Bromomethane | | 4200 | U |
| 75-00-3 | Chloroethane | | 4200 | U |
| 75-69-4 | Trichlorofluoromethane | | 4200 | U |
| 75-35-4 | 1,1-Dichloroethene | | 4200 | U |
| 75-09-2 | Methylene chloride | | 4200 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 4200 | U |
| 75-34-3 | 1,1-Dichloroethane | | 4200 | U |
| 594-20-7 | 2,2-Dichloropropane | | 4200 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 4200 | U |
| 74-97-5 | Bromochloromethane | | 4200 | U |
| 67-66-3 | Chloroform | | 4200 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 4200 | U |
| 56-23-5 | Carbon tetrachloride | | 4200 | U |
| 563-58-6 | 1,1-Dichloropropene | | 4200 | U |
| 71-43-2 | Benzene | | 4200 | U |
| 107-06-2 | 1,2-Dichloroethane | | 4200 | U |
| 79-01-6 | Trichloroethene | | 4200 | U |
| 78-87-5 | 1,2-Dichloropropane | | 4200 | U |
| 74-95-3 | Dibromomethane | | 4200 | U |
| 75-27-4 | Bromodichloromethane | | 4200 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 4200 | U |
| 108-88-3 | Toluene | | 4200 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 4200 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 4200 | U |
| 127-18-4 | Tetrachloroethene | | 16000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 4200 | U |
| 124-48-1 | Dibromochloromethane | | 4200 | U |
| 106-93-4 | 1,2-Dibromoethane | | 4200 | U |
| 108-90-7 | Chlorobenzene | | 4200 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 4200 | U |
| 100-41-4 | Ethylbenzene | | 4200 | U |
| 1330-20-7 | Xylene (para & meta) | | 4200 | U |
| 95-47-6 | Xylene (Ortho) | | 4200 | U |
| 100-42-5 | Styrene | | 4200 | U |
| 75-25-2 | Bromoform | | 4200 | U |
| 98-82-8 | Isopropylbenzene | | 4200 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE

7504-12

Lab Name: ATL Analyst: Tony Soto
 NJDEP: 11001 Calib date 11-16-06 GC/MS: Two Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7504_12 0.050 ml
 Sample wt/vol: 0.050 (g/ml) ML Lab File ID: 7504-12.D
 Level: (low/med) LOW Date Received: 01/04/06
 % Moisture: not dec. _____ Date Analyzed: 01/06/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 840.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 108-86-1 | Bromobenzene | | 4200 | U |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 4200 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 4200 | U |
| 103-65-1 | n-Propylbenzene | | 4200 | U |
| 95-49-8 | 2-Chlorotoluene | | 4200 | U |
| 106-43-4 | 4-Chlorotoluene | | 4200 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 4200 | U |
| 98-06-6 | tert-Butylbenzene | | 4200 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 4200 | U |
| 135-98-8 | sec-Butylbenzene | | 4200 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 4200 | U |
| 99-87-6 | 4-Isopropyltoluene | | 4200 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 4200 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 4200 | U |
| 104-51-8 | n-Butylbenzene | | 4200 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 4200 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 4200 | U |
| 87-68-3 | Hexachlorobutadiene | | 4200 | U |
| 91-20-3 | Naphthalene | | 4200 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 4200 | U |
| 1634-04-4 | MTBE | | 4200 | U |
| 67-64-1 | Acetone | | 8400 | U |
| 75-15-0 | Carbon disulfide | | 4200 | U |
| 78-93-3 | 2-Butanone | | 8400 | U |
| 109-99-9 | Tetrahydrofuran | | 8400 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 8400 | U |
| 591-78-6 | 2-Hexanone | | 8400 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 8400 | U |

3.0 Raw Data and Chromatograms

Raw data and Chromatograms are attached.

Data File : C:\HPCHEM\1\DATA\010507\7504-01.D
 Acq On : 6 Jan 2007 6:47 am
 Sample : 7504_01 0.050 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 17 15:33 2007

Vial: 31
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Wed Jan 10 11:01:47 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.13 | 96 | 153260 | 25.00 | ug/L | 0.03 |
| 36) Chlorobenzene-d5 | 15.97 | 117 | 107178 | 25.00 | ug/L | 0.02 |
| 58) 1,4-Dichlorobenzene-d4 | 20.38 | 152 | 43070 | 25.00 | ug/L | 0.02 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|--------|----------|------|---------|
| 16) Dibromofluoromethane | 8.44 | 113 | 43613 | 29.44 | ug/L | 0.04 |
| Spiked Amount | 25.000 | | | Recovery | = | 117.76% |
| 19) 1,2-Dichloroethane-d4 | 9.46 | 65 | 35130 | 27.39 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | | Recovery | = | 109.56% |
| 28) Toluene-d8 | 13.19 | 98 | 134744 | 26.23 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | | Recovery | = | 104.92% |
| 45) 4-Bromofluorobenzene | 18.22 | 95 | 43419 | 21.86 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | | Recovery | = | 87.44% |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\010507\7504-01.D
 Acq On : 6 Jan 2007 6:47 am
 Sample : 7504_01 0.050 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 17 15:33 2007

Vial: 31
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Wed Jan 10 11:01:47 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN1

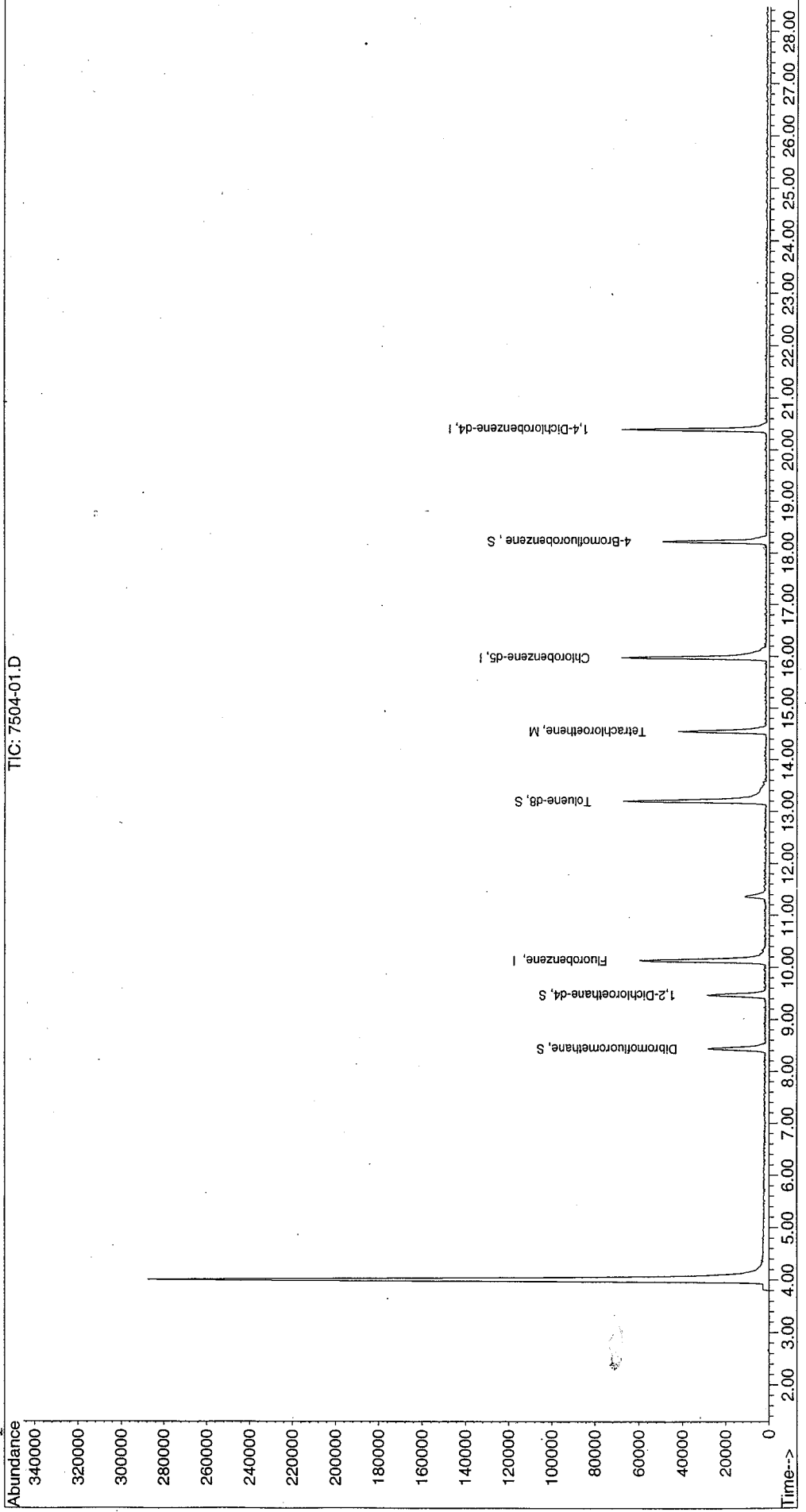
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.54 | 166 | 29405 | 23.77 | ug/L | 96 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | d | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\010507\7504-01.D
Acq On : 6 Jan 2007 6:47 am Vial: 31
Sample : 7504_01 0.050 ml Operator: TS
Misc : Inst : GC/MS #2
MS Integration Params: ODD.P Multiplr: 1.00
Quant Time: Jan 17 15:33 2007 Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\010507\7504-02.D
 Acq On : 6 Jan 2007 7:22 am
 Sample : 7504_02 0.050 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 17 15:33 2007

Vial: 32
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Wed Jan 10 11:01:47 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.13 | 96 | 150114 | 25.00 | ug/L | 0.03 |
| 36) Chlorobenzene-d5 | 15.97 | 117 | 106453 | 25.00 | ug/L | 0.02 |
| 58) 1,4-Dichlorobenzene-d4 | 20.38 | 152 | 44079 | 25.00 | ug/L | 0.02 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|--------|----------|------|---------|
| 16) Dibromofluoromethane | 8.44 | 113 | 42987 | 29.62 | ug/L | 0.04 |
| Spiked Amount | 25.000 | | | Recovery | = | 118.48% |
| 19) 1,2-Dichloroethane-d4 | 9.46 | 65 | 34557 | 27.51 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | | Recovery | = | 110.04% |
| 28) Toluene-d8 | 13.19 | 98 | 123872 | 24.62 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | | Recovery | = | 98.48% |
| 45) 4-Bromofluorobenzene | 18.21 | 95 | 45314 | 22.97 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | | Recovery | = | 91.88% |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\010507\7504-02.D
 Acq On : 6 Jan 2007 7:22 am
 Sample : 7504_02 0.050 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 17 15:33 2007

Vial: 32
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

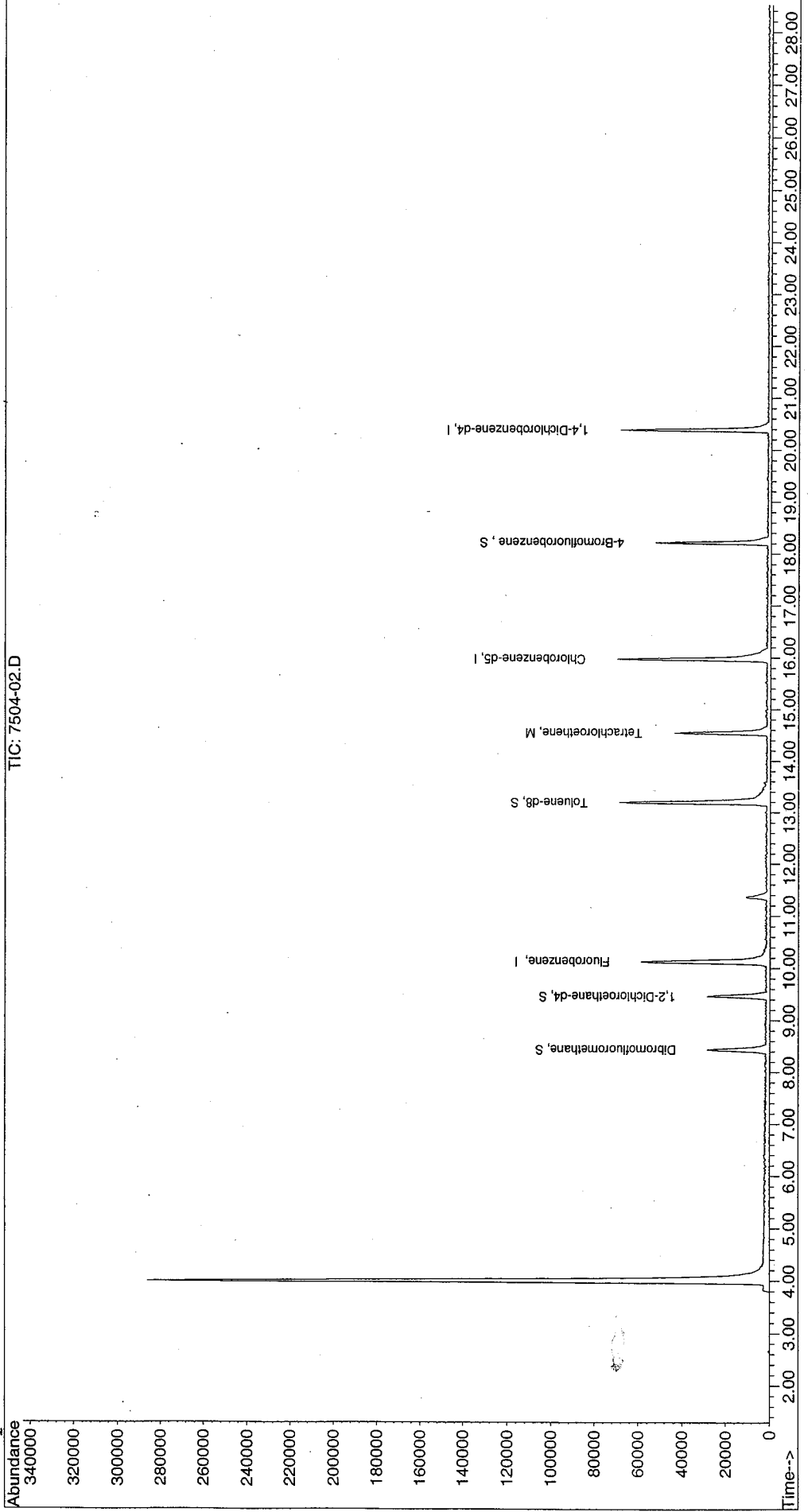
Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Wed Jan 10 11:01:47 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.54 | 166 | 30427 | 25.11 | ug/L | 91 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | d | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

Quantitation Report

Data File : C:\HPCHEM\1\DATA\010507\7504-02.D
Acq On : 6 Jan 2007 7:22 am
Sample : 7504_02 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 17 15:33 2007
Vial: 32
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\010507\7504-03.D
Acq On : 6 Jan 2007 7:57 am
Sample : 7504_03 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 17 15:34 2007

Vial: 33
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration
DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.12 | 96 | 149385 | 25.00 | ug/L | 0.02 |
| 36) Chlorobenzene-d5 | 15.97 | 117 | 99630 | 25.00 | ug/L | 0.02 |
| 58) 1,4-Dichlorobenzene-d4 | 20.38 | 152 | 42339 | 25.00 | ug/L | 0.02 |

| System Monitoring Compounds | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------|--------|------|----------|-------|---------|----------|
| 16) Dibromofluoromethane | 8.44 | 113 | 43298 | 29.98 | ug/L | 0.04 |
| Spiked Amount | 25.000 | | Recovery | = | 119.92% | |
| 19) 1,2-Dichloroethane-d4 | 9.46 | 65 | 34697 | 27.76 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 111.04% | |
| 28) Toluene-d8 | 13.19 | 98 | 116794 | 23.33 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 93.32% | |
| 45) 4-Bromofluorobenzene | 18.21 | 95 | 41723 | 22.60 | ug/L | 0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 90.40% | |

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration
7504-03.D 71NV16_6.M Thu Jan 18 09:24:08 2007 GC#2

Data File : C:\HPCHEM\1\DATA\010507\7504-03.D
Acq On : 6 Jan 2007 7:57 am
Sample : 7504_03 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 17 15:34 2007

Vial: 33
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration
DataAcq Meth : RUN1

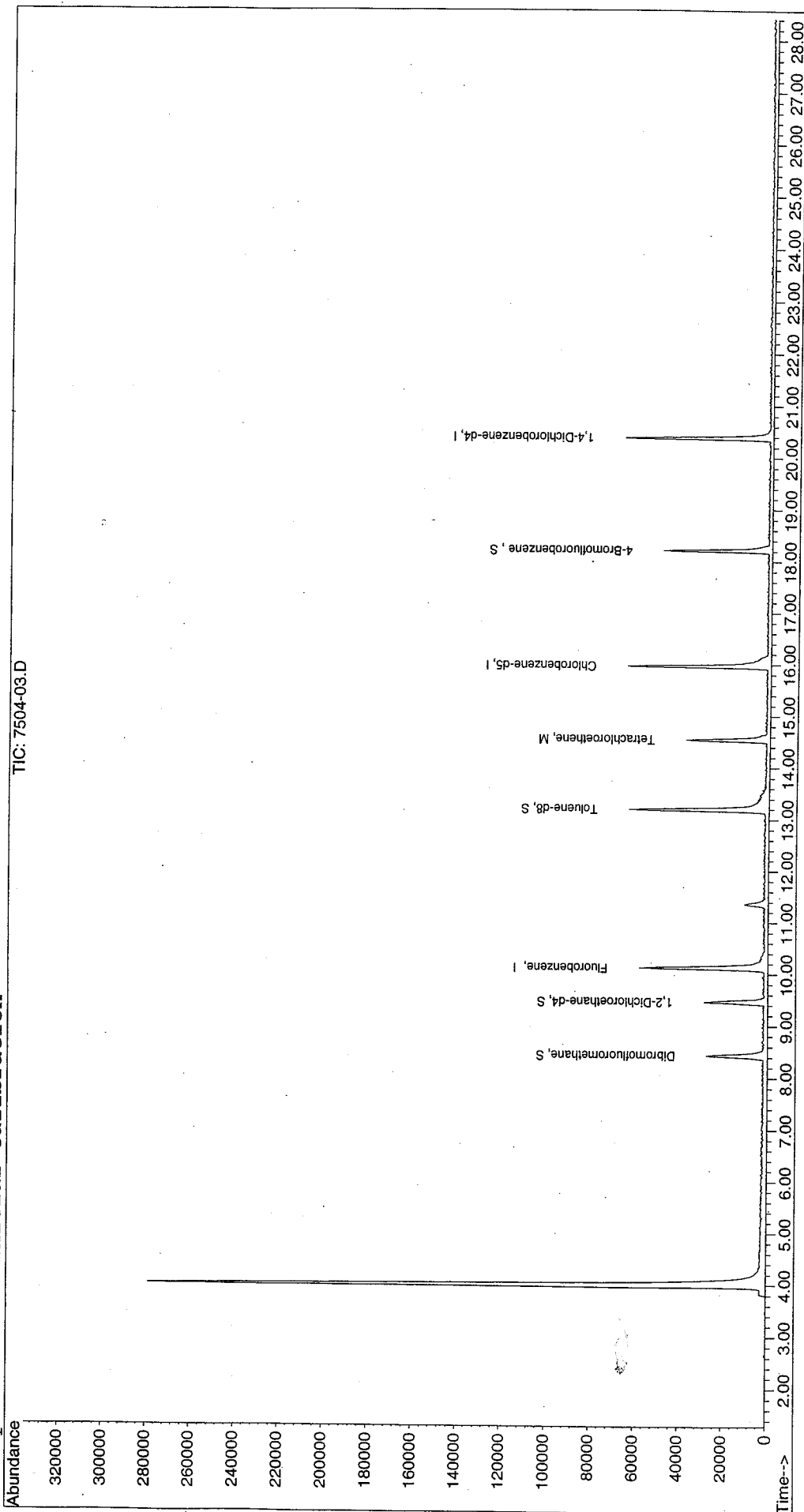
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.54 | 166 | 27960 | 23.19 | ug/L | 94 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | d | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration
7504-03.D 71NV16_6.M Thu Jan 18 09:24:09 2007 GC#2

Quantitation Report

Data File : C:\HPCHEM\1\DATA\010507\7504-03.D
Acq On : 6 Jan 2007 7:57 am Vial: 33
Sample : 7504_03 0.050 ml Operator: TS
Misc : GC/MS #2
MS Integration Params: ODD.P Inst : GC/MS #2
Quant Time: Jan 17 15:34 2007 Multiplr: 1.00
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\010507\7504-04.D
Acq On : 6 Jan 2007 8:33 am
Sample : 7504_04 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 17 15:34 2007

Vial: 34
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration
DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.12 | 96 | 151395 | 25.00 | ug/L | 0.02 |
| 36) Chlorobenzene-d5 | 15.97 | 117 | 99759 | 25.00 | ug/L | 0.02 |
| 58) 1,4-Dichlorobenzene-d4 | 20.38 | 152 | 42056 | 25.00 | ug/L | 0.02 |

| System Monitoring Compounds | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------|--------|------|----------|-------|---------|----------|
| 16) Dibromofluoromethane | 8.43 | 113 | 42694 | 29.17 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 116.68% | |
| 19) 1,2-Dichloroethane-d4 | 9.46 | 65 | 35026 | 27.65 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 110.60% | |
| 28) Toluene-d8 | 13.19 | 98 | 122479 | 24.14 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 96.56% | |
| 45) 4-Bromofluorobenzene | 18.20 | 95 | 40811 | 22.08 | ug/L | 0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 88.32% | |

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\010507\7504-04.D
 Acq On : 6 Jan 2007 8:33 am
 Sample : 7504_04 0.050 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 17 15:34 2007

Vial: 34
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Wed Jan 10 11:01:47 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN1

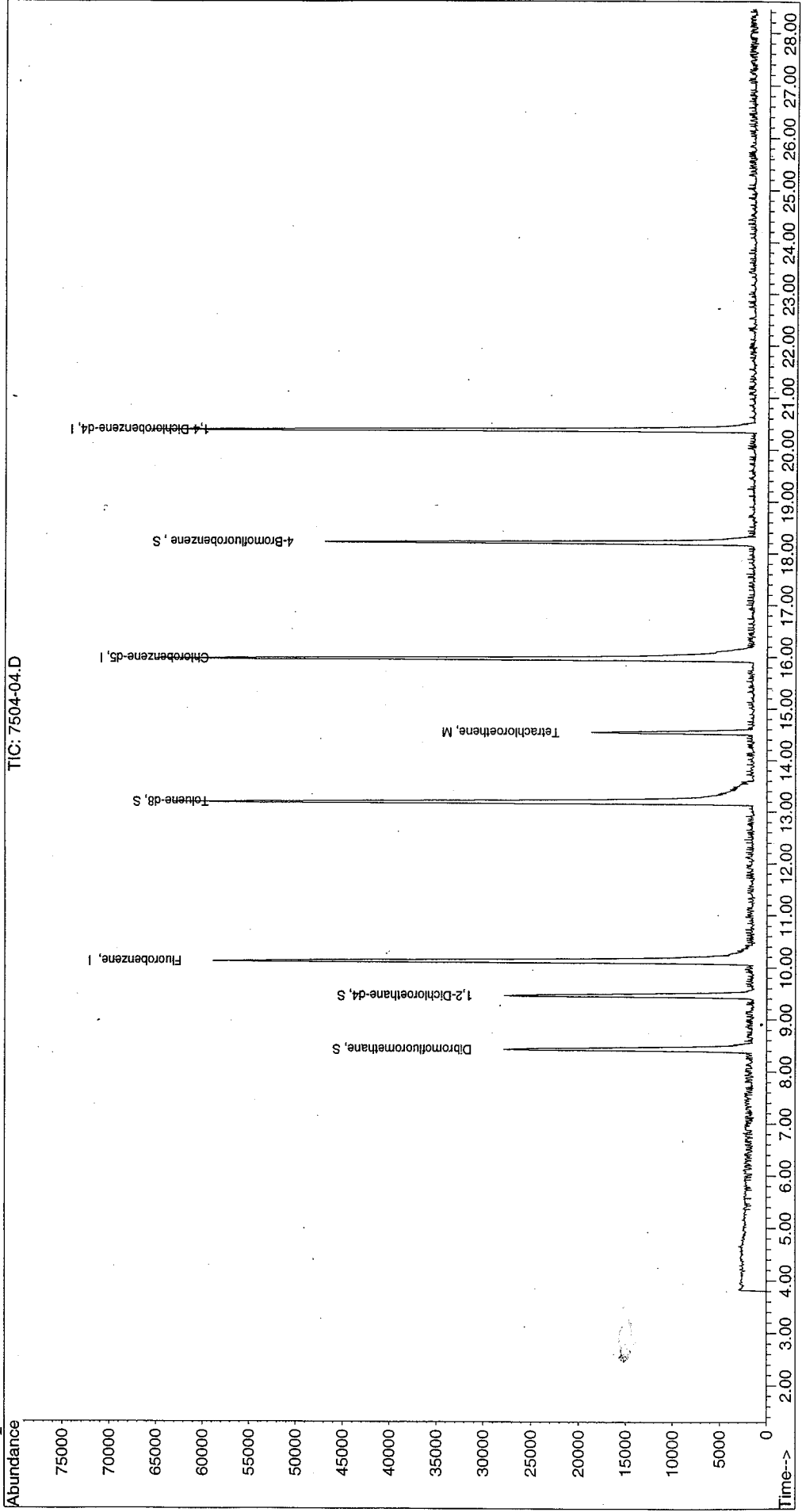
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.55 | 166 | 12777 | 10.46 | ug/L | 96 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration.

Quantitation Report

Data File : C:\HPCHEM\1\DATA\010507\7504-04.D Vial: 34
Acq On : 6 Jan 2007 8:33 am Operator: TS
Sample : 7504_04 0.050 ml Inst : GC/MS #2
Misc : Multiplr: 1.00
MS Integration Params: ODD.P
Quant Time: Jan 17 15:34 2007 Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\010507\7504-05.D

Vial: 35

Acq On : 6 Jan 2007 9:08 am

Operator: TS

Sample : 7504_05 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 17 15:35 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Wed Jan 10 11:01:47 2007

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.12 | 96 | 147269 | 25.00 | ug/L | 0.02 |
| 36) Chlorobenzene-d5 | 15.97 | 117 | 99147 | 25.00 | ug/L | 0.02 |
| 58) 1,4-Dichlorobenzene-d4 | 20.38 | 152 | 39603 | 25.00 | ug/L | 0.02 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|------|
| 16) Dibromofluoromethane | 8.43 | 113 | 41107 | 28.87 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 115.48% | |
| 19) 1,2-Dichloroethane-d4 | 9.46 | 65 | 34009 | 27.60 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 110.40% | |
| 28) Toluene-d8 | 13.19 | 98 | 111722 | 22.64 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 90.56% | |
| 45) 4-Bromofluorobenzene | 18.21 | 95 | 41314 | 22.49 | ug/L | 0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 89.96% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\010507\7504-05.D

Vial: 35

Acq On : 6 Jan 2007 9:08 am

Operator: TS

Sample : 7504_05 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 17 15:35 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Wed Jan 10 11:01:47 2007

Response via : Initial Calibration

DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.53 | 166 | 23512 | 19.78 | ug/L | 97 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7504-05.D 71NV16_6.M

Thu Jan 18 09:24:15 2007

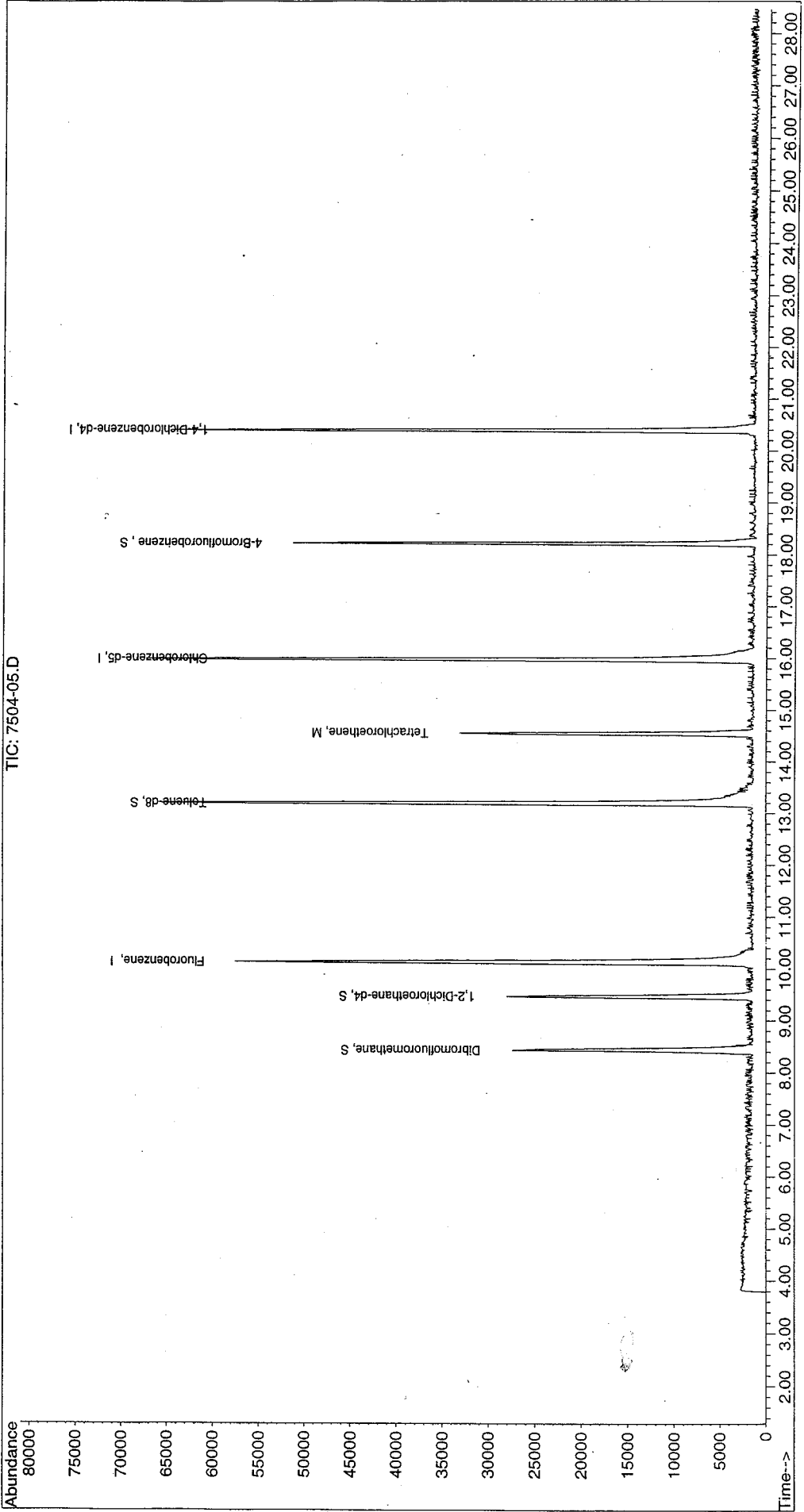
GC#2

Page 2

Quantitation Report

Data File : C:\HPCHEM\1\DATA\010507\7504-05.D
Acq On : 6 Jan 2007 9:08 am
Sample : 7504_05 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 17 15:35 2007
Vial: 35
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\010507\7504-06.D

Vial: 36

Acq On : 6 Jan 2007 9:44 am

Operator: TS

Sample : 7504_06 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 17 15:36 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Wed Jan 10 11:01:47 2007

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.12 | 96 | 144203 | 25.00 | ug/L | 0.02 |
| 36) Chlorobenzene-d5 | 15.96 | 117 | 99154 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 20.38 | 152 | 41549 | 25.00 | ug/L | 0.02 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|------|
| 16) Dibromofluoromethane | 8.43 | 113 | 41336 | 29.65 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 118.60% | |
| 19) 1,2-Dichloroethane-d4 | 9.46 | 65 | 34350 | 28.47 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 113.88% | |
| 28) Toluene-d8 | 13.18 | 98 | 118015 | 24.42 | ug/L | 0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 97.68% | |
| 45) 4-Bromofluorobenzene | 18.21 | 95 | 43432 | 23.64 | ug/L | 0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 94.56% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\010507\7504-06.D

Vial: 36

Acq On : 6 Jan 2007 9:44 am

Operator: TS

Sample : 7504_06 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 17 15:36 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Wed Jan 10 11:01:47 2007

Response via : Initial Calibration

DataAcq Meth : RUN1

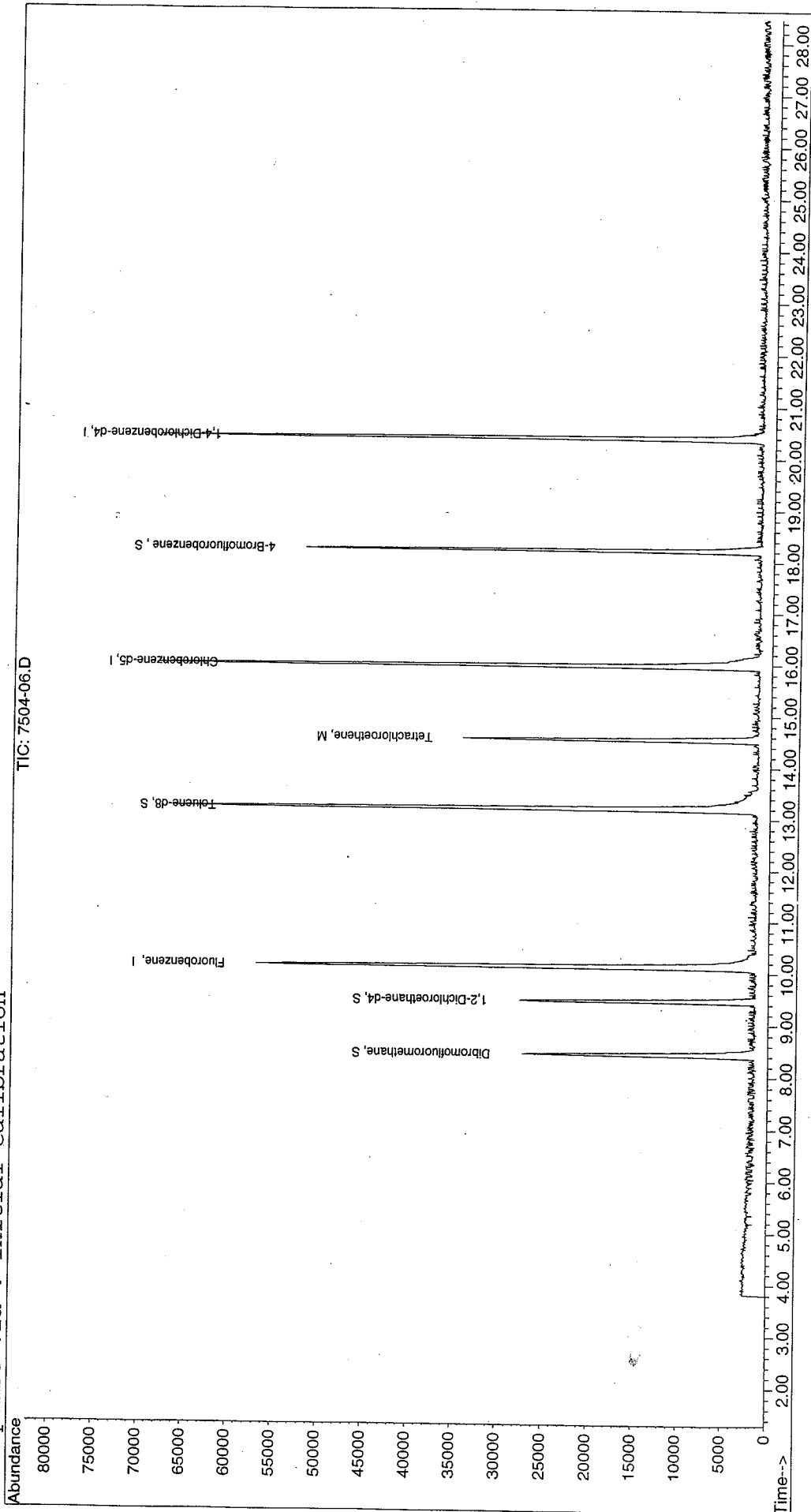
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.54 | 166 | 24704 | 21.22 | ug/L | 96 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\010507\7504-06.D
Acq On : 6 Jan 2007 9:44 am
Sample : 7504_06 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 17 15:36 2007
Vial: 36
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\010507\7504-07.D

Vial: 37

Acq On : 6 Jan 2007 10:19 am

Operator: TS

Sample : 7504_07 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 17 15:36 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Wed Jan 10 11:01:47 2007

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.12 | 96 | 144582 | 25.00 | ug/L | 0.02 |
| 36) Chlorobenzene-d5 | 15.97 | 117 | 93605 | 25.00 | ug/L | 0.02 |
| 58) 1,4-Dichlorobenzene-d4 | 20.38 | 152 | 40239 | 25.00 | ug/L | 0.02 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|------|
| 16) Dibromofluoromethane | 8.43 | 113 | 41384 | 29.61 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 118.44% | |
| 19) 1,2-Dichloroethane-d4 | 9.45 | 65 | 34315 | 28.37 | ug/L | 0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 113.48% | |
| 28) Toluene-d8 | 13.20 | 98 | 112366 | 23.19 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 92.76% | |
| 45) 4-Bromofluorobenzene | 18.21 | 95 | 37742 | 21.76 | ug/L | 0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 87.04% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\010507\7504-07.D

Vial: 37

Acq On : 6 Jan 2007 10:19 am

Operator: TS

Sample : 7504_07 0.050 ml

Inst : GC/MS #2

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 17 15:36 2007

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Wed Jan 10 11:01:47 2007

Response via : Initial Calibration

DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.54 | 166 | 26568 | 22.76 | ug/L | 97 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7504-07.D 71NV16_6.M

Thu Jan 18 09:24:21 2007

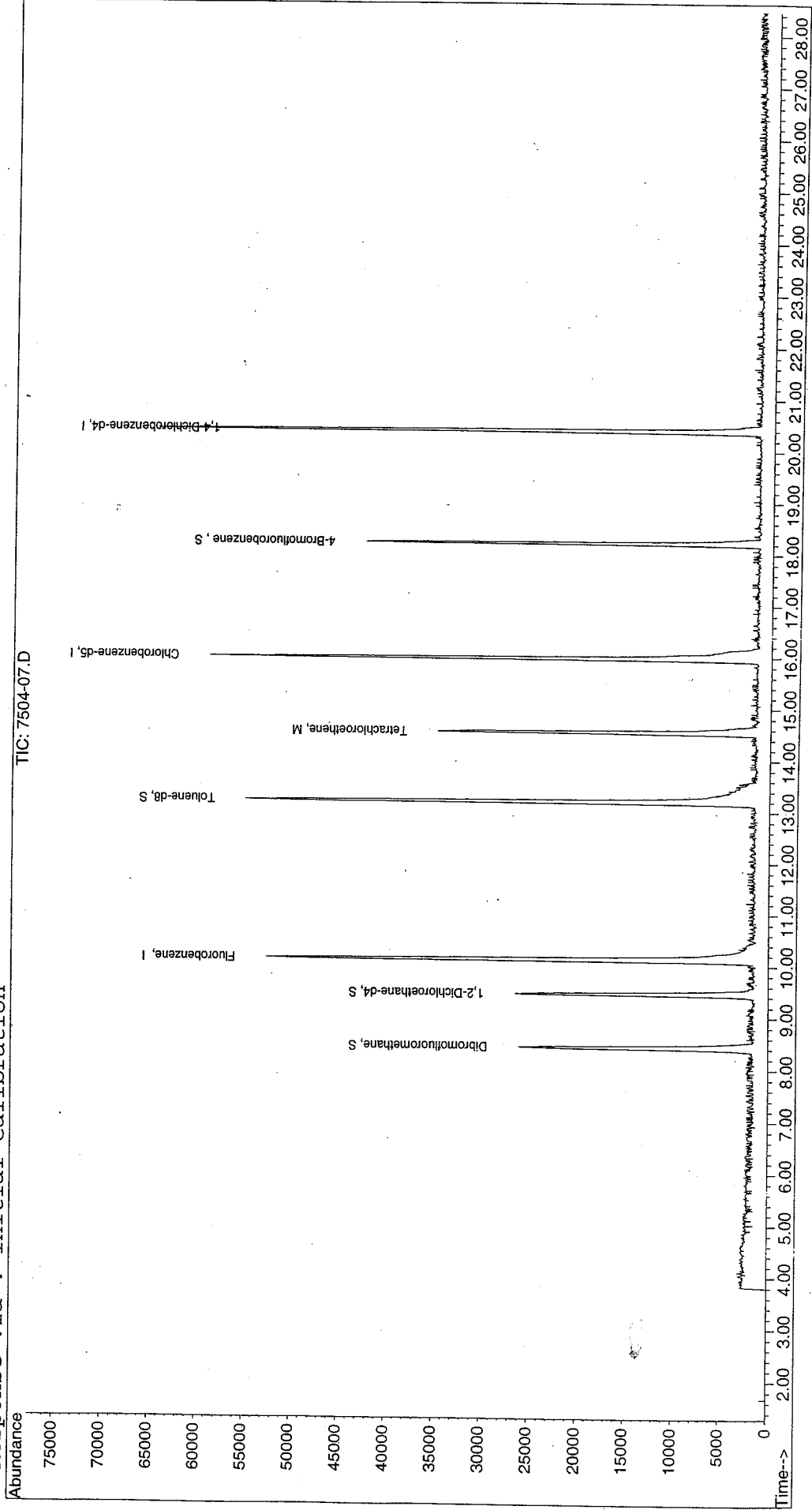
GC#2

Page 2

Quantitation report

Data File : C:\HPCHEM\1\DATA\010507\7504-07.D
Acq On : 6 Jan 2007 10:19 am
Sample : 7504_07 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 17 15:36 2007
Vial: 37
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\010507\7504-08.D
 Acq On : 6 Jan 2007 10:55 am
 Sample : 7504_08 0.050 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 17 15:37 2007

Vial: 38
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Wed Jan 10 11:01:47 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.12 | 96 | 148536 | 25.00 | ug/L | 0.02 |
| 36) Chlorobenzene-d5 | 15.97 | 117 | 102699 | 25.00 | ug/L | 0.02 |
| 58) 1,4-Dichlorobenzene-d4 | 20.38 | 152 | 42471 | 25.00 | ug/L | 0.02 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|------|
| 16) Dibromofluoromethane | 8.43 | 113 | 42157 | 29.36 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 117.44% | |
| 19) 1,2-Dichloroethane-d4 | 9.46 | 65 | 34273 | 27.58 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 110.32% | |
| 28) Toluene-d8 | 13.18 | 98 | 127405 | 25.59 | ug/L | 0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 102.36% | |
| 45) 4-Bromofluorobenzene | 18.21 | 95 | 43678 | 22.95 | ug/L | 0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 91.80% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\O10507\7504-08.D
Acq On : 6 Jan 2007 10:55 am
Sample : 7504_08 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 17 15:37 2007

Vial: 38
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration
DataAcq Meth : RUN1

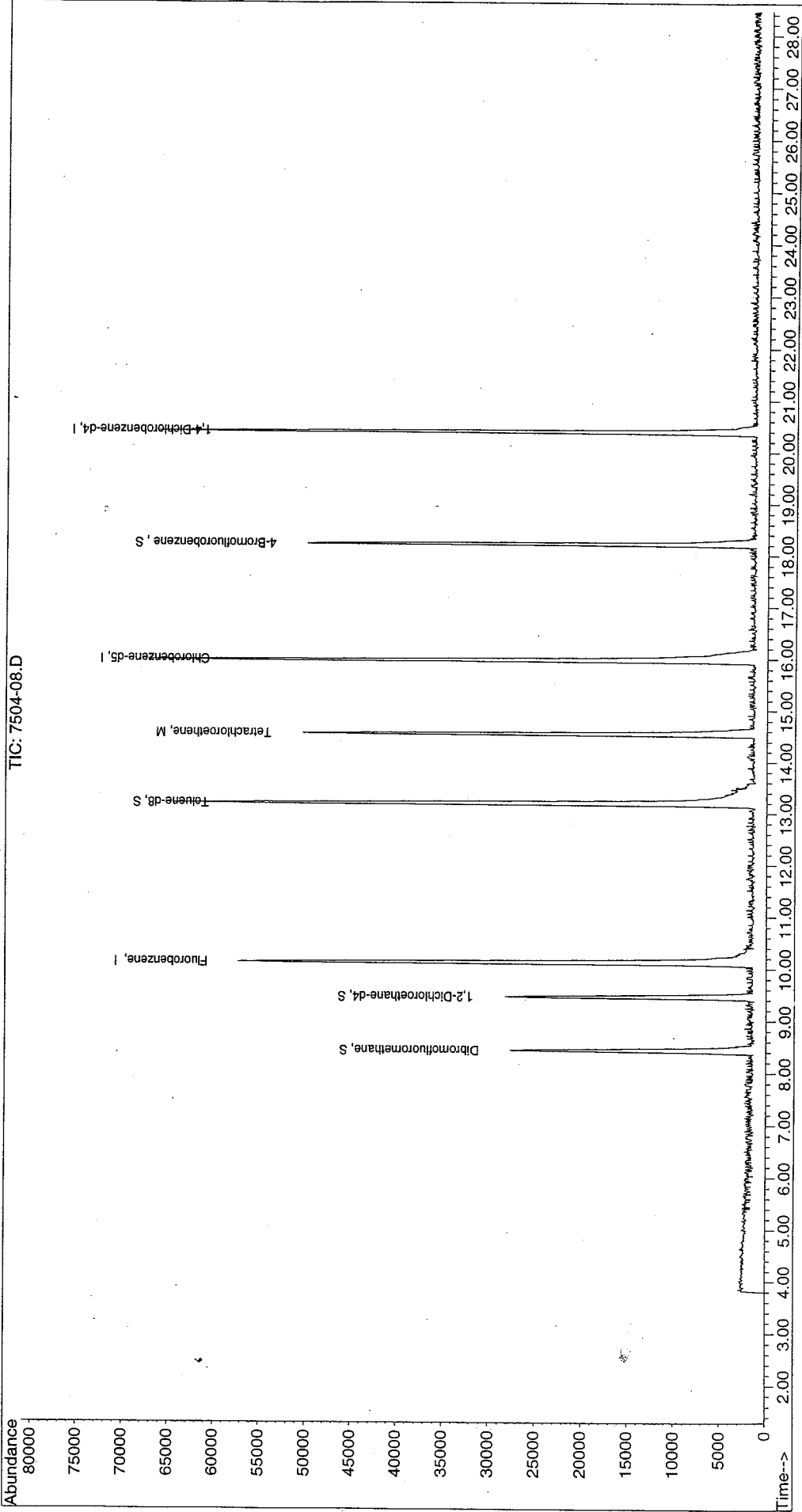
| Compound | R.T. | QIon | Response | Conc Unit | Qvalue |
|---------------------------------|-------|------|----------|------------|--------|
| 32) Tetrachloroethene | 14.54 | 166 | 37474 | 31.26 ug/L | 96 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | |

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\010507\7504-08.D
Acq On : 6 Jan 2007 10:55 am
Sample : 7504_08 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 17 15:37 2007
Vial: 38
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\O10507\7504-09.D
 Acq On : 6 Jan 2007 11:30 am
 Sample : 7504_09 0.050 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 17 15:38 2007

Vial: 39
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Wed Jan 10 11:01:47 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|-------|------|----------|----------|-------|----------|
| 1) Fluorobenzene | 10.12 | 96 | 146438 | 25.00 | ug/L | 0.02 |
| 36) Chlorobenzene-d5 | 15.97 | 117 | 99741 | 25.00 | ug/L | 0.02 |
| 58) 1,4-Dichlorobenzene-d4 | 20.38 | 152 | 40808 | 25.00 | ug/L | 0.02 |
| System Monitoring Compounds | | | | | | |
| 16) Dibromofluoromethane | 8.43 | 113 | 41946 | 29.63 | ug/L | 0.03 |
| Spiked Amount | | | | 25.000 | | |
| | | | | Recovery | = | 118.52% |
| 19) 1,2-Dichloroethane-d4 | 9.47 | 65 | 34162 | 27.88 | ug/L | 0.04 |
| Spiked Amount | | | | 25.000 | | |
| | | | | Recovery | = | 111.52% |
| 28) Toluene-d8 | 13.19 | 98 | 125227 | 25.52 | ug/L | 0.03 |
| Spiked Amount | | | | 25.000 | | |
| | | | | Recovery | = | 102.08% |
| 45) 4-Bromofluorobenzene | 18.22 | 95 | 42318 | 22.90 | ug/L | 0.03 |
| Spiked Amount | | | | 25.000 | | |
| | | | | Recovery | = | 91.60% |
| Target Compounds | | | | | | |
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | Qvalue |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. | d |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration
 7504-09.D 71NV16_6.M Thu Jan 18 09:24:26 2007 GC#2

Data File : C:\HPCHEM\1\DATA\010507\7504-09.D
 Acq On : 6 Jan 2007 11:30 am
 Sample : 7504_09 0.050 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 17 15:38 2007

Vial: 39
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Wed Jan 10 11:01:47 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN1

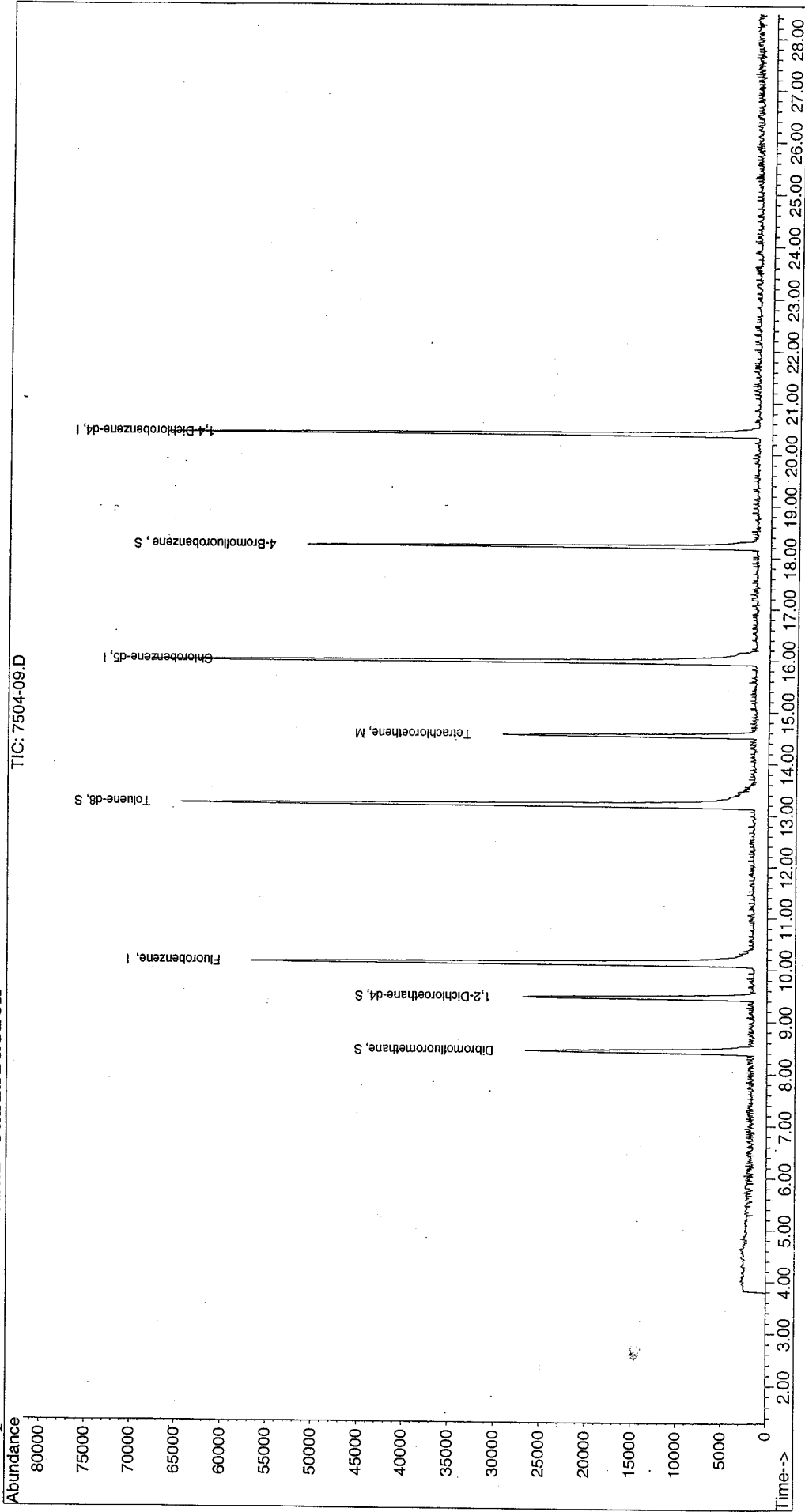
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|--------|--------|
| 32) Tetrachloroethene | 14.53 | 166 | 20340 | 17.21 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | | N.D. | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | | N.D. d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | | N.D. | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | | N.D. | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | | N.D. | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | | N.D. | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | | N.D. | |
| 42) Styrene | 0.00 | 104 | 0 | | N.D. | |
| 43) Bromoform | 0.00 | 173 | 0 | | N.D. | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 46) Bromobenzene | 0.00 | 156 | 0 | | N.D. | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | | N.D. | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | | N.D. | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | | N.D. | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | | N.D. | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | | N.D. | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | | N.D. | |
| 65) Naphthalene | 0.00 | 128 | 0 | | N.D. | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | | N.D. | |
| 68) Acetone | 0.00 | 43 | 0 | | N.D. | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | | N.D. | |
| 70) 2-Butanone | 0.00 | 43 | 0 | | N.D. | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | | N.D. | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | | N.D. | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | | N.D. | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

Quantification report

Data File : C:\HPCHEM\1\DATA\010507\7504-09.D
Acq On : 6 Jan 2007 11:30 am
Sample : 7504_09 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 17 15:38 2007
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\O10507\7504-10.D
 Acq On : 6 Jan 2007 12:05 pm
 Sample : 7504_10 0.050 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 17 15:39 2007

Vial: 40
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Wed Jan 10 11:01:47 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.12 | 96 | 142492 | 25.00 | ug/L | 0.02 |
| 36) Chlorobenzene-d5 | 15.97 | 117 | 94034 | 25.00 | ug/L | 0.02 |
| 58) 1,4-Dichlorobenzene-d4 | 20.38 | 152 | 40242 | 25.00 | ug/L | 0.02 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|------|
| 16) Dibromofluoromethane | 8.43 | 113 | 41808 | 30.35 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 121.40% | |
| 19) 1,2-Dichloroethane-d4 | 9.45 | 65 | 34134 | 28.63 | ug/L | 0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 114.52% | |
| 28) Toluene-d8 | 13.19 | 98 | 114567 | 23.99 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | Recovery | = | 95.96% | |
| 45) 4-Bromofluorobenzene | 18.21 | 95 | 38093 | 21.86 | ug/L | 0.02 |
| Spiked Amount | 25.000 | | Recovery | = | 87.44% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. | d |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\O10507\7504-10.D
 Acq On : 6 Jan 2007 12:05 pm
 Sample : 7504_10 0.050 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 17 15:39 2007

Vial: 40
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Wed Jan 10 11:01:47 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN1

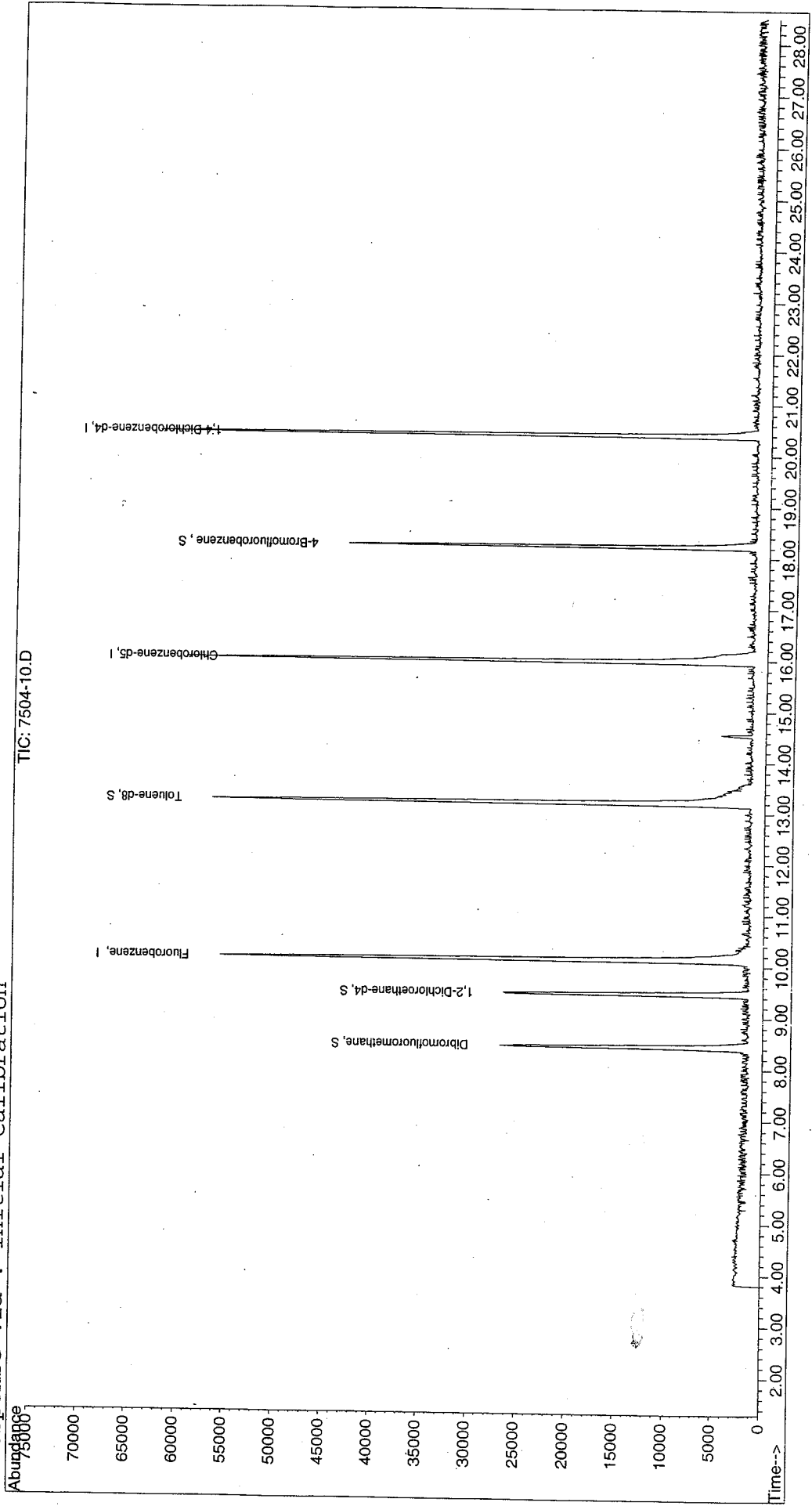
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|------|------|----------|------|------|--------|
| 32) Tetrachloroethene | 0.00 | 166 | 0 | | N.D. | |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | | N.D. | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | | N.D. | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | | N.D. | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | | N.D. | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | | N.D. | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | | N.D. | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | | N.D. | |
| 42) Styrene | 0.00 | 104 | 0 | | N.D. | |
| 43) Bromoform | 0.00 | 173 | 0 | | N.D. | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 46) Bromobenzene | 0.00 | 156 | 0 | | N.D. | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | | N.D. | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | | N.D. | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | | N.D. | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | | N.D. | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | | N.D. | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | | N.D. | |
| 65) Naphthalene | 0.00 | 128 | 0 | | N.D. | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | | N.D. | |
| 68) Acetone | 0.00 | 43 | 0 | | N.D. | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | | N.D. | |
| 70) 2-Butanone | 0.00 | 43 | 0 | | N.D. | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | | N.D. | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | | N.D. | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | | N.D. | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

Quantification report

Data File : C:\HPCHEM\1\DATA\010507\7504-10.D
Acq On : 6 Jan 2007 12:05 pm Vial: 40
Sample : 7504_10 0.050 ml Operator: TS
Misc : Inst : GC/MS #2
MS Integration Params: ODD.P Multiplr: 1.00
Quant Time: Jan 17 15:39 2007 Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\010507\7504-11.D
 Acq On : 6 Jan 2007 12:40 pm
 Sample : 7504_11 0.050 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 17 15:40 2007

Vial: 41
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Wed Jan 10 11:01:47 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.13 | 96 | 142887 | 25.00 | ug/L | 0.03 |
| 36) Chlorobenzene-d5 | 15.97 | 117 | 89766 | 25.00 | ug/L | 0.02 |
| 58) 1,4-Dichlorobenzene-d4 | 20.38 | 152 | 41265 | 25.00 | ug/L | 0.02 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|--------|-------|------------|---------|
| 16) Dibromofluoromethane | 8.43 | 113 | 41211 | 29.83 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | Recovery = | 119.32% |
| 19) 1,2-Dichloroethane-d4 | 9.46 | 65 | 33996 | 28.43 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | Recovery = | 113.72% |
| 28) Toluene-d8 | 13.19 | 98 | 106771 | 22.30 | ug/L | 0.03 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | Recovery = | 89.20% |
| 45) 4-Bromofluorobenzene | 18.21 | 95 | 40341 | 24.25 | ug/L | 0.02 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | Recovery = | 97.00% |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. | d |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\010507\7504-11.D
 Acq On : 6 Jan 2007 12:40 pm
 Sample : 7504_11 0.050 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 17 15:40 2007

Vial: 41
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Wed Jan 10 11:01:47 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN1

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.54 | 166 | 26525 | 23.00 | ug/L | 93 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

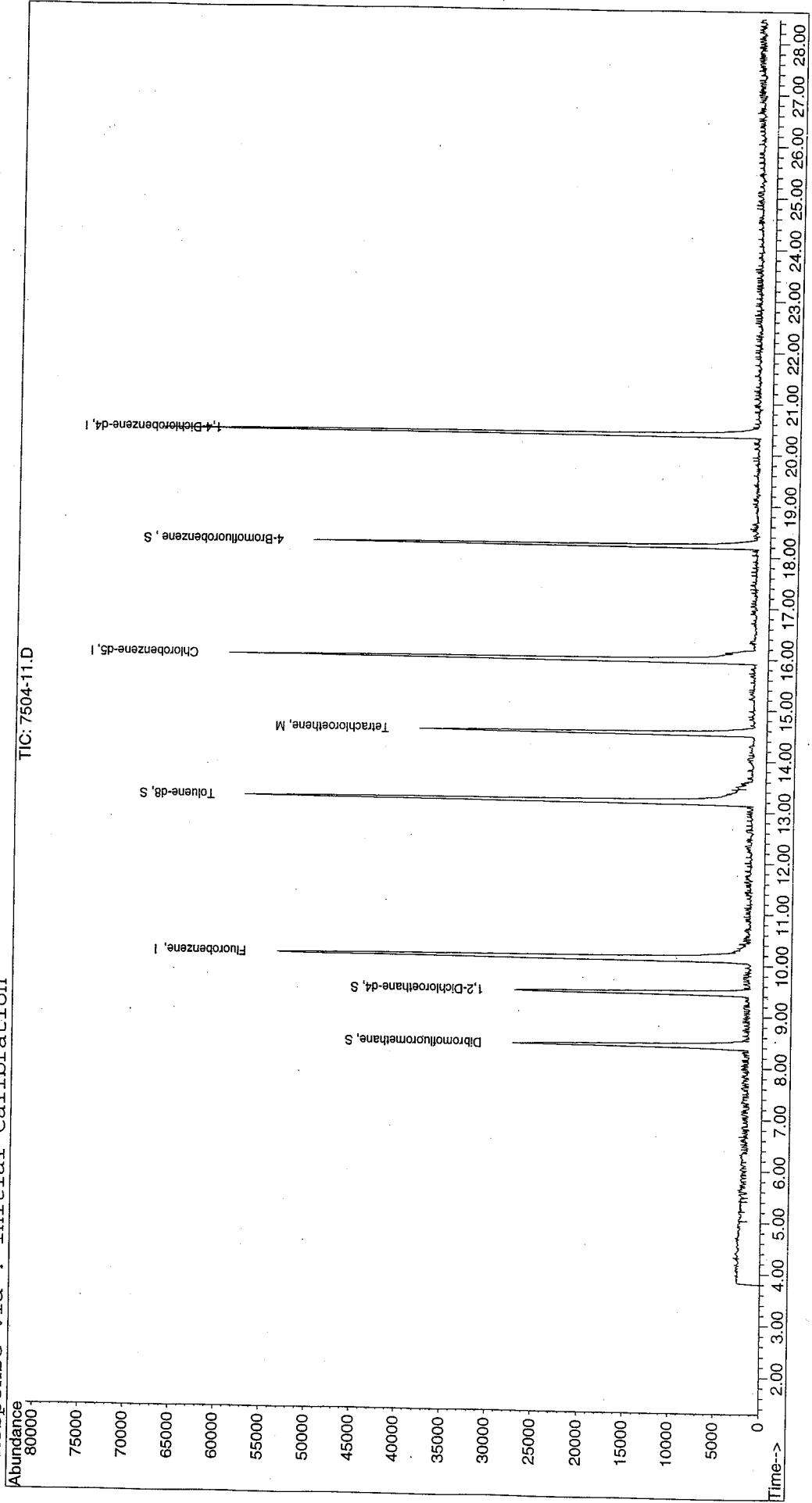
(#) = qualifier out of range (m) = manual integration
 7504-11.D 71NV16_6.M Thu Jan 18 09:24:33 2007

GC#2

Quantitation report

Data File : C:\HPCHEM\1\DATA\010507\7504-11.D
Acq On : 6 Jan 2007 12:40 pm
Sample : 7504_11 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 17 15:40 2007
Vial: 41
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\010507\7504-12.D

Acq On : 6 Jan 2007 1:15 pm

Sample : 7504_12 0.050 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Jan 17 15:40 2007

Vial: 42

Operator: TS

Inst : GC/MS #2

Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Wed Jan 10 11:01:47 2007

Response via : Initial Calibration

DataAcq Meth : RUN1

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 10.12 | 96 | 145528 | 25.00 | ug/L | 0.02 |
| 36) Chlorobenzene-d5 | 15.97 | 117 | 93977 | 25.00 | ug/L | 0.02 |
| 58) 1,4-Dichlorobenzene-d4 | 20.38 | 152 | 39553 | 25.00 | ug/L | 0.02 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|-------|-----|--------|------------|---------|------|
| 16) Dibromofluoromethane | 8.43 | 113 | 41230 | 29.31 | ug/L | 0.03 |
| Spiked Amount | | | | | | |
| | | | | Recovery = | 117.24% | |
| 19) 1,2-Dichloroethane-d4 | 9.46 | 65 | 34273 | 28.15 | ug/L | 0.03 |
| Spiked Amount | | | | | | |
| | | | | Recovery = | 112.60% | |
| 28) Toluene-d8 | 13.19 | 98 | 115035 | 23.59 | ug/L | 0.02 |
| Spiked Amount | | | | | | |
| | | | | Recovery = | 94.36% | |
| 45) 4-Bromofluorobenzene | 18.21 | 95 | 37429 | 21.49 | ug/L | 0.02 |
| Spiked Amount | | | | | | |
| | | | | Recovery = | 85.96% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7504-12.D 71NV16_6.M

Thu Jan 18 09:24:36 2007

GC#2

Page 1

Data File : C:\HPCHEM\1\DATA\010507\7504-12.D
 Acq On : 6 Jan 2007 1:15 pm
 Sample : 7504_12 0.050 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 17 15:40 2007

Vial: 42
 Operator: TS
 Inst : GC/MS #2
 Multiplr: 1.00

Quant Results File: 71NV16_6.RES

Quant Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Wed Jan 10 11:01:47 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN1

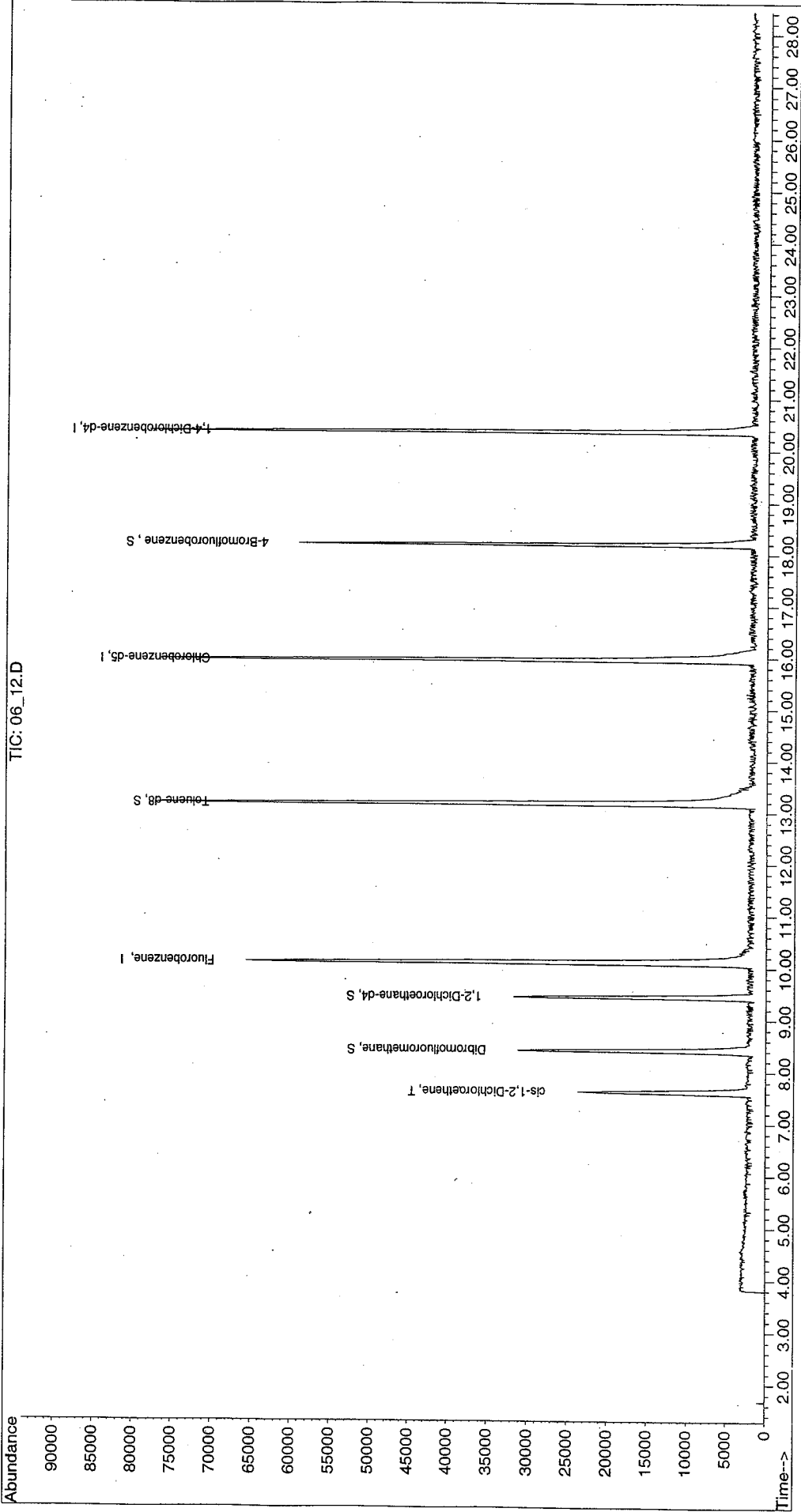
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 14.54 | 166 | 22337 | 19.02 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromoethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration
 7504-12.D 71NV16_6.M Thu Jan 18 09:24:36 2007

Quantitation Report

Data File : C:\HPCHEM\1\DATA\010507\06_12.D
Acq On : 5 Jan 2007 8:49 pm
Sample : 7506-12 0.40 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 17 16:31 2007
Vial: 14
Operator: TS
Inst : GC/MS #2
Multiplr: 1.00
Quant Results File: 71NV16_6.RES

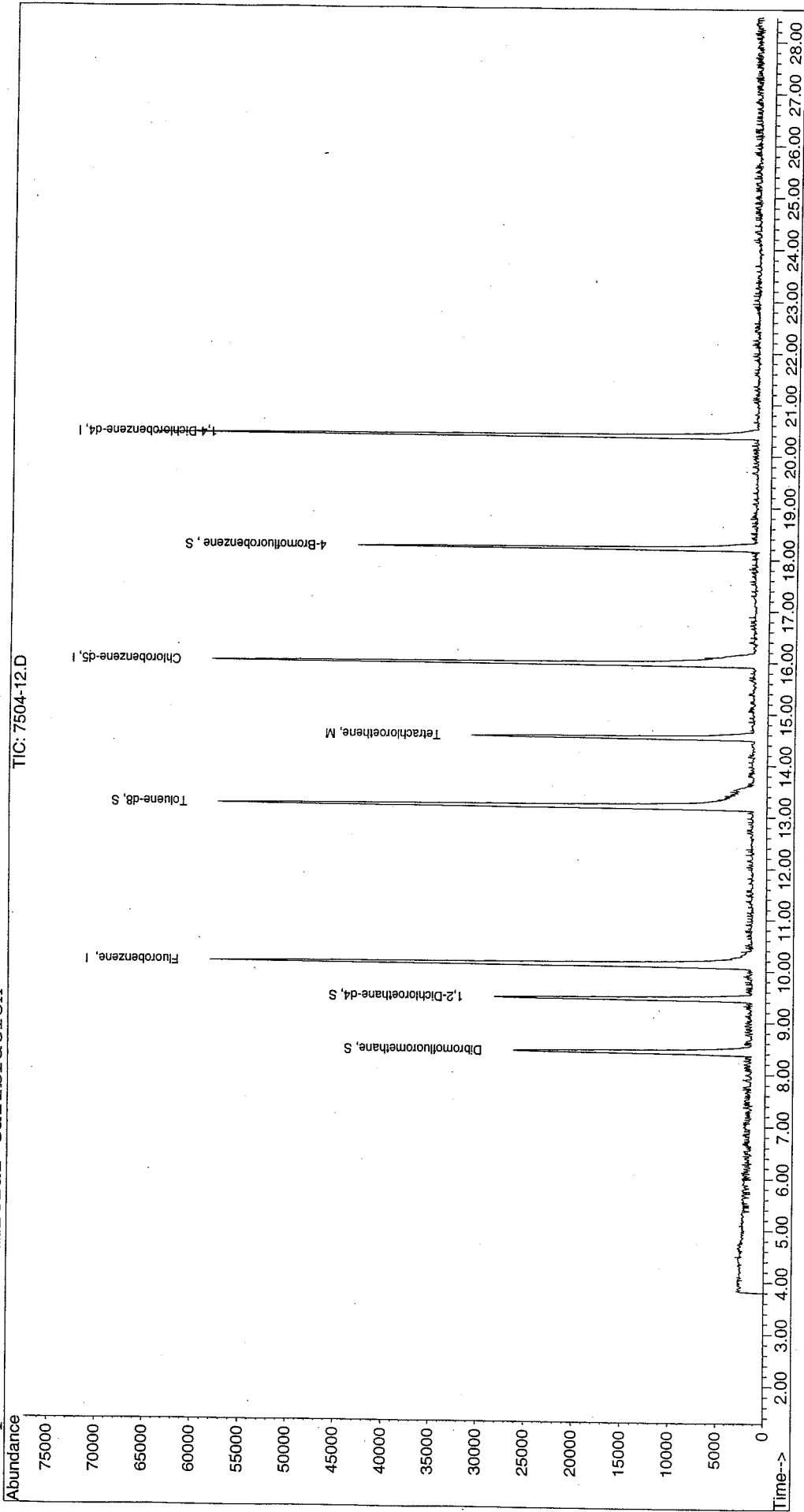
Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration

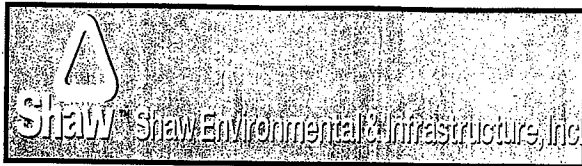


Validation report

Data File : C:\HPCHEM\1\DATA\010507\7504-12.D
Acq On : 6 Jan 2007 1:15 pm
Sample : 7504_12 0.050 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 17 15:40 2007
Quant Results File: 71NV16_6.RES

Method : C:\HPCHEM\1\METHODS\71NV16_6.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Wed Jan 10 11:01:47 2007
Response via : Initial Calibration





17 Princess Rd
Lawrenceville, New Jersey 08648
Tel: 609/895-5370
Fax: 609/895-1858

Volatile Organic Compound Data Summary Package

Prepared for
Gorham Textron

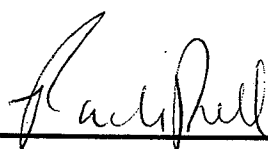
Lab ID
7519

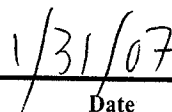
Project Number 101960 02000000

Samples Received
18-Jan-07

Reported
31-Jan-07

NJDEP Certified Lab 11001


Randi K Rothmel, PhD
Laboratory Director


Date

1.0 Chain of Custody

2.0. Sample Summaries

Sample summaries are enclosed

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-1

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519-1 0.020ml
 Sample wt/vol: 0.020 (g/ml) ML Lab File ID: 7519_1.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 2500 | U |
| 74-87-3 | Chloromethane | | 2500 | U |
| 75-01-4 | Vinyl chloride | | 2500 | U |
| 74-83-9 | Bromomethane | | 2500 | U |
| 75-00-3 | Chloroethane | | 2500 | U |
| 75-69-4 | Trichlorofluoromethane | | 2500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 2500 | U |
| 75-09-2 | Methylene chloride | | 2500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 2500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 2500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 2500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 2500 | U |
| 74-97-5 | Bromochloromethane | | 2500 | U |
| 67-66-3 | Chloroform | | 2500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 2500 | U |
| 56-23-5 | Carbon tetrachloride | | 2500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 2500 | U |
| 71-43-2 | Benzene | | 2500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 2500 | U |
| 79-01-6 | Trichloroethene | | 2500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 2500 | U |
| 74-95-3 | Dibromomethane | | 2500 | U |
| 75-27-4 | Bromodichloromethane | | 2500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 2500 | U |
| 108-88-3 | Toluene | | 2500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 2500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 2500 | U |
| 127-18-4 | Tetrachloroethene | | 27000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 2500 | U |
| 124-48-1 | Dibromochloromethane | | 2500 | U |
| 108-90-7 | Chlorobenzene | | 2500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 2500 | U |
| 100-41-4 | Ethylbenzene | | 2500 | U |
| 1330-20-7 | Xylene (para & meta) | | 2500 | U |
| 95-47-6 | Xylene (Ortho) | | 2500 | U |
| 100-42-5 | Styrene | | 2500 | U |
| 75-25-2 | Bromoform | | 2500 | U |
| 98-82-8 | Isopropylbenzene | | 2500 | U |
| 108-86-1 | Bromobenzene | | 2500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-1

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519-1 0.020ml
 Sample wt/vol: 0.020 (g/ml) ML Lab File ID: 7519_1.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 2500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 2500 | U |
| 103-65-1 | n-Propylbenzene | | 2500 | U |
| 95-49-8 | 2-Chlorotoluene | | 2500 | U |
| 106-43-4 | 4-Chlorotoluene | | 2500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 2500 | U |
| 98-06-6 | tert-Butylbenzene | | 2500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 2500 | U |
| 135-98-8 | sec-Butylbenzene | | 2500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 2500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 2500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 2500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 2500 | U |
| 104-51-8 | n-Butylbenzene | | 2500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 2500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 2500 | U |
| 87-68-3 | Hexachlorobutadiene | | 2500 | U |
| 91-20-3 | Naphthalene | | 2500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 2500 | U |
| 1634-04-4 | MTBE | | 2500 | U |
| 67-64-1 | Acetone | | 5000 | U |
| 75-15-0 | Carbon disulfide | | 2500 | U |
| 78-93-3 | 2-Butanone | | 5000 | U |
| 109-99-9 | Tetrahydrofuran | | 5000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 5000 | U |
| 591-78-6 | 2-Hexanone | | 5000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 5000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-2

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519-2 0.020ml
 Sample wt/vol: 0.020 (g/ml) ML Lab File ID: 7519_2.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 2500 | U |
| 74-87-3 | Chloromethane | | 2500 | U |
| 75-01-4 | Vinyl chloride | | 2500 | U |
| 74-83-9 | Bromomethane | | 2500 | U |
| 75-00-3 | Chloroethane | | 2500 | U |
| 75-69-4 | Trichlorofluoromethane | | 2500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 2500 | U |
| 75-09-2 | Methylene chloride | | 2500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 2500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 2500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 2500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 2500 | U |
| 74-97-5 | Bromochloromethane | | 2500 | U |
| 67-66-3 | Chloroform | | 2500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 2500 | U |
| 56-23-5 | Carbon tetrachloride | | 2500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 2500 | U |
| 71-43-2 | Benzene | | 2500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 2500 | U |
| 79-01-6 | Trichloroethene | | 2500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 2500 | U |
| 74-95-3 | Dibromomethane | | 2500 | U |
| 75-27-4 | Bromodichloromethane | | 2500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 2500 | U |
| 108-88-3 | Toluene | | 2500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 2500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 2500 | U |
| 127-18-4 | Tetrachloroethene | | 28000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 2500 | U |
| 124-48-1 | Dibromochloromethane | | 2500 | U |
| 108-90-7 | Chlorobenzene | | 2500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 2500 | U |
| 100-41-4 | Ethylbenzene | | 2500 | U |
| 1330-20-7 | Xylene (para & meta) | | 2500 | U |
| 95-47-6 | Xylene (Ortho) | | 2500 | U |
| 100-42-5 | Styrene | | 2500 | U |
| 75-25-2 | Bromoform | | 2500 | U |
| 98-82-8 | Isopropylbenzene | | 2500 | U |
| 108-86-1 | Bromobenzene | | 2500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-2

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519-2 0.020ml
 Sample wt/vol: 0.020 (g/ml) ML Lab File ID: 7519_2.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

| CAS NO. | COMPOUND | CONCENTRATION UNITS: | | Q |
|-----------|-----------------------------|----------------------|------|---|
| | | (ug/L or ug/Kg) | UG/L | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 2500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 2500 | U |
| 103-65-1 | n-Propylbenzene | | 2500 | U |
| 95-49-8 | 2-Chlorotoluene | | 2500 | U |
| 106-43-4 | 4-Chlorotoluene | | 2500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 2500 | U |
| 98-06-6 | tert-Butylbenzene | | 2500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 2500 | U |
| 135-98-8 | sec-Butylbenzene | | 2500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 2500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 2500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 2500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 2500 | U |
| 104-51-8 | n-Butylbenzene | | 2500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 2500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 2500 | U |
| 87-68-3 | Hexachlorobutadiene | | 2500 | U |
| 91-20-3 | Naphthalene | | 2500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 2500 | U |
| 1634-04-4 | MTBE | | 2500 | U |
| 67-64-1 | Acetone | | 5000 | U |
| 75-15-0 | Carbon disulfide | | 2500 | U |
| 78-93-3 | 2-Butanone | | 5000 | U |
| 109-99-9 | Tetrahydrofuran | | 5000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 5000 | U |
| 591-78-6 | 2-Hexanone | | 5000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 5000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-3

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519-3 0.020ml
 Sample wt/vol: 0.020 (g/ml) ML Lab File ID: 7519_3.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 2500 | U |
| 74-87-3 | Chloromethane | | 2500 | U |
| 75-01-4 | Vinyl chloride | | 2500 | U |
| 74-83-9 | Bromomethane | | 2500 | U |
| 75-00-3 | Chloroethane | | 2500 | U |
| 75-69-4 | Trichlorofluoromethane | | 2500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 2500 | U |
| 75-09-2 | Methylene chloride | | 2500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 2500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 2500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 2500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 2500 | U |
| 74-97-5 | Bromochloromethane | | 2500 | U |
| 67-66-3 | Chloroform | | 2500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 2500 | U |
| 56-23-5 | Carbon tetrachloride | | 2500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 2500 | U |
| 71-43-2 | Benzene | | 2500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 2500 | U |
| 79-01-6 | Trichloroethene | | 2500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 2500 | U |
| 74-95-3 | Dibromomethane | | 2500 | U |
| 75-27-4 | Bromodichloromethane | | 2500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 2500 | U |
| 108-88-3 | Toluene | | 2500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 2500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 2500 | U |
| 127-18-4 | Tetrachloroethene | | 26000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 2500 | U |
| 124-48-1 | Dibromochloromethane | | 2500 | U |
| 108-90-7 | Chlorobenzene | | 2500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 2500 | U |
| 100-41-4 | Ethylbenzene | | 2500 | U |
| 1330-20-7 | Xylene (para & meta) | | 2500 | U |
| 95-47-6 | Xylene (Ortho) | | 2500 | U |
| 100-42-5 | Styrene | | 2500 | U |
| 75-25-2 | Bromoform | | 2500 | U |
| 98-82-8 | Isopropylbenzene | | 2500 | U |
| 108-86-1 | Bromobenzene | | 2500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-3

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519-3 0.020ml
 Sample wt/vol: 0.020 (g/ml) ML Lab File ID: 7519_3.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 2500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 2500 | U |
| 103-65-1 | n-Propylbenzene | | 2500 | U |
| 95-49-8 | 2-Chlorotoluene | | 2500 | U |
| 106-43-4 | 4-Chlorotoluene | | 2500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 2500 | U |
| 98-06-6 | tert-Butylbenzene | | 2500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 2500 | U |
| 135-98-8 | sec-Butylbenzene | | 2500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 2500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 2500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 2500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 2500 | U |
| 104-51-8 | n-Butylbenzene | | 2500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 2500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 2500 | U |
| 87-68-3 | Hexachlorobutadiene | | 2500 | U |
| 91-20-3 | Naphthalene | | 2500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 2500 | U |
| 1634-04-4 | MTBE | | 2500 | U |
| 67-64-1 | Acetone | | 5000 | U |
| 75-15-0 | Carbon disulfide | | 2500 | U |
| 78-93-3 | 2-Butanone | | 5000 | U |
| 109-99-9 | Tetrahydrofuran | | 5000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 5000 | U |
| 591-78-6 | 2-Hexanone | | 5000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 5000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-4

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519-4 0.020ml
 Sample wt/vol: 0.020 (g/ml) ML Lab File ID: 7519_4.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 2500 | U |
| 74-87-3 | Chloromethane | | 2500 | U |
| 75-01-4 | Vinyl chloride | | 2500 | U |
| 74-83-9 | Bromomethane | | 2500 | U |
| 75-00-3 | Chloroethane | | 2500 | U |
| 75-69-4 | Trichlorofluoromethane | | 2500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 2500 | U |
| 75-09-2 | Methylene chloride | | 2500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 2500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 2500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 2500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 2500 | U |
| 74-97-5 | Bromochloromethane | | 2500 | U |
| 67-66-3 | Chloroform | | 2500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 2500 | U |
| 56-23-5 | Carbon tetrachloride | | 2500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 2500 | U |
| 71-43-2 | Benzene | | 2500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 2500 | U |
| 79-01-6 | Trichloroethene | | 2500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 2500 | U |
| 74-95-3 | Dibromomethane | | 2500 | U |
| 75-27-4 | Bromodichloromethane | | 2500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 2500 | U |
| 108-88-3 | Toluene | | 2500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 2500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 2500 | U |
| 127-18-4 | Tetrachloroethene | | 16000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 2500 | U |
| 124-48-1 | Dibromochloromethane | | 2500 | U |
| 108-90-7 | Chlorobenzene | | 2500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 2500 | U |
| 100-41-4 | Ethylbenzene | | 2500 | U |
| 1330-20-7 | Xylene (para & meta) | | 2500 | U |
| 95-47-6 | Xylene (Ortho) | | 2500 | D |
| 100-42-5 | Styrene | | 2500 | U |
| 75-25-2 | Bromoform | | 2500 | U |
| 98-82-8 | Isopropylbenzene | | 2500 | U |
| 108-86-1 | Bromobenzene | | 2500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-4

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519-4 0.020ml
 Sample wt/vol: 0.020 (g/ml) ML Lab File ID: 7519_4.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 2500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 2500 | U |
| 103-65-1 | n-Propylbenzene | | 2500 | U |
| 95-49-8 | 2-Chlorotoluene | | 2500 | U |
| 106-43-4 | 4-Chlorotoluene | | 2500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 2500 | U |
| 98-06-6 | tert-Butylbenzene | | 2500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 2500 | U |
| 135-98-8 | sec-Butylbenzene | | 2500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 2500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 2500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 2500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 2500 | U |
| 104-51-8 | n-Butylbenzene | | 2500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 2500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 2500 | U |
| 87-68-3 | Hexachlorobutadiene | | 2500 | U |
| 91-20-3 | Naphthalene | | 2500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 2500 | U |
| 1634-04-4 | MTBE | | 2500 | U |
| 67-64-1 | Acetone | | 5000 | U |
| 75-15-0 | Carbon disulfide | | 2500 | U |
| 78-93-3 | 2-Butanone | | 5000 | U |
| 109-99-9 | Tetrahydrofuran | | 5000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 5000 | U |
| 591-78-6 | 2-Hexanone | | 5000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 5000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-5

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519-5 0.020ml
 Sample wt/vol: 0.020 (g/ml) ML Lab File ID: 7519_5.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 2500 | U |
| 74-87-3 | Chloromethane | | 2500 | U |
| 75-01-4 | Vinyl chloride | | 2500 | U |
| 74-83-9 | Bromomethane | | 2500 | U |
| 75-00-3 | Chloroethane | | 2500 | U |
| 75-69-4 | Trichlorofluoromethane | | 2500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 2500 | U |
| 75-09-2 | Methylene chloride | | 2500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 2500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 2500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 2500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 2500 | U |
| 74-97-5 | Bromochloromethane | | 2500 | U |
| 67-66-3 | Chloroform | | 2500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 2500 | U |
| 56-23-5 | Carbon tetrachloride | | 2500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 2500 | U |
| 71-43-2 | Benzene | | 2500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 2500 | U |
| 79-01-6 | Trichloroethene | | 2500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 2500 | U |
| 74-95-3 | Dibromomethane | | 2500 | U |
| 75-27-4 | Bromodichloromethane | | 2500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 2500 | U |
| 108-88-3 | Toluene | | 2500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 2500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 2500 | U |
| 127-18-4 | Tetrachloroethene | | 22000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 2500 | U |
| 124-48-1 | Dibromochloromethane | | 2500 | U |
| 108-90-7 | Chlorobenzene | | 2500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 2500 | U |
| 100-41-4 | Ethylbenzene | | 2500 | U |
| 1330-20-7 | Xylene (para & meta) | | 2500 | U |
| 95-47-6 | Xylene (Ortho) | | 2500 | U |
| 100-42-5 | Styrene | | 2500 | U |
| 75-25-2 | Bromoform | | 2500 | U |
| 98-82-8 | Isopropylbenzene | | 2500 | U |
| 108-86-1 | Bromobenzene | | 2500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-5

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519-5 0.020ml
 Sample wt/vol: 0.020 (g/ml) ML Lab File ID: 7519_5.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 2500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 2500 | U |
| 103-65-1 | n-Propylbenzene | | 2500 | U |
| 95-49-8 | 2-Chlorotoluene | | 2500 | U |
| 106-43-4 | 4-Chlorotoluene | | 2500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 2500 | U |
| 98-06-6 | tert-Butylbenzene | | 2500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 2500 | U |
| 135-98-8 | sec-Butylbenzene | | 2500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 2500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 2500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 2500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 2500 | U |
| 104-51-8 | n-Butylbenzene | | 2500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 2500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 2500 | U |
| 87-68-3 | Hexachlorobutadiene | | 2500 | U |
| 91-20-3 | Naphthalene | | 2500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 2500 | U |
| 1634-04-4 | MTBE | | 2500 | U |
| 67-64-1 | Acetone | | 5000 | U |
| 75-15-0 | Carbon disulfide | | 2500 | U |
| 78-93-3 | 2-Butanone | | 5000 | U |
| 109-99-9 | Tetrahydrofuran | | 5000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 5000 | U |
| 591-78-6 | 2-Hexanone | | 5000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 5000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-6

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519-6 0.020ml
 Sample wt/vol: 0.020 (g/ml) ML Lab File ID: 7519_6.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-----------------------|---|
| 75-71-8 | Dichlorodifluoromethane | | 2500 | U |
| 74-87-3 | Chloromethane | | 2500 | U |
| 75-01-4 | Vinyl chloride | | 2500 | U |
| 74-83-9 | Bromomethane | | 2500 | U |
| 75-00-3 | Chloroethane | | 2500 | U |
| 75-69-4 | Trichlorofluoromethane | | 2500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 2500 | U |
| 75-09-2 | Methylene chloride | | 2500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 2500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 2500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 2500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 2500 | U |
| 74-97-5 | Bromochloromethane | | 2500 | U |
| 67-66-3 | Chloroform | | 2500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 2500 | U |
| 56-23-5 | Carbon tetrachloride | | 2500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 2500 | U |
| 71-43-2 | Benzene | | 2500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 2500 | U |
| 79-01-6 | Trichloroethene | | 2500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 2500 | U |
| 74-95-3 | Dibromomethane | | 2500 | U |
| 75-27-4 | Bromodichloromethane | | 2500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 2500 | U |
| 108-88-3 | Toluene | | 2500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 2500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 2500 | U |
| 127-18-4 | Tetrachloroethene | | 18000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 2500 | U |
| 124-48-1 | Dibromochloromethane | | 2500 24000 | D |
| 108-90-7 | Chlorobenzene | | 2500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 2500 | U |
| 100-41-4 | Ethylbenzene | | 2500 | U |
| 1330-20-7 | Xylene (para & meta) | | 2500 | U |
| 95-47-6 | Xylene (Ortho) | | 2500 | U |
| 100-42-5 | Styrene | | 2500 | U |
| 75-25-2 | Bromoform | | 2500 | U |
| 98-82-8 | Isopropylbenzene | | 2500 | U |
| 108-86-1 | Bromobenzene | | 2500 | U |

- NOT a peak
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1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-6

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519-6 0.020ml
 Sample wt/vol: 0.020 (g/ml) ML Lab File ID: 7519_6.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 2500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 2500 | U |
| 103-65-1 | n-Propylbenzene | | 2500 | U |
| 95-49-8 | 2-Chlorotoluene | | 2500 | U |
| 106-43-4 | 4-Chlorotoluene | | 2500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 2500 | U |
| 98-06-6 | tert-Butylbenzene | | 2500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 2500 | U |
| 135-98-8 | sec-Butylbenzene | | 2500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 2500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 2500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 2500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 2500 | U |
| 104-51-8 | n-Butylbenzene | | 2500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 2500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 2500 | U |
| 87-68-3 | Hexachlorobutadiene | | 2500 | U |
| 91-20-3 | Naphthalene | | 2500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 2500 | U |
| 1634-04-4 | MTBE | | 2500 | U |
| 67-64-1 | Acetone | | 5000 | U |
| 75-15-0 | Carbon disulfide | | 2500 | U |
| 78-93-3 | 2-Butanone | | 5000 | U |
| 109-99-9 | Tetrahydrofuran | | 5000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 5000 | U |
| 591-78-6 | 2-Hexanone | | 5000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 5000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-07

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_07 0.020 ml
 Sample wt/vol: 0.0 (g/ml) ML Lab File ID: 7519_07.D
 Level: (low/med) LOW Date Received: 01/11/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 2500 | U |
| 74-87-3 | Chloromethane | | 2500 | U |
| 75-01-4 | Vinyl chloride | | 2500 | U |
| 74-83-9 | Bromomethane | | 2500 | U |
| 75-00-3 | Chloroethane | | 2500 | U |
| 75-69-4 | Trichlorofluoromethane | | 2500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 2500 | U |
| 75-09-2 | Methylene chloride | | 2500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 2500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 2500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 2500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 2500 | U |
| 74-97-5 | Bromochloromethane | | 2500 | U |
| 67-66-3 | Chloroform | | 2500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 2500 | U |
| 56-23-5 | Carbon tetrachloride | | 2500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 2500 | U |
| 71-43-2 | Benzene | | 2500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 2500 | U |
| 79-01-6 | Trichloroethene | | 2500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 2500 | U |
| 74-95-3 | Dibromomethane | | 2500 | U |
| 75-27-4 | Bromodichloromethane | | 2500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 2500 | U |
| 108-88-3 | Toluene | | 2500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 2500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 2500 | U |
| 127-18-4 | Tetrachloroethene | | 24000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 2500 | U |
| 124-48-1 | Dibromochloromethane | | 2500 | U |
| 108-90-7 | Chlorobenzene | | 2500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 2500 | U |
| 100-41-4 | Ethylbenzene | | 2500 | U |
| 1330-20-7 | Xylene (para & meta) | | 2500 | U |
| 95-47-6 | Xylene (Ortho) | | 2500 | U |
| 100-42-5 | Styrene | | 2500 | U |
| 75-25-2 | Bromoform | | 2500 | U |
| 98-82-8 | Isopropylbenzene | | 2500 | U |
| 108-86-1 | Bromobenzene | | 2500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-07

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_07 0.020 ml
 Sample wt/vol: 0.0 (g/ml) ML Lab File ID: 7519_07.D
 Level: (low/med) LOW Date Received: 01/11/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 2500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 2500 | U |
| 103-65-1 | n-Propylbenzene | | 2500 | U |
| 95-49-8 | 2-Chlorotoluene | | 2500 | U |
| 106-43-4 | 4-Chlorotoluene | | 2500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 2500 | U |
| 98-06-6 | tert-Butylbenzene | | 2500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 2500 | U |
| 135-98-8 | sec-Butylbenzene | | 2500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 2500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 2500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 2500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 2500 | U |
| 104-51-8 | n-Butylbenzene | | 2500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 2500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 2500 | U |
| 87-68-3 | Hexachlorobutadiene | | 2500 | U |
| 91-20-3 | Naphthalene | | 2500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 2500 | U |
| 1634-04-4 | MTBE | | 2500 | U |
| 67-64-1 | Acetone | | 5000 | U |
| 75-15-0 | Carbon disulfide | | 2500 | U |
| 78-93-3 | 2-Butanone | | 5000 | U |
| 109-99-9 | Tetrahydrofuran | | 5000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 5000 | U |
| 591-78-6 | 2-Hexanone | | 5000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 5000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-8

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA.
 Matrix: (soil/water) WATER Lab Sample ID: 7519_8 0.020 ml
 Sample wt/vol: 0.0 (g/ml) ML Lab File ID: 7519_8.D
 Level: (low/med) LOW Date Received: 01/11/05
 % Moisture: not dec: _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 2500 | U |
| 74-87-3 | Chloromethane | | 2500 | U |
| 75-01-4 | Vinyl chloride | | 2500 | U |
| 74-83-9 | Bromomethane | | 2500 | U |
| 75-00-3 | Chloroethane | | 2500 | U |
| 75-69-4 | Trichlorofluoromethane | | 2500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 2500 | U |
| 75-09-2 | Methylene chloride | | 2500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 2500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 2500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 2500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 2500 | U |
| 74-97-5 | Bromochloromethane | | 2500 | U |
| 67-66-3 | Chloroform | | 2500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 2500 | U |
| 56-23-5 | Carbon tetrachloride | | 2500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 2500 | U |
| 71-43-2 | Benzene | | 2500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 2500 | U |
| 79-01-6 | Trichloroethene | | 2500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 2500 | U |
| 74-95-3 | Dibromomethane | | 2500 | U |
| 75-27-4 | Bromodichloromethane | | 2500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 2500 | U |
| 108-88-3 | Toluene | | 2500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 2500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 2500 | U |
| 127-18-4 | Tetrachloroethene | | 32000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 2500 | U |
| 124-48-1 | Dibromochloromethane | | 2500 | U |
| 108-90-7 | Chlorobenzene | | 2500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 2500 | U |
| 100-41-4 | Ethylbenzene | | 2500 | U |
| 1330-20-7 | Xylene (para & meta) | | 2500 | U |
| 95-47-6 | Xylene (Ortho) | | 2500 | U |
| 100-42-5 | Styrene | | 2500 | U |
| 75-25-2 | Bromoform | | 2500 | U |
| 98-82-8 | Isopropylbenzene | | 2500 | U |
| 108-86-1 | Bromobenzene | | 2500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-8

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_8 0.020 ml
 Sample wt/vol: 0.0 (g/ml) ML Lab File ID: 7519_8.D
 Level: (low/med) LOW Date Received: 01/11/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 2500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 2500 | U |
| 103-65-1 | n-Propylbenzene | | 2500 | U |
| 95-49-8 | 2-Chlorotoluene | | 2500 | U |
| 106-43-4 | 4-Chlorotoluene | | 2500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 2500 | U |
| 98-06-6 | tert-Butylbenzene | | 2500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 2500 | U |
| 135-98-8 | sec-Butylbenzene | | 2500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 2500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 2500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 2500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 2500 | U |
| 104-51-8 | n-Butylbenzene | | 2500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 2500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 2500 | U |
| 87-68-3 | Hexachlorobutadiene | | 2500 | U |
| 91-20-3 | Naphthalene | | 2500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 2500 | U |
| 1634-04-4 | MTBE | | 2500 | U |
| 67-64-1 | Acetone | | 5000 | U |
| 75-15-0 | Carbon disulfide | | 2500 | U |
| 78-93-3 | 2-Butanone | | 5000 | U |
| 109-99-9 | Tetrahydrofuran | | 5000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 5000 | U |
| 591-78-6 | 2-Hexanone | | 5000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 5000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-09

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_09 0.020 ml
 Sample wt/vol: 0.0 (g/ml) ML Lab File ID: 7519_09.D
 Level: (low/med) LOW Date Received: 01/11/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 2500 | U |
| 74-87-3 | Chloromethane | | 2500 | U |
| 75-01-4 | Vinyl chloride | | 2500 | U |
| 74-83-9 | Bromomethane | | 2500 | U |
| 75-00-3 | Chloroethane | | 2500 | U |
| 75-69-4 | Trichlorofluoromethane | | 2500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 2500 | U |
| 75-09-2 | Methylene chloride | | 2500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 2500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 2500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 2500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 2500 | U |
| 74-97-5 | Bromochloromethane | | 2500 | U |
| 67-66-3 | Chloroform | | 2500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 2500 | U |
| 56-23-5 | Carbon tetrachloride | | 2500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 2500 | U |
| 71-43-2 | Benzene | | 2500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 2500 | U |
| 79-01-6 | Trichloroethene | | 2500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 2500 | U |
| 74-95-3 | Dibromomethane | | 2500 | U |
| 75-27-4 | Bromodichloromethane | | 2500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 2500 | U |
| 108-88-3 | Toluene | | 2500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 2500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 2500 | U |
| 127-18-4 | Tetrachloroethene | | 22000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 2500 | U |
| 124-48-1 | Dibromochloromethane | | 2500 | U |
| 108-90-7 | Chlorobenzene | | 2500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 2500 | U |
| 100-41-4 | Ethylbenzene | | 2500 | U |
| 1330-20-7 | Xylene (para & meta) | | 2500 | U |
| 95-47-6 | Xylene (Ortho) | | 2500 | U |
| 100-42-5 | Styrene | | 2500 | U |
| 75-25-2 | Bromoform | | 2500 | U |
| 98-82-8 | Isopropylbenzene | | 2500 | U |
| 108-86-1 | Bromobenzene | | 2500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-09

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_09 0.020 ml
 Sample wt/vol: 0.0 (g/ml) ML Lab File ID: 7519_09.D
 Level: (low/med) LOW Date Received: 01/11/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 2500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 2500 | U |
| 103-65-1 | n-Propylbenzene | | 2500 | U |
| 95-49-8 | 2-Chlorotoluene | | 2500 | U |
| 106-43-4 | 4-Chlorotoluene | | 2500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 2500 | U |
| 98-06-6 | tert-Butylbenzene | | 2500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 2500 | U |
| 135-98-8 | sec-Butylbenzene | | 2500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 2500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 2500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 2500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 2500 | U |
| 104-51-8 | n-Butylbenzene | | 2500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 2500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 2500 | U |
| 87-68-3 | Hexachlorobutadiene | | 2500 | U |
| 91-20-3 | Naphthalene | | 2500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 2500 | U |
| 1634-04-4 | MTBE | | 2500 | U |
| 67-64-1 | Acetone | | 5000 | U |
| 75-15-0 | Carbon disulfide | | 2500 | U |
| 78-93-3 | 2-Butanone | | 5000 | U |
| 109-99-9 | Tetrahydrofuran | | 5000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 5000 | U |
| 591-78-6 | 2-Hexanone | | 5000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 5000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-10

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_10 0.020 ml
 Sample wt/vol: 0.0 (g/ml) ML Lab File ID: 7519_10.D
 Level: (low/med) LOW Date Received: 01/11/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|----|
| 75-71-8 | Dichlorodifluoromethane | | 2500 | U |
| 74-87-3 | Chloromethane | | 2500 | U |
| 75-01-4 | Vinyl chloride | | 2500 | U |
| 74-83-9 | Bromomethane | | 2500 | U |
| 75-00-3 | Chloroethane | | 2500 | U |
| 75-69-4 | Trichlorofluoromethane | | 2500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 2500 | U |
| 75-09-2 | Methylene chloride | | 2500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 2500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 2500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 2500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 2500 | U |
| 74-97-5 | Bromochloromethane | | 2500 | U |
| 67-66-3 | Chloroform | | 2500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 2500 | U |
| 56-23-5 | Carbon tetrachloride | | 2500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 2500 | U |
| 71-43-2 | Benzene | | 2500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 2500 | U |
| 79-01-6 | Trichloroethene | | 2500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 2500 | U |
| 74-95-3 | Dibromomethane | | 2500 | U |
| 75-27-4 | Bromodichloromethane | | 2500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 2500 | U |
| 108-88-3 | Toluene | | 2500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 2500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 2500 | U |
| 127-18-4 | Tetrachloroethene | | 2100 | JD |
| 142-28-9 | 1,3-Dichloropropane | | 2500 | U |
| 124-48-1 | Dibromochloromethane | | 2500 | U |
| 108-90-7 | Chlorobenzene | | 2500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 2500 | U |
| 100-41-4 | Ethylbenzene | | 2500 | U |
| 1330-20-7 | Xylene (para & meta) | | 2500 | U |
| 95-47-6 | Xylene (Ortho) | | 2500 | U |
| 100-42-5 | Styrene | | 2500 | U |
| 75-25-2 | Bromoform | | 2500 | U |
| 98-82-8 | Isopropylbenzene | | 2500 | U |
| 108-86-1 | Bromobenzene | | 2500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-10

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_10 0.020 ml
 Sample wt/vol: 0.0 (g/ml) ML Lab File ID: 7519_10.D
 Level: (low/med) LOW Date Received: 01/11/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 2500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 2500 | U |
| 103-65-1 | n-Propylbenzene | | 2500 | U |
| 95-49-8 | 2-Chlorotoluene | | 2500 | U |
| 106-43-4 | 4-Chlorotoluene | | 2500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 2500 | U |
| 98-06-6 | tert-Butylbenzene | | 2500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 2500 | U |
| 135-98-8 | sec-Butylbenzene | | 2500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 2500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 2500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 2500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 2500 | U |
| 104-51-8 | n-Butylbenzene | | 2500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 2500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 2500 | U |
| 87-68-3 | Hexachlorobutadiene | | 2500 | U |
| 91-20-3 | Naphthalene | | 2500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 2500 | U |
| 1634-04-4 | MTBE | | 2500 | U |
| 67-64-1 | Acetone | | 5000 | U |
| 75-15-0 | Carbon disulfide | | 2500 | U |
| 78-93-3 | 2-Butanone | | 5000 | U |
| 109-99-9 | Tetrahydrofuran | | 5000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 5000 | U |
| 591-78-6 | 2-Hexanone | | 5000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 5000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-11

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_11 0.020 ml
 Sample wt/vol: 0.0 (g/ml) ML Lab File ID: 7519_11.D
 Level: (low/med) LOW Date Received: 01/11/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 2500 | U |
| 74-87-3 | Chloromethane | | 2500 | U |
| 75-01-4 | Vinyl chloride | | 2500 | U |
| 74-83-9 | Bromomethane | | 2500 | U |
| 75-00-3 | Chloroethane | | 2500 | U |
| 75-69-4 | Trichlorofluoromethane | | 2500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 2500 | U |
| 75-09-2 | Methylene chloride | | 2500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 2500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 2500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 2500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 2500 | U |
| 74-97-5 | Bromochloromethane | | 2500 | U |
| 67-66-3 | Chloroform | | 2500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 2500 | U |
| 56-23-5 | Carbon tetrachloride | | 2500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 2500 | U |
| 71-43-2 | Benzene | | 2500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 2500 | U |
| 79-01-6 | Trichloroethene | | 2500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 2500 | U |
| 74-95-3 | Dibromomethane | | 2500 | U |
| 75-27-4 | Bromodichloromethane | | 2500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 2500 | U |
| 108-88-3 | Toluene | | 2500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 2500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 2500 | U |
| 127-18-4 | Tetrachloroethene | | 17000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 2500 | U |
| 124-48-1 | Dibromochloromethane | | 2500 | U |
| 108-90-7 | Chlorobenzene | | 2500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 2500 | U |
| 100-41-4 | Ethylbenzene | | 2500 | U |
| 1330-20-7 | Xylene (para & meta) | | 2500 | U |
| 95-47-6 | Xylene (Ortho) | | 2500 | U |
| 100-42-5 | Styrene | | 2500 | U |
| 75-25-2 | Bromoform | | 2500 | U |
| 98-82-8 | Isopropylbenzene | | 2500 | U |
| 108-86-1 | Bromobenzene | | 2500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-11

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_11 0.020 ml
 Sample wt/vol: 0.0 (g/ml) ML Lab File ID: 7519_11.D
 Level: (low/med) LOW Date Received: 01/11/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 2500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 2500 | U |
| 103-65-1 | n-Propylbenzene | | 2500 | U |
| 95-49-8 | 2-Chlorotoluene | | 2500 | U |
| 106-43-4 | 4-Chlorotoluene | | 2500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 2500 | U |
| 98-06-6 | tert-Butylbenzene | | 2500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 2500 | U |
| 135-98-8 | sec-Butylbenzene | | 2500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 2500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 2500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 2500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 2500 | U |
| 104-51-8 | n-Butylbenzene | | 2500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 2500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 2500 | U |
| 87-68-3 | Hexachlorobutadiene | | 2500 | U |
| 91-20-3 | Naphthalene | | 2500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 2500 | U |
| 1634-04-4 | MTBE | | 2500 | U |
| 67-64-1 | Acetone | | 5000 | U |
| 75-15-0 | Carbon disulfide | | 2500 | U |
| 78-93-3 | 2-Butanone | | 5000 | U |
| 109-99-9 | Tetrahydrofuran | | 5000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 5000 | U |
| 591-78-6 | 2-Hexanone | | 5000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 5000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7119-12

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_12 0.020 ml
 Sample wt/vol: 0.0 (g/ml) ML Lab File ID: 7519_12.D
 Level: (low/med) LOW Date Received: 01/11/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 2500 | U |
| 74-87-3 | Chloromethane | | 2500 | U |
| 75-01-4 | Vinyl chloride | | 2500 | U |
| 74-83-9 | Bromomethane | | 2500 | U |
| 75-00-3 | Chloroethane | | 2500 | U |
| 75-69-4 | Trichlorofluoromethane | | 2500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 2500 | U |
| 75-09-2 | Methylene chloride | | 2500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 2500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 2500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 2500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 2500 | U |
| 74-97-5 | Bromochloromethane | | 2500 | U |
| 67-66-3 | Chloroform | | 2500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 2500 | U |
| 56-23-5 | Carbon tetrachloride | | 2500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 2500 | U |
| 71-43-2 | Benzene | | 2500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 2500 | U |
| 79-01-6 | Trichloroethene | | 2500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 2500 | U |
| 74-95-3 | Dibromomethane | | 2500 | U |
| 75-27-4 | Bromodichloromethane | | 2500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 2500 | U |
| 108-88-3 | Toluene | | 2500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 2500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 2500 | U |
| 127-18-4 | Tetrachloroethene | | 13000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 2500 | U |
| 124-48-1 | Dibromochloromethane | | 2500 | U |
| 108-90-7 | Chlorobenzene | | 2500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 2500 | U |
| 100-41-4 | Ethylbenzene | | 2500 | U |
| 1330-20-7 | Xylene (para & meta) | | 2500 | U |
| 95-47-6 | Xylene (Ortho) | | 2500 | U |
| 100-42-5 | Styrene | | 2500 | U |
| 75-25-2 | Bromoform | | 2500 | U |
| 98-82-8 | Isopropylbenzene | | 2500 | U |
| 108-86-1 | Bromobenzene | | 2500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7119-12

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_12 0.020 ml
 Sample wt/vol: 0.0 (g/ml) ML Lab File ID: 7519_12.D
 Level: (low/med) LOW Date Received: 01/11/05
 % Moisture: not dec. _____ Date Analyzed: 01/23/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 500.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 2500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 2500 | U |
| 103-65-1 | n-Propylbenzene | | 2500 | U |
| 95-49-8 | 2-Chlorotoluene | | 2500 | U |
| 106-43-4 | 4-Chlorotoluene | | 2500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 2500 | U |
| 98-06-6 | tert-Butylbenzene | | 2500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 2500 | U |
| 135-98-8 | sec-Butylbenzene | | 2500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 2500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 2500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 2500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 2500 | U |
| 104-51-8 | n-Butylbenzene | | 2500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 2500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 2500 | U |
| 87-68-3 | Hexachlorobutadiene | | 2500 | U |
| 91-20-3 | Naphthalene | | 2500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 2500 | U |
| 1634-04-4 | MTBE | | 2500 | U |
| 67-64-1 | Acetone | | 5000 | U |
| 75-15-0 | Carbon disulfide | | 2500 | U |
| 78-93-3 | 2-Butanone | | 5000 | U |
| 109-99-9 | Tetrahydrofuran | | 5000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 5000 | U |
| 591-78-6 | 2-Hexanone | | 5000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 5000 | U |

3.0 Raw Data and Chromatograms

Raw data and Chromatograms are attached.

Acq On : 23 Jan 2007 3:57 pm

Operator: RR/AS

Sample : 7519_10 0.020 ml

Inst : GC/MS Ins

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 25 11:25 2007

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:51:14 2007

Response via : Initial Calibration

DataAcq Meth : RUN

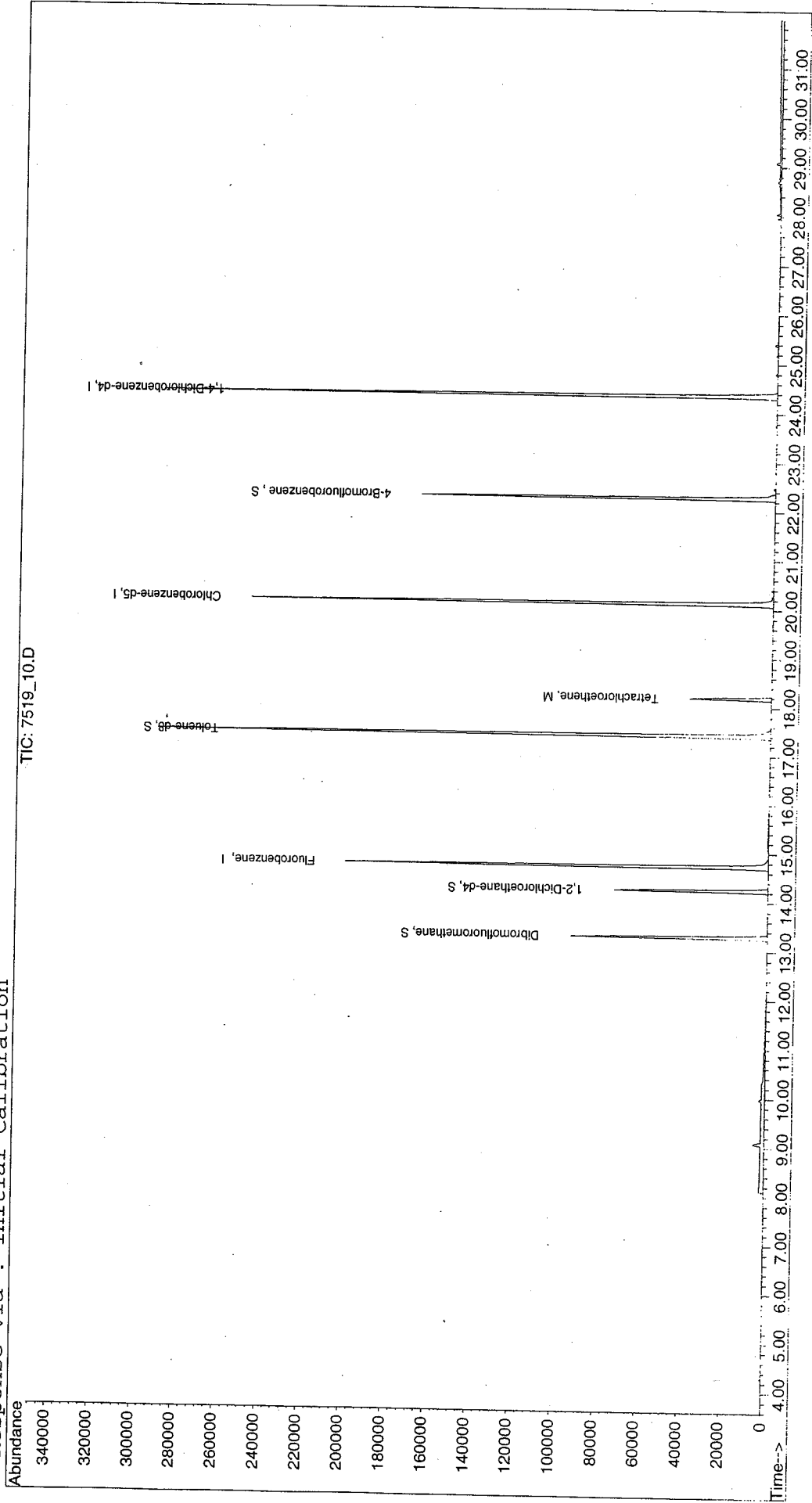
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|------|------|--------|
| 32) Tetrachloroethene | 18.18 | 166 | 21833 | 4.29 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 20.15 | 112 | 6795 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7519_10.D 01_12_07.M Thu Jan 25 16:58:54 2007

Quantitation Report

Data File : C:\HPCHEM\1\DATA\0102307\7519_10.D
Acq On : 23 Jan 2007 3:57 pm
Sample : 7519_10 0.020 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 25 11:25 2007
Vial: 11
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: 01_12_07.RES
Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:51:14 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\0102307\7519_11.D
 Acq On : 23 Jan 2007 4:54 pm
 Sample : 7519_11 0.020 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 25 11:26 2007

Vial: 1
 Operator: RR/AS
 Inst : GC/MS Ins
 Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Mon Jan 15 14:51:14 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.73 | 96 | 249514 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 20.12 | 117 | 215188 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.35 | 152 | 117041 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|------|
| 16) Dibromofluoromethane | 13.29 | 113 | 74495 | 27.51 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 110.04% | |
| 19) 1,2-Dichloroethane-d4 | 14.24 | 65 | 70987 | 25.87 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 103.48% | |
| 28) Toluene-d8 | 17.41 | 98 | 279061 | 24.76 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 99.04% | |
| 45) 4-Bromofluorobenzene | 22.27 | 95 | 75750 | 20.61 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 82.44% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration
 7519_11.D 01_12_07.M Thu Jan 25 16:59:01 2007

Data File : C:\HPCHEM\1\DATA\0102307\7519_11.D
Acq On : 23 Jan 2007 4:54 pm
Sample : 7519_11 0.020 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 25 11:26 2007

Vial: 1
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:51:14 2007
Response via : Initial Calibration
DataAcq Meth : RUN

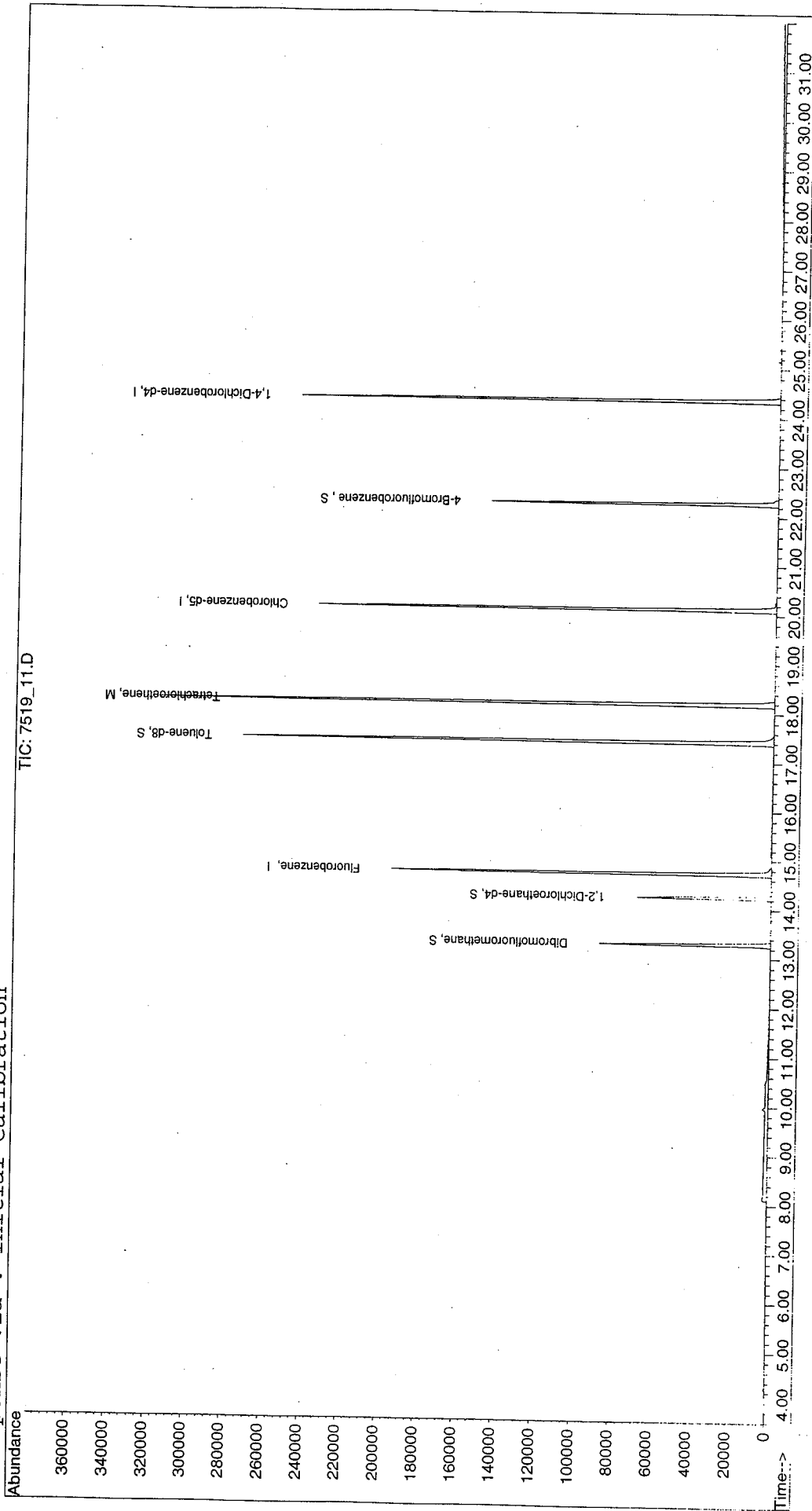
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 164634 | 33.77 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration
7519_11.D 01_12_07.M Thu Jan 25 16:59:03 2007

Quantitation Report

Data File : C:\HPCHEM\1\DATA\0102307\7519_11.D
Acq On : 23 Jan 2007 4:54 pm
Sample : 7519_11 0.020 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 25 11:26 2007
Vial: 1
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: 01_12_07.RES

Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:51:14 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\0102307\7519_12.D
 Acq On : 23 Jan 2007 5:33 pm
 Sample : 7519_12 0.020 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 25 11:26 2007

Vial: 2
 Operator: RR/AS
 Inst : GC/MS Ins
 Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Mon Jan 15 14:51:14 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN

Internal Standards R.T. QIon Response Conc Units Dev(Min)

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.74 | 96 | 261609 | 25.00 | ug/L | 0.01 |
| 36) Chlorobenzene-d5 | 20.12 | 117 | 219160 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.36 | 152 | 119172 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|--------|----------|------|---------|
| 16) Dibromofluoromethane | 13.29 | 113 | 79620 | 28.04 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | Recovery | = | 112.16% |
| 19) 1,2-Dichloroethane-d4 | 14.25 | 65 | 75339 | 26.19 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | Recovery | = | 104.76% |
| 28) Toluene-d8 | 17.41 | 98 | 286796 | 24.27 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | Recovery | = | 97.08% |
| 45) 4-Bromofluorobenzene | 22.28 | 95 | 77077 | 20.60 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | Recovery | = | 82.40% |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\0102307\7519_12.D

Acq On : 23 Jan 2007 5:33 pm

Sample : 7519_12 0.020 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Jan 25 11:26 2007

Vial: 2

Operator: RR/AS

Inst : GC/MS Ins

Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:51:14 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 134983 | 26.41 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 20.14 | 112 | 5184 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7519_12.D 01_12_07.M

Thu Jan 25 16:59:11 2007

Page 2

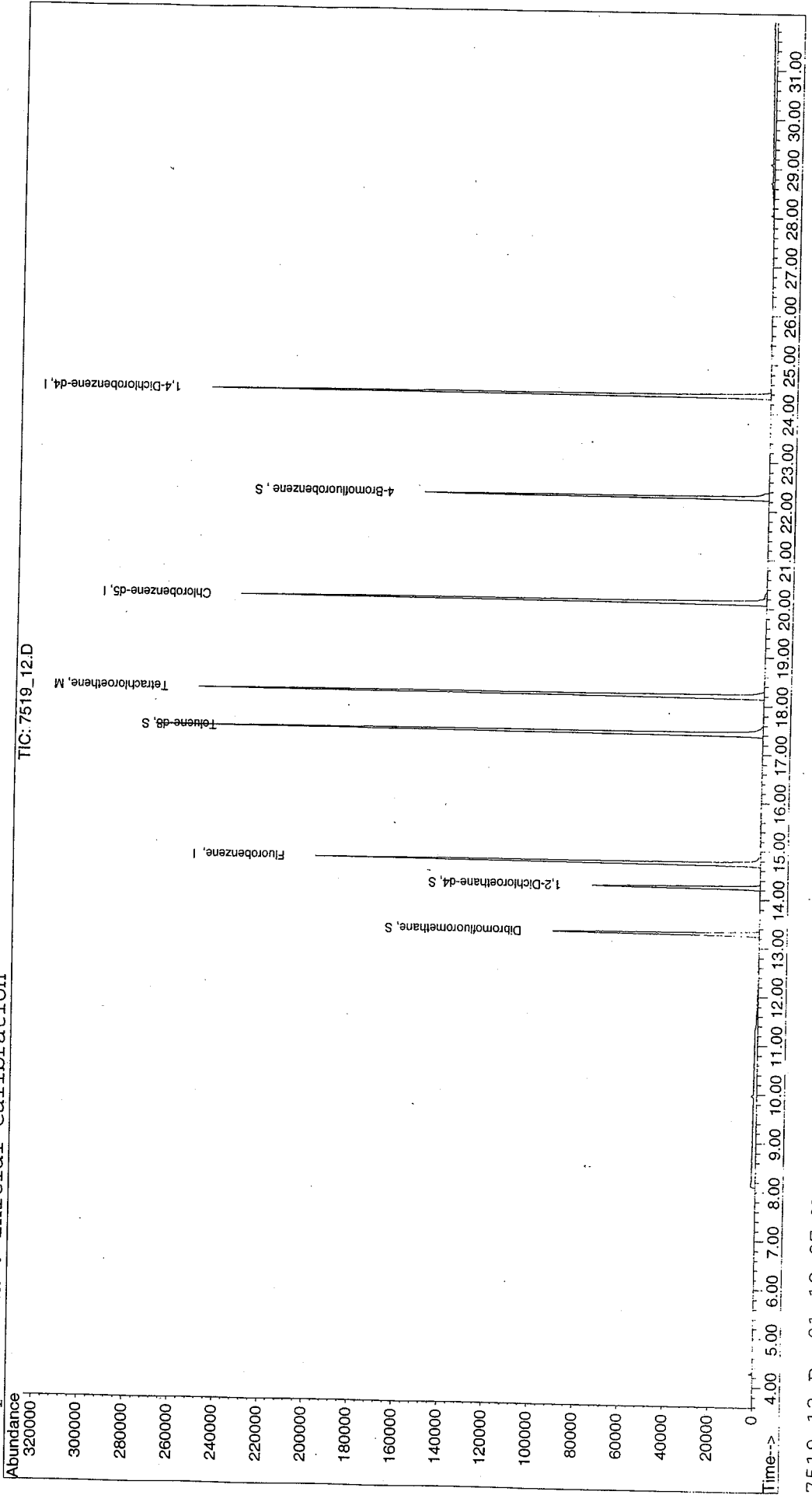
Quantitation Report

Data File : C:\HPCHEM\1\DATA\0102307\7519_12.D
Acq On : 23 Jan 2007 5:33 pm
Sample : 7519_12 0.020 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 25 11:26 2007

Vial: 2
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00

Quant Results File: 01_12_07.RE5

Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:51:14 2007
Response via : Initial Calibration



1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-01

Lab Name: ATL. Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA.
 Matrix: (soil/water) WATER Lab Sample ID: 7519_01 0.10 ml
 Sample wt/vol: 0.1 (g/ml) ML Lab File ID: 7519_01.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 100.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|----|
| 75-71-8 | Dichlorodifluoromethane | | 500 | U |
| 74-87-3 | Chloromethane | | 500 | U |
| 75-01-4 | Vinyl chloride | | 500 | U |
| 74-83-9 | Bromomethane | | 500 | U |
| 75-00-3 | Chloroethane | | 500 | U |
| 75-69-4 | Trichlorofluoromethane | | 500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 500 | U |
| 75-09-2 | Methylene chloride | | 500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 500 | U |
| 74-97-5 | Bromochloromethane | | 500 | U |
| 67-66-3 | Chloroform | | 500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 500 | U |
| 56-23-5 | Carbon tetrachloride | | 500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 500 | U |
| 71-43-2 | Benzene | | 500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 500 | U |
| 79-01-6 | Trichloroethene | | 500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 500 | U |
| 74-95-3 | Dibromomethane | | 500 | U |
| 75-27-4 | Bromodichloromethane | | 500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 500 | U |
| 108-88-3 | Toluene | | 500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 500 | U |
| 127-18-4 | Tetrachloroethene | | 39000 | ED |
| 142-28-9 | 1,3-Dichloropropane | | 500 | U |
| 124-48-1 | Dibromochloromethane | | 500 | U |
| 108-90-7 | Chlorobenzene | | 500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 500 | U |
| 100-41-4 | Ethylbenzene | | 500 | U |
| 1330-20-7 | Xylene (para & meta) | | 500 | U |
| 95-47-6 | Xylene (Ortho) | | 500 | U |
| 100-42-5 | Styrene | | 500 | U |
| 75-25-2 | Bromoform | | 500 | U |
| 98-82-8 | Isopropylbenzene | | 500 | U |
| 108-86-1 | Bromobenzene | | 500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-01

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_01 0.10 ml
 Sample wt/vol: 0.1 (g/ml) ML Lab File ID: 7519_01.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 100.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 500 | U |
| 103-65-1 | n-Propylbenzene | | 500 | U |
| 95-49-8 | 2-Chlorotoluene | | 500 | U |
| 106-43-4 | 4-Chlorotoluene | | 500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 500 | U |
| 98-06-6 | tert-Butylbenzene | | 500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 500 | U |
| 135-98-8 | sec-Butylbenzene | | 500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 500 | U |
| 104-51-8 | n-Butylbenzene | | 500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 500 | U |
| 87-68-3 | Hexachlorobutadiene | | 500 | U |
| 91-20-3 | Naphthalene | | 500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 500 | U |
| 1634-04-4 | MTBE | | 500 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 500 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-02

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_02 0.10 ml
 Sample wt/vol: 0.1 (g/ml) ML Lab File ID: 7519_02.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 100.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|----|
| 75-71-8 | Dichlorodifluoromethane | | 500 | U |
| 74-87-3 | Chloromethane | | 500 | U |
| 75-01-4 | Vinyl chloride | | 500 | U |
| 74-83-9 | Bromomethane | | 500 | U |
| 75-00-3 | Chloroethane | | 500 | U |
| 75-69-4 | Trichlorofluoromethane | | 500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 500 | U |
| 75-09-2 | Methylene chloride | | 500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 500 | U |
| 74-97-5 | Bromochloromethane | | 500 | U |
| 67-66-3 | Chloroform | | 500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 500 | U |
| 56-23-5 | Carbon tetrachloride | | 500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 500 | U |
| 71-43-2 | Benzene | | 500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 500 | U |
| 79-01-6 | Trichloroethene | | 500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 500 | U |
| 74-95-3 | Dibromomethane | | 500 | U |
| 75-27-4 | Bromodichloromethane | | 500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 500 | U |
| 108-88-3 | Toluene | | 500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 500 | U |
| 127-18-4 | Tetrachloroethene | | 42000 | ED |
| 142-28-9 | 1,3-Dichloropropane | | 500 | U |
| 124-48-1 | Dibromochloromethane | | 500 | U |
| 108-90-7 | Chlorobenzene | | 500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 500 | U |
| 100-41-4 | Ethylbenzene | | 500 | U |
| 1330-20-7 | Xylene (para & meta) | | 500 | U |
| 95-47-6 | Xylene (Ortho) | | 500 | U |
| 100-42-5 | Styrene | | 500 | U |
| 75-25-2 | Bromoform | | 500 | U |
| 98-82-8 | Isopropylbenzene | | 500 | U |
| 108-86-1 | Bromobenzene | | 500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-02

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_02 0.10 ml
 Sample wt/vol: 0.1 (g/ml) ML Lab File ID: 7519_02.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 100.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 500 | U |
| 103-65-1 | n-Propylbenzene | | 500 | U |
| 95-49-8 | 2-Chlorotoluene | | 500 | U |
| 106-43-4 | 4-Chlorotoluene | | 500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 500 | U |
| 98-06-6 | tert-Butylbenzene | | 500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 500 | U |
| 135-98-8 | sec-Butylbenzene | | 500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 500 | U |
| 104-51-8 | n-Butylbenzene | | 500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 500 | U |
| 87-68-3 | Hexachlorobutadiene | | 500 | U |
| 91-20-3 | Naphthalene | | 500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 500 | U |
| 1634-04-4 | MTBE | | 500 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 500 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-03

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_03 0.10 ml
 Sample wt/vol: 0.1 (g/ml) ML Lab File ID: 7519_03.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 100.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|----|
| 75-71-8 | Dichlorodifluoromethane | | 500 | U |
| 74-87-3 | Chloromethane | | 500 | U |
| 75-01-4 | Vinyl chloride | | 500 | U |
| 74-83-9 | Bromomethane | | 500 | U |
| 75-00-3 | Chloroethane | | 500 | U |
| 75-69-4 | Trichlorofluoromethane | | 500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 500 | U |
| 75-09-2 | Methylene chloride | | 500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 500 | U |
| 74-97-5 | Bromochloromethane | | 500 | U |
| 67-66-3 | Chloroform | | 500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 500 | U |
| 56-23-5 | Carbon tetrachloride | | 500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 500 | U |
| 71-43-2 | Benzene | | 500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 500 | U |
| 79-01-6 | Trichloroethene | | 150 | JD |
| 78-87-5 | 1,2-Dichloropropane | | 500 | U |
| 74-95-3 | Dibromomethane | | 500 | U |
| 75-27-4 | Bromodichloromethane | | 500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 500 | U |
| 108-88-3 | Toluene | | 500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 500 | U |
| 127-18-4 | Tetrachloroethene | | 40000 | ED |
| 142-28-9 | 1,3-Dichloropropane | | 500 | U |
| 124-48-1 | Dibromochloromethane | | 500 | U |
| 108-90-7 | Chlorobenzene | | 500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 500 | U |
| 100-41-4 | Ethylbenzene | | 500 | U |
| 1330-20-7 | Xylene (para & meta) | | 500 | U |
| 95-47-6 | Xylene (Ortho) | | 500 | U |
| 100-42-5 | Styrene | | 500 | U |
| 75-25-2 | Bromoform | | 500 | U |
| 98-82-8 | Isopropylbenzene | | 500 | U |
| 108-86-1 | Bromobenzene | | 500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-03

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_03 0.10 ml
 Sample wt/vol: 0.1 (g/ml) ML Lab File ID: 7519_03.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 100.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 500 | U |
| 103-65-1 | n-Propylbenzene | | 500 | U |
| 95-49-8 | 2-Chlorotoluene | | 500 | U |
| 106-43-4 | 4-Chlorotoluene | | 500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 500 | U |
| 98-06-6 | tert-Butylbenzene | | 500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 500 | U |
| 135-98-8 | sec-Butylbenzene | | 500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 500 | U |
| 104-51-8 | n-Butylbenzene | | 500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 500 | U |
| 87-68-3 | Hexachlorobutadiene | | 500 | U |
| 91-20-3 | Naphthalene | | 500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 500 | U |
| 1634-04-4 | MTBE | | 500 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 500 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

Data File : C:\HPCHEM\1\DATA\0102207\7519_01.D

Acq On : 22 Jan 2007 12:27 pm

Sample : 7519_01 0.10 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Jan 23 14:57 2007

Vial: 5

Operator: RR/AS

Inst : GC/MS Ins

Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:54:07 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.73 | 96 | 293670 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 0.00 | 117 | 0 | 0.00 | ug/L | -20.12 |
| 58) 1,4-Dichlorobenzene-d4 | 0.00 | 152 | 0 | 0.00 | ug/L | -24.35 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|------|
| 16) Dibromofluoromethane | 13.28 | 113 | 85050 | 26.69 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 106.76% | |
| 19) 1,2-Dichloroethane-d4 | 14.25 | 65 | 80179 | 24.83 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 99.32% | |
| 28) Toluene-d8 | 17.42 | 98 | 341762 | 25.77 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 103.08% | |
| 45) 4-Bromofluorobenzene | 0.00 | 95 | 0 | 0.00 | ug/L | |
| Spiked Amount | 25.000 | | Recovery | = | 0.00% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | d | |

(#) = qualifier out of range (m) = manual integration

7519_01.D 01_12_07.M

Thu Jan 25 14:18:06 2007

Data File : C:\HPCHEM\1\DATA\0102207\7519_01.D

Vial: 5

Acq On : 22 Jan 2007 12:27 pm

Operator: RR/AS

Sample : 7519_01 0.10 ml

Inst : GC/MS Ins

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 23 14:57 2007

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:54:07 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|--------|------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 2244882 | 391.29 | ug/L | 100 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 9.06 | 43 | 70745 | N.D. | | |
| 69) Carbon disulfide | 9.06 | 76 | 248583 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

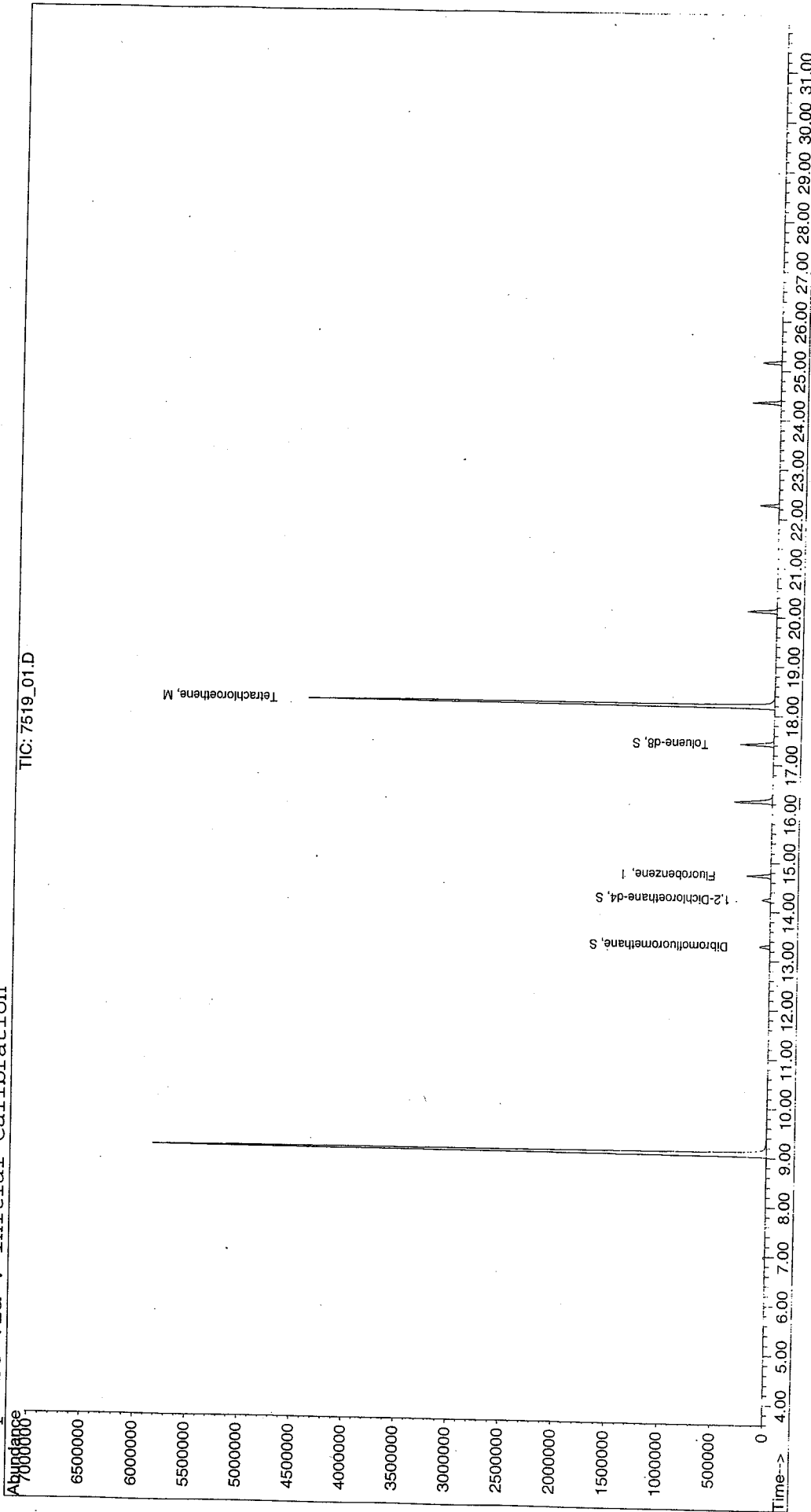
(#) = qualifier out of range (m) = manual integration

7519_01.D 01_12_07.M Thu Jan 25 14:18:09 2007

Quantitation Report

Data File : C:\HPCHEM\1\DATA\0102207\7519_01.D
Acq On : 22 Jan 2007 12:27 pm
Sample : 7519_01 0.10 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 23 14:57 2007
Vial: 5
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: 01_12_07.RES

Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:54:07 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\0102207\7519_02.D
 Acq On : 22 Jan 2007 1:06 pm
 Sample : 7519_02 0.10 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 23 14:59 2007

Vial: 6
 Operator: RR/AS
 Inst : GC/MS Ins
 Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Mon Jan 15 14:54:07 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.73 | 96 | 295750 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 20.12 | 117 | 261159 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.36 | 152 | 130228 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|--------|-------|------|--------------------|
| 16) Dibromofluoromethane | 13.29 | 113 | 85491 | 26.63 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 106.52% |
| 19) 1,2-Dichloroethane-d4 | 14.25 | 65 | 81692 | 25.12 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 100.48% |
| 28) Toluene-d8 | 17.42 | 98 | 348016 | 26.06 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 104.24% |
| 45) 4-Bromofluorobenzene | 22.27 | 95 | 92442 | 20.73 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 82.92% |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | d | |

(#) = qualifier out of range (m) = manual integration
 7519_02.D 01_12_07.M Thu Jan 25 14:18:17 2007

Data File : C:\HPCHEM\1\DATA\0102207\7519_02.D
 Acq On : 22 Jan 2007 1:06 pm
 Sample : 7519_02 0.10 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 23 14:59 2007

Vial: 6
 Operator: RR/AS
 Inst : GC/MS Ins
 Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Mon Jan 15 14:54:07 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN

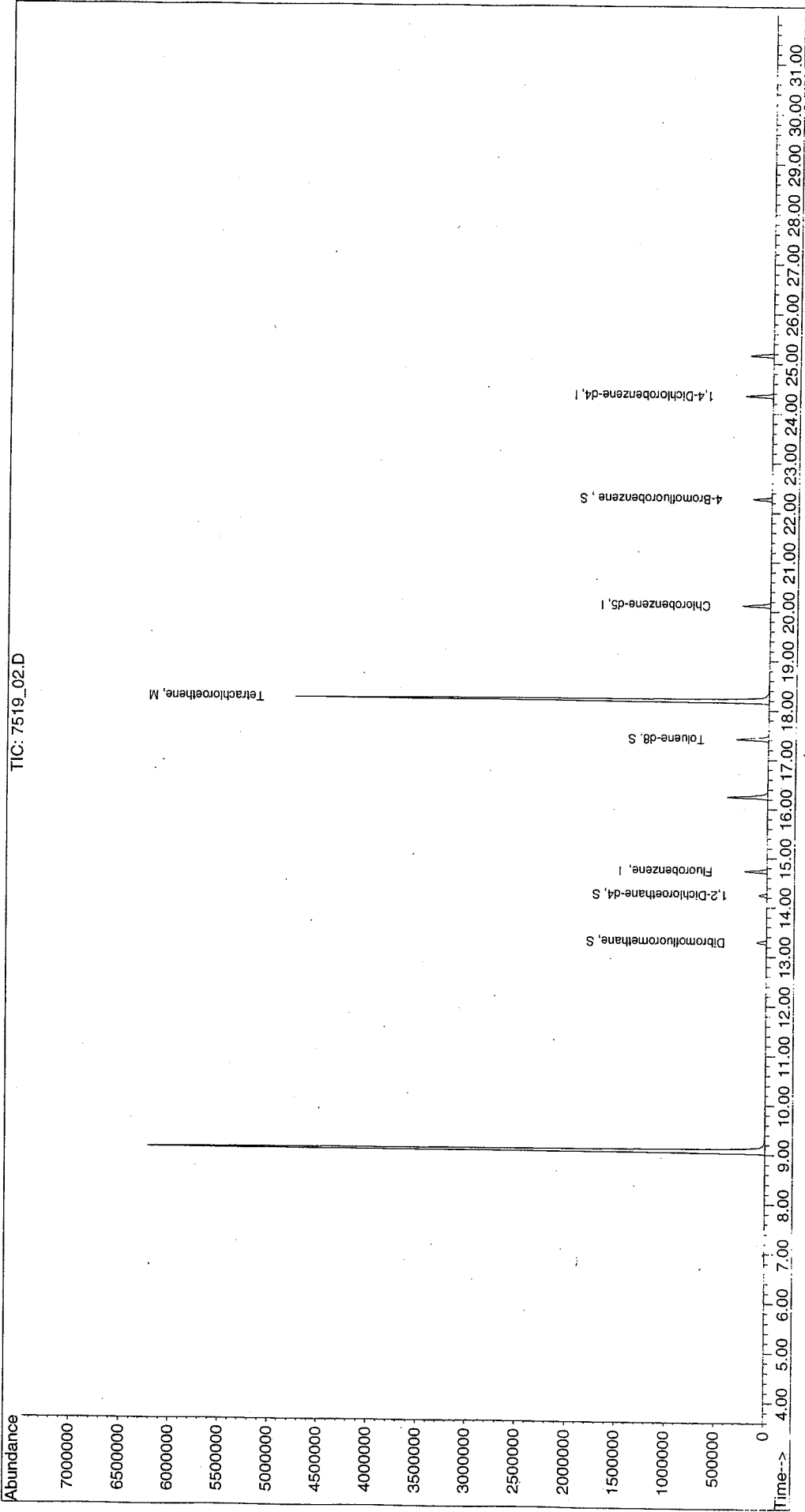
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|--------|------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 2439524 | 422.23 | ug/L | 99 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | d | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | d | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | d | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration
 7519_02.D 01_12_07.M Thu Jan 25 14:18:19 2007

Quantitation Report

Data File : C:\HPCHEM\1\DATA\0102207\7519_02.D
Acq On : 22 Jan 2007 1:06 pm
Sample : 7519_02 0.10 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 23 14:59 2007
Vial: 6
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: 01_12_07.REB

Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:54:07 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\0102207\7519_03.D
 Acq On : 22 Jan 2007 1:44 pm
 Sample : 7519_03 0.10 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 24 11:11 2007

Vial: 7
 Operator: RR/AS
 Inst : GC/MS Ins
 Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Mon Jan 15 14:54:07 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN

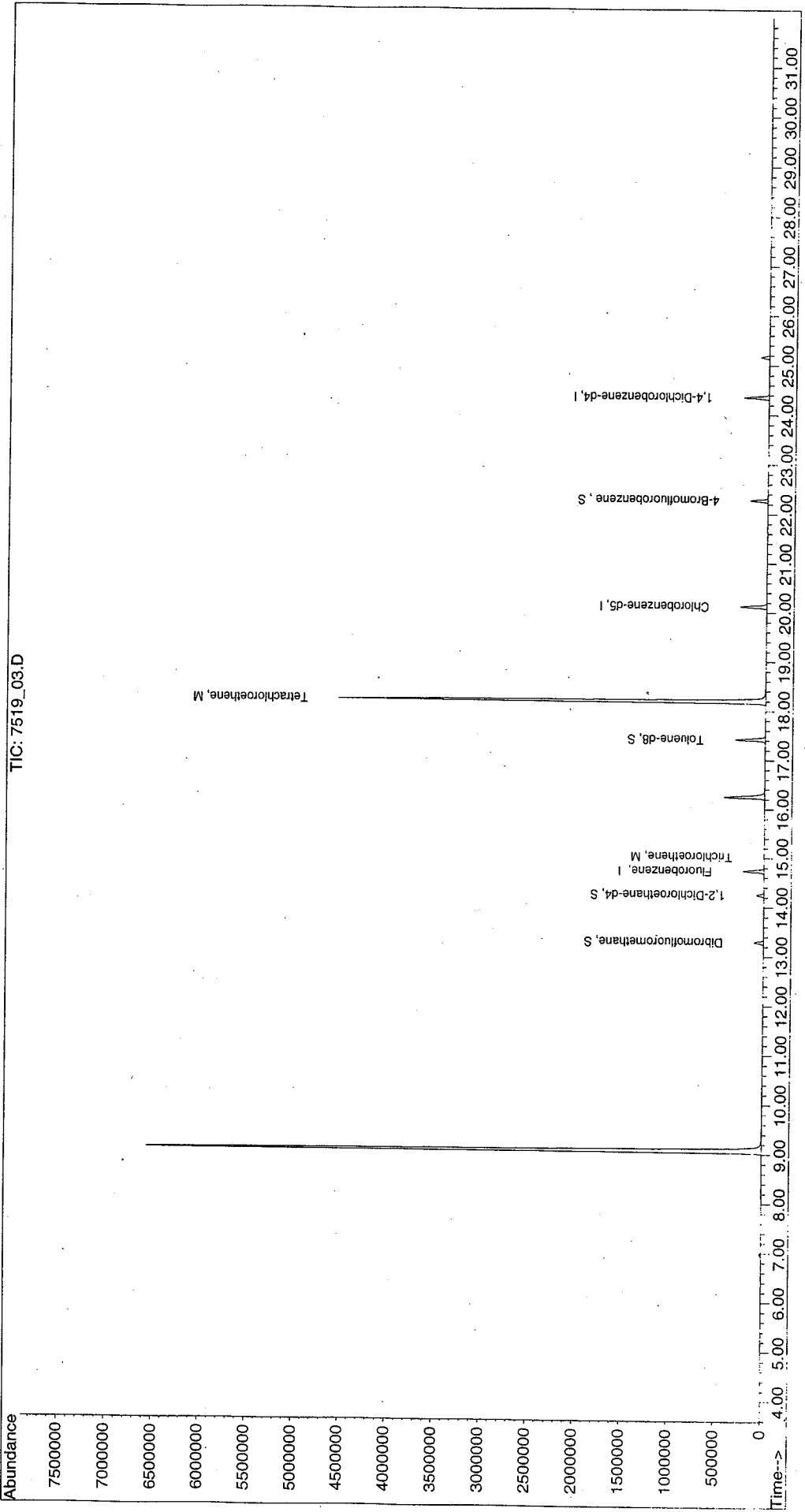
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|--------|------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 2326948 | 402.24 | ug/L | 99 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | d | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | d | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | d | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration
 7519_03.D 01_12_07.M Thu Jan 25 14:18:30 2007

Quantitation Report

Data File : C:\HPCHEM\1\DATA\0102207\7519_03.D
Acq On : 22 Jan 2007 1:44 pm Vial: 7
Sample : 7519_03 0.10 ml Operator: RR/AS
Misc : Inst : GC/MS Ins
MS Integration Params: ODD.P Multiplr: 1.00
Quant Time: Jan 24 11:11 2007 Quant Results File: 01_12_07.RES

Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:54:07 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\0102207\7519_3.D

Vial: 9

Acq On : 22 Jan 2007 10:07 pm

Operator: RR/AS

Sample : 7519-3 0.020ml

Inst : GC/MS Ins

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 24 11:13 2007

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:54:07 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.73 | 96 | 295454 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 20.12 | 117 | 245383 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.35 | 152 | 117674 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|------|
| 16) Dibromofluoromethane | 13.29 | 113 | 77548 | 24.18 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 96.72% | |
| 19) 1,2-Dichloroethane-d4 | 14.24 | 65 | 74269 | 22.86 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 91.44% | |
| 28) Toluene-d8 | 17.42 | 98 | 335295 | 25.13 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 100.52% | |
| 45) 4-Bromofluorobenzene | 22.28 | 95 | 81836 | 19.53 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 78.12% | |

Target Compounds

| | | | | | Qvalue |
|-------------------------------|------|-----|---|------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | d |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | |

(#) = qualifier out of range (m) = manual integration

7519_3.D 01_12_07.M

Thu Jan 25 14:18:59 2007

MEEP Results with non-heated samples
Using Henry's law

| Run Date | Sample | Liquid Volume (mL) | Headspace Volume (mL) | Room temp. (°C) | Room temp. (K) | Obs | | | Obs Measured Ethane (ppmV) | Obs Measured Ethene (ppmV) | Obs Measured Propane (ppmV) | dilution | Calculated Aq. Concentrations | | | Calculated umolar Conc. | | |
|-----------|-------------|--------------------|-----------------------|-----------------|----------------|----------------|---------------|----------------|----------------------------|----------------------------|-----------------------------|----------|-------------------------------|---------------|----------------|-------------------------|-------------|--------------|
| | | | | | | Methane (ppmV) | Ethane (ppmV) | Propane (ppmV) | | | | | Methane (ug/L) | Ethane (ug/L) | Propane (ug/L) | Methane (uM) | Ethane (uM) | Propane (uM) |
| 1/29/2007 | blank | 32 | 10 | 21.67 | 294.7 | 32.374 | 0 | 0 | 0 | 0 | 0 | 1.0 | 8.09 | 0.00 | 0.00 | 0.51 | 0.00 | 0.00 |
| | 100ppmv | 32 | 10 | 21.67 | 294.7 | 112.436 | 97.921 | 97.61 | 95.408 | 0 | 0 | 1.0 | 26.09 | 43.85 | 50.65 | 1.76 | 1.46 | 1.81 |
| | 7519-1 | 3.75 | 1.25 | 21.67 | 294.7 | 74.427 | 4.428 | 0 | 0 | 0 | 0 | 1.0 | 19.82 | 2.10 | 0.00 | 1.23 | 0.07 | 0.00 |
| | 7519-2 | 3.75 | 1.25 | 21.67 | 294.7 | 78.804 | 4.502 | 0 | 0 | 0 | 0 | 1.0 | 20.77 | 2.13 | 0.00 | 1.30 | 0.07 | 0.00 |
| | 7519-3 | 3.75 | 1.25 | 21.67 | 294.7 | 85.485 | 4.131 | 0 | 0 | 0 | 0 | 1.0 | 22.54 | 1.96 | 0.00 | 1.41 | 0.07 | 0.00 |
| | 7519-4 | 3.75 | 1.25 | 21.67 | 294.7 | 82.104 | 4.217 | 0 | 0 | 0 | 0 | 1.0 | 21.64 | 2.00 | 0.00 | 1.35 | 0.07 | 0.00 |
| | 7519-5 | 3.75 | 1.25 | 21.67 | 294.7 | 85.04 | 4.35 | 0 | 0 | 0 | 0 | 1.0 | 25.05 | 2.08 | 0.00 | 1.57 | 0.07 | 0.00 |
| | 7519-6 | 3.75 | 1.25 | 21.67 | 294.7 | 85.248 | 4.511 | 0 | 0 | 0 | 0 | 1.0 | 22.47 | 2.14 | 0.00 | 1.40 | 0.07 | 0.00 |
| | 7519-7 | 3.75 | 1.25 | 21.67 | 294.7 | 83.3 | 4.386 | 0 | 0 | 0 | 0 | 1.0 | 21.96 | 2.08 | 0.00 | 1.37 | 0.07 | 0.00 |
| | 7519-8 | 3.75 | 1.25 | 21.67 | 294.7 | 39.381 | 5.082 | 0 | 0 | 0 | 0 | 1.0 | 10.38 | 2.41 | 0.00 | 0.65 | 0.08 | 0.00 |
| | 7519-9 | 3.75 | 1.25 | 21.67 | 294.7 | 60.906 | 4.447 | 0 | 0 | 0 | 0 | 1.0 | 16.06 | 2.11 | 0.00 | 1.00 | 0.07 | 0.00 |
| | 7519-10 | 3.75 | 1.25 | 21.67 | 294.7 | 60.222 | 9.003 | 0 | 0 | 0 | 0 | 1.0 | 15.88 | 4.26 | 0.00 | 0.99 | 0.14 | 0.00 |
| | blank | 32 | 10 | 21.67 | 294.7 | 9.154 | 0 | 0 | 0 | 0 | 0 | 1.0 | 2.29 | 0.00 | 0.00 | 0.14 | 0.00 | 0.00 |
| | 50ppmv | 32 | 10 | 21.67 | 294.7 | 51.205 | 52.166 | 50.58 | 51.733 | 0 | 0 | 1.0 | 12.79 | 23.36 | 26.24 | 0.80 | 0.78 | 0.94 |
| | 7519-11 | 3.75 | 1.25 | 21.67 | 294.7 | 68.538 | 8.052 | 0 | 0 | 0 | 0 | 1.0 | 18.07 | 3.81 | 0.00 | 1.13 | 0.13 | 0.00 |
| | 7519-12 | 3.75 | 1.25 | 21.67 | 294.7 | 70.623 | 12.135 | 0 | 1.8 | 0 | 0 | 1.0 | 18.62 | 5.75 | 0.00 | 1.16 | 0.19 | 0.00 |
| | 7519-12 MS | 3.75 | 1.25 | 21.67 | 294.7 | 170.238 | 105.873 | 97.222 | 95.157 | 0 | 0 | 1.0 | 44.88 | 50.14 | 52.79 | 2.80 | 1.67 | 1.89 |
| | 7519-12 MSD | 3.75 | 1.25 | 21.67 | 294.7 | 171.865 | 104.782 | 97 | 96.545 | 0 | 0 | 1.0 | 45.33 | 49.63 | 52.67 | 2.83 | 1.65 | 1.88 |
| | blank | 32 | 10 | 21.67 | 294.7 | 12.793 | 0 | 0 | 0 | 0 | 0 | 1.0 | 3.20 | 0.00 | 0.00 | 0.20 | 0.00 | 0.00 |

MEE- henry's constant adjusted to temperature

Propane - henry's constant at 25 C

Acq On : 23 Jan 2007 3:57 pm

Operator: RR/AS

Sample : 7519_10 0.020 ml

Inst : GC/MS Ins

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 25 11:25 2007

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:51:14 2007

Response via : Initial Calibration

DataAcq Meth : RUN

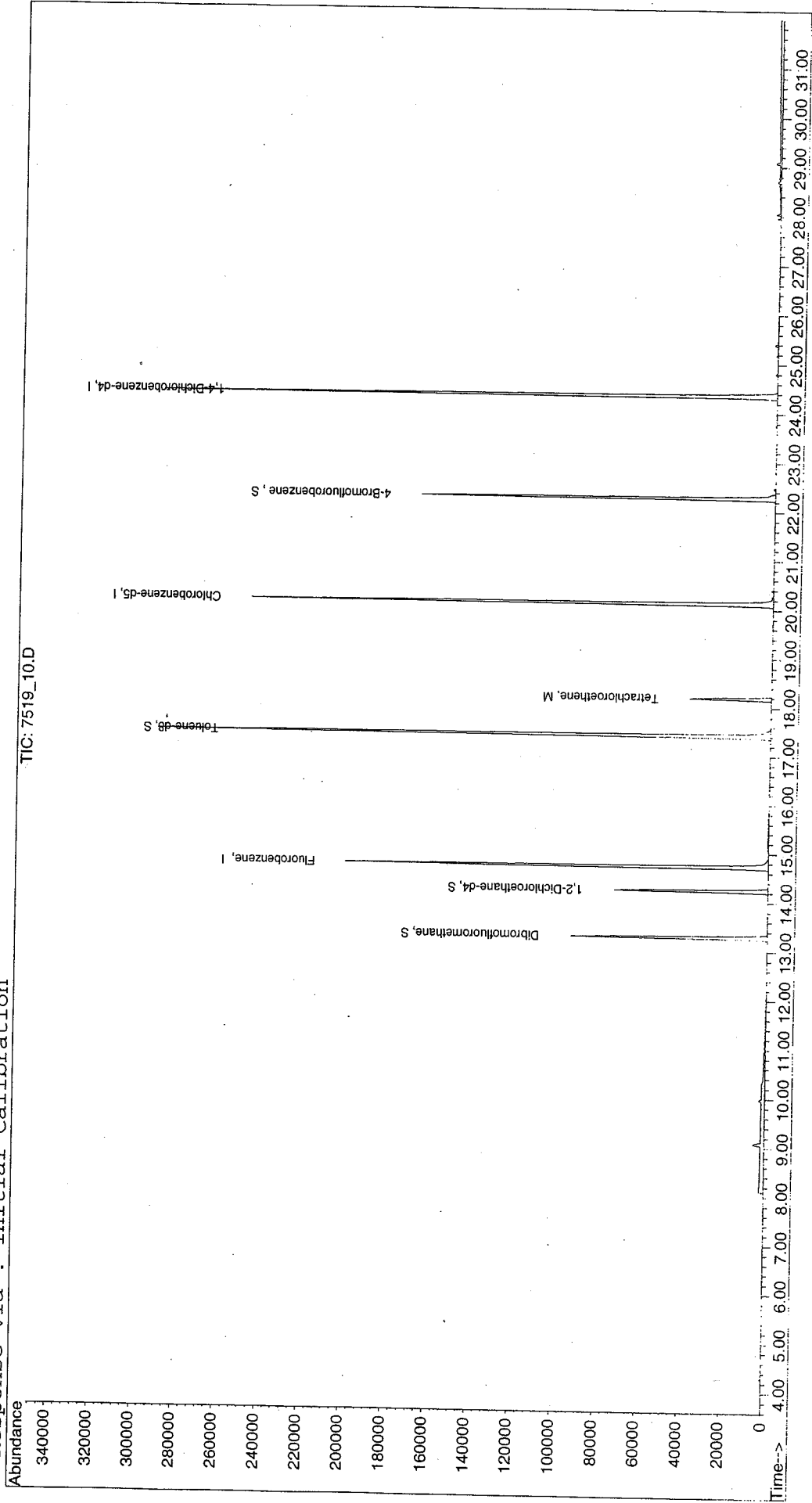
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|------|------|--------|
| 32) Tetrachloroethene | 18.18 | 166 | 21833 | 4.29 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 20.15 | 112 | 6795 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7519_10.D 01_12_07.M Thu Jan 25 16:58:54 2007

Quantitation Report

Data File : C:\HPCHEM\1\DATA\0102307\7519_10.D
Acq On : 23 Jan 2007 3:57 pm
Sample : 7519_10 0.020 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 25 11:25 2007
Vial: 11
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: 01_12_07.RES
Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:51:14 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\0102307\7519_11.D
 Acq On : 23 Jan 2007 4:54 pm
 Sample : 7519_11 0.020 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 25 11:26 2007

Vial: 1
 Operator: RR/AS
 Inst : GC/MS Ins
 Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Mon Jan 15 14:51:14 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.73 | 96 | 249514 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 20.12 | 117 | 215188 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.35 | 152 | 117041 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|------|
| 16) Dibromofluoromethane | 13.29 | 113 | 74495 | 27.51 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 110.04% | |
| 19) 1,2-Dichloroethane-d4 | 14.24 | 65 | 70987 | 25.87 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 103.48% | |
| 28) Toluene-d8 | 17.41 | 98 | 279061 | 24.76 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 99.04% | |
| 45) 4-Bromofluorobenzene | 22.27 | 95 | 75750 | 20.61 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 82.44% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration
 7519_11.D 01_12_07.M Thu Jan 25 16:59:01 2007

Data File : C:\HPCHEM\1\DATA\0102307\7519_11.D
Acq On : 23 Jan 2007 4:54 pm
Sample : 7519_11 0.020 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 25 11:26 2007

Vial: 1
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:51:14 2007
Response via : Initial Calibration
DataAcq Meth : RUN

| Compound | R.T. | QIon | Response | Conc Unit | Qvalue |
|---------------------------------|-------|------|----------|------------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 164634 | 33.77 ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | |

(#) = qualifier out of range (m) = manual integration

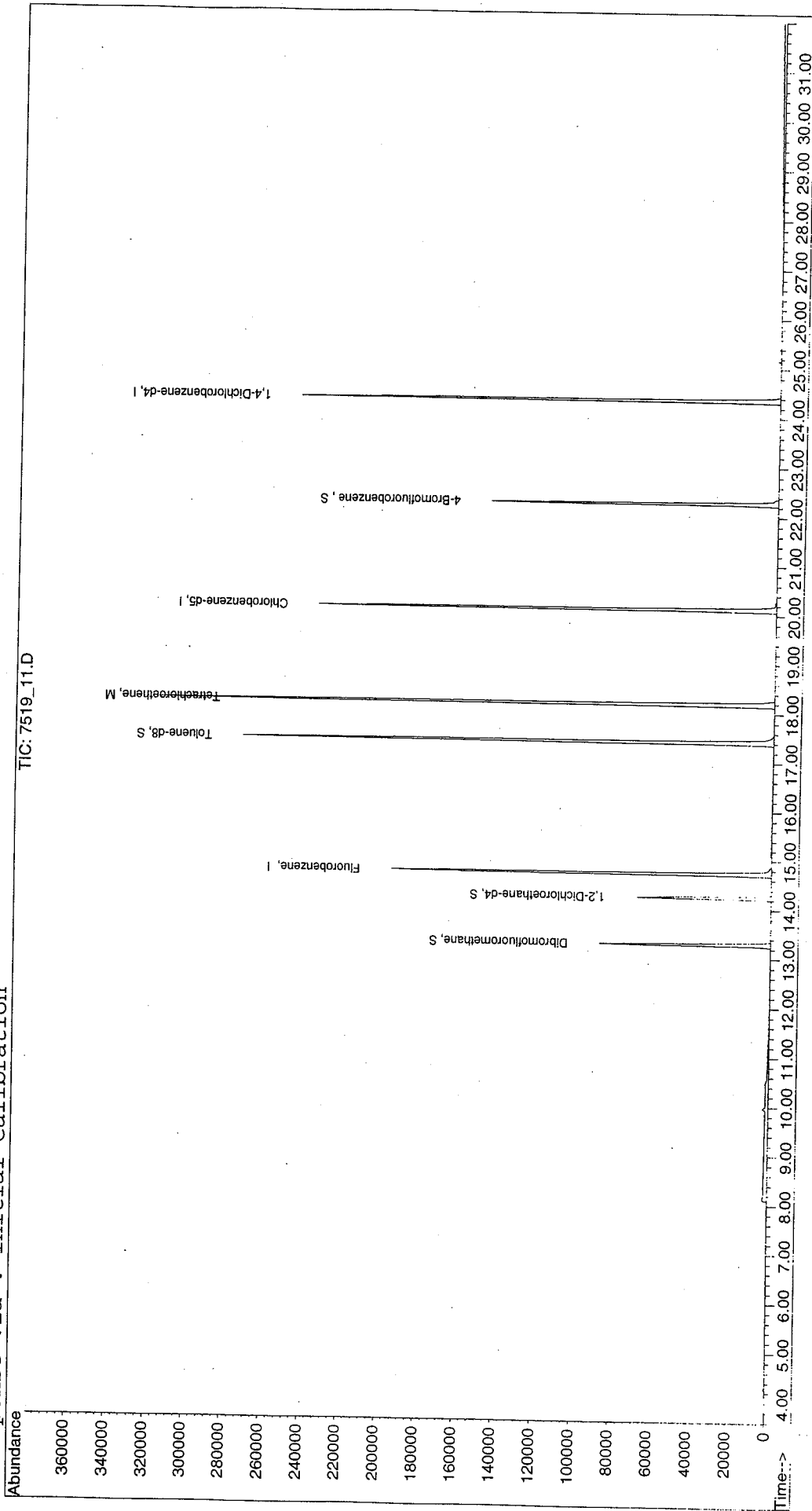
7519_11.D 01_12_07.M

Thu Jan 25 16:59:03 2007

Quantitation Report

Data File : C:\HPCHEM\1\DATA\0102307\7519_11.D
Acq On : 23 Jan 2007 4:54 pm
Sample : 7519_11 0.020 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 25 11:26 2007
Vial: 1
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: 01_12_07.RES

Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:51:14 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\0102307\7519_12.D
 Acq On : 23 Jan 2007 5:33 pm
 Sample : 7519_12 0.020 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 25 11:26 2007

Vial: 2
 Operator: RR/AS
 Inst : GC/MS Ins
 Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Mon Jan 15 14:51:14 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN

Internal Standards R.T. QIon Response Conc Units Dev(Min)

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.74 | 96 | 261609 | 25.00 | ug/L | 0.01 |
| 36) Chlorobenzene-d5 | 20.12 | 117 | 219160 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.36 | 152 | 119172 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|--------|----------|------|---------|
| 16) Dibromofluoromethane | 13.29 | 113 | 79620 | 28.04 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | Recovery | = | 112.16% |
| 19) 1,2-Dichloroethane-d4 | 14.25 | 65 | 75339 | 26.19 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | Recovery | = | 104.76% |
| 28) Toluene-d8 | 17.41 | 98 | 286796 | 24.27 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | Recovery | = | 97.08% |
| 45) 4-Bromofluorobenzene | 22.28 | 95 | 77077 | 20.60 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | Recovery | = | 82.40% |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\0102307\7519_12.D

Acq On : 23 Jan 2007 5:33 pm

Sample : 7519_12 0.020 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Jan 25 11:26 2007

Vial: 2

Operator: RR/AS

Inst : GC/MS Ins

Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:51:14 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 134983 | 26.41 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 20.14 | 112 | 5184 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

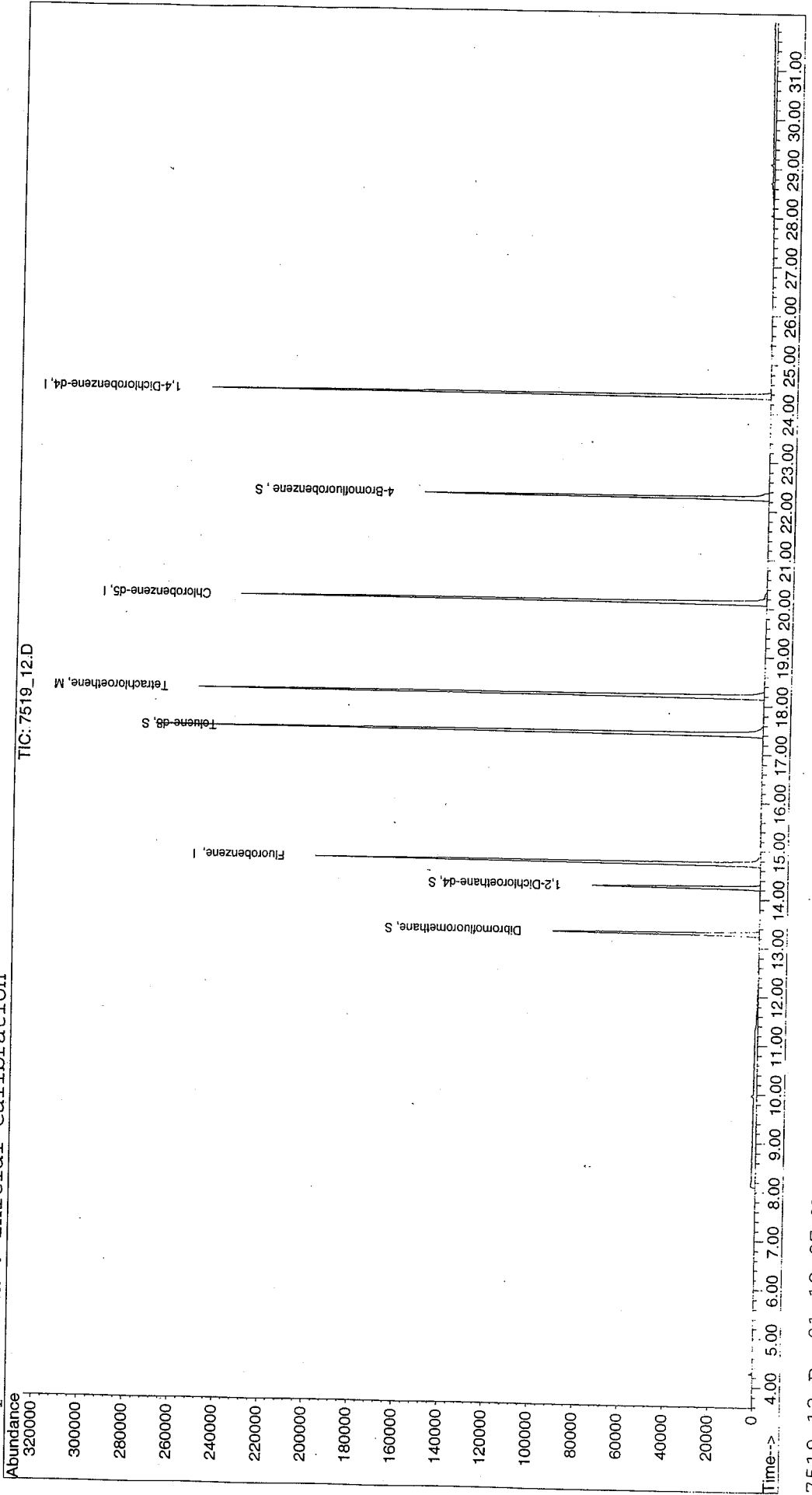
(#) = qualifier out of range (m) = manual integration

7519_12.D 01_12_07.M

Thu Jan 25 16:59:11 2007

Quantitation Report

Data File : C:\HPCHEM\1\DATA\0102307\7519_12.D
Acq On : 23 Jan 2007 5:33 pm
Sample : 7519_12 0.020 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 25 11:26 2007
Vial: 2
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: 01_12_07.RE5
Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:51:14 2007
Response via : Initial Calibration



1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-01

Lab Name: ATL. Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA.
 Matrix: (soil/water) WATER Lab Sample ID: 7519_01 0.10 ml
 Sample wt/vol: 0.1 (g/ml) ML Lab File ID: 7519_01.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 100.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|----|
| 75-71-8 | Dichlorodifluoromethane | | 500 | U |
| 74-87-3 | Chloromethane | | 500 | U |
| 75-01-4 | Vinyl chloride | | 500 | U |
| 74-83-9 | Bromomethane | | 500 | U |
| 75-00-3 | Chloroethane | | 500 | U |
| 75-69-4 | Trichlorofluoromethane | | 500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 500 | U |
| 75-09-2 | Methylene chloride | | 500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 500 | U |
| 74-97-5 | Bromochloromethane | | 500 | U |
| 67-66-3 | Chloroform | | 500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 500 | U |
| 56-23-5 | Carbon tetrachloride | | 500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 500 | U |
| 71-43-2 | Benzene | | 500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 500 | U |
| 79-01-6 | Trichloroethene | | 500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 500 | U |
| 74-95-3 | Dibromomethane | | 500 | U |
| 75-27-4 | Bromodichloromethane | | 500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 500 | U |
| 108-88-3 | Toluene | | 500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 500 | U |
| 127-18-4 | Tetrachloroethene | | 39000 | ED |
| 142-28-9 | 1,3-Dichloropropane | | 500 | U |
| 124-48-1 | Dibromochloromethane | | 500 | U |
| 108-90-7 | Chlorobenzene | | 500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 500 | U |
| 100-41-4 | Ethylbenzene | | 500 | U |
| 1330-20-7 | Xylene (para & meta) | | 500 | U |
| 95-47-6 | Xylene (Ortho) | | 500 | U |
| 100-42-5 | Styrene | | 500 | U |
| 75-25-2 | Bromoform | | 500 | U |
| 98-82-8 | Isopropylbenzene | | 500 | U |
| 108-86-1 | Bromobenzene | | 500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-01

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_01 0.10 ml
 Sample wt/vol: 0.1 (g/ml) ML Lab File ID: 7519_01.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 100.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 500 | U |
| 103-65-1 | n-Propylbenzene | | 500 | U |
| 95-49-8 | 2-Chlorotoluene | | 500 | U |
| 106-43-4 | 4-Chlorotoluene | | 500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 500 | U |
| 98-06-6 | tert-Butylbenzene | | 500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 500 | U |
| 135-98-8 | sec-Butylbenzene | | 500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 500 | U |
| 104-51-8 | n-Butylbenzene | | 500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 500 | U |
| 87-68-3 | Hexachlorobutadiene | | 500 | U |
| 91-20-3 | Naphthalene | | 500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 500 | U |
| 1634-04-4 | MTBE | | 500 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 500 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-02

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_02 0.10 ml
 Sample wt/vol: 0.1 (g/ml) ML Lab File ID: 7519_02.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 100.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|----|
| 75-71-8 | Dichlorodifluoromethane | | 500 | U |
| 74-87-3 | Chloromethane | | 500 | U |
| 75-01-4 | Vinyl chloride | | 500 | U |
| 74-83-9 | Bromomethane | | 500 | U |
| 75-00-3 | Chloroethane | | 500 | U |
| 75-69-4 | Trichlorofluoromethane | | 500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 500 | U |
| 75-09-2 | Methylene chloride | | 500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 500 | U |
| 74-97-5 | Bromochloromethane | | 500 | U |
| 67-66-3 | Chloroform | | 500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 500 | U |
| 56-23-5 | Carbon tetrachloride | | 500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 500 | U |
| 71-43-2 | Benzene | | 500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 500 | U |
| 79-01-6 | Trichloroethene | | 500 | U |
| 78-87-5 | 1,2-Dichloropropane | | 500 | U |
| 74-95-3 | Dibromomethane | | 500 | U |
| 75-27-4 | Bromodichloromethane | | 500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 500 | U |
| 108-88-3 | Toluene | | 500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 500 | U |
| 127-18-4 | Tetrachloroethene | | 42000 | ED |
| 142-28-9 | 1,3-Dichloropropane | | 500 | U |
| 124-48-1 | Dibromochloromethane | | 500 | U |
| 108-90-7 | Chlorobenzene | | 500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 500 | U |
| 100-41-4 | Ethylbenzene | | 500 | U |
| 1330-20-7 | Xylene (para & meta) | | 500 | U |
| 95-47-6 | Xylene (Ortho) | | 500 | U |
| 100-42-5 | Styrene | | 500 | U |
| 75-25-2 | Bromoform | | 500 | U |
| 98-82-8 | Isopropylbenzene | | 500 | U |
| 108-86-1 | Bromobenzene | | 500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-02

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA.
 Matrix: (soil/water) WATER Lab Sample ID: 7519_02 0.10 ml
 Sample wt/vol: 0.1 (g/ml) ML Lab File ID: 7519_02.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 100.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 500 | U |
| 103-65-1 | n-Propylbenzene | | 500 | U |
| 95-49-8 | 2-Chlorotoluene | | 500 | U |
| 106-43-4 | 4-Chlorotoluene | | 500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 500 | U |
| 98-06-6 | tert-Butylbenzene | | 500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 500 | U |
| 135-98-8 | sec-Butylbenzene | | 500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 500 | U |
| 104-51-8 | n-Butylbenzene | | 500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 500 | U |
| 87-68-3 | Hexachlorobutadiene | | 500 | U |
| 91-20-3 | Naphthalene | | 500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 500 | U |
| 1634-04-4 | MTBE | | 500 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 500 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-03

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_03 0.10 ml
 Sample wt/vol: 0.1 (g/ml) ML Lab File ID: 7519_03.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 100.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|----|
| 75-71-8 | Dichlorodifluoromethane | | 500 | U |
| 74-87-3 | Chloromethane | | 500 | U |
| 75-01-4 | Vinyl chloride | | 500 | U |
| 74-83-9 | Bromomethane | | 500 | U |
| 75-00-3 | Chloroethane | | 500 | U |
| 75-69-4 | Trichlorofluoromethane | | 500 | U |
| 75-35-4 | 1,1-Dichloroethene | | 500 | U |
| 75-09-2 | Methylene chloride | | 500 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 500 | U |
| 75-34-3 | 1,1-Dichloroethane | | 500 | U |
| 594-20-7 | 2,2-Dichloropropane | | 500 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 500 | U |
| 74-97-5 | Bromochloromethane | | 500 | U |
| 67-66-3 | Chloroform | | 500 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 500 | U |
| 56-23-5 | Carbon tetrachloride | | 500 | U |
| 563-58-6 | 1,1-Dichloropropene | | 500 | U |
| 71-43-2 | Benzene | | 500 | U |
| 107-06-2 | 1,2-Dichloroethane | | 500 | U |
| 79-01-6 | Trichloroethene | | 150 | JD |
| 78-87-5 | 1,2-Dichloropropane | | 500 | U |
| 74-95-3 | Dibromomethane | | 500 | U |
| 75-27-4 | Bromodichloromethane | | 500 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 500 | U |
| 108-88-3 | Toluene | | 500 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 500 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 500 | U |
| 127-18-4 | Tetrachloroethene | | 40000 | ED |
| 142-28-9 | 1,3-Dichloropropane | | 500 | U |
| 124-48-1 | Dibromochloromethane | | 500 | U |
| 108-90-7 | Chlorobenzene | | 500 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 500 | U |
| 100-41-4 | Ethylbenzene | | 500 | U |
| 1330-20-7 | Xylene (para & meta) | | 500 | U |
| 95-47-6 | Xylene (Ortho) | | 500 | U |
| 100-42-5 | Styrene | | 500 | U |
| 75-25-2 | Bromoform | | 500 | U |
| 98-82-8 | Isopropylbenzene | | 500 | U |
| 108-86-1 | Bromobenzene | | 500 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7519-03

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7519_03 0.10 ml
 Sample wt/vol: 0.1 (g/ml) ML Lab File ID: 7519_03.D
 Level: (low/med) LOW Date Received: 01/10/05
 % Moisture: not dec. _____ Date Analyzed: 01/22/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 100.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 500 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 500 | U |
| 103-65-1 | n-Propylbenzene | | 500 | U |
| 95-49-8 | 2-Chlorotoluene | | 500 | U |
| 106-43-4 | 4-Chlorotoluene | | 500 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 500 | U |
| 98-06-6 | tert-Butylbenzene | | 500 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 500 | U |
| 135-98-8 | sec-Butylbenzene | | 500 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 500 | U |
| 99-87-6 | 4-Isopropyltoluene | | 500 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 500 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 500 | U |
| 104-51-8 | n-Butylbenzene | | 500 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 500 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 500 | U |
| 87-68-3 | Hexachlorobutadiene | | 500 | U |
| 91-20-3 | Naphthalene | | 500 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 500 | U |
| 1634-04-4 | MTBE | | 500 | U |
| 67-64-1 | Acetone | | 1000 | U |
| 75-15-0 | Carbon disulfide | | 500 | U |
| 78-93-3 | 2-Butanone | | 1000 | U |
| 109-99-9 | Tetrahydrofuran | | 1000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 1000 | U |
| 591-78-6 | 2-Hexanone | | 1000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 1000 | U |

Data File : C:\HPCHEM\1\DATA\0102207\7519_01.D

Acq On : 22 Jan 2007 12:27 pm

Sample : 7519_01 0.10 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Jan 23 14:57 2007

Vial: 5

Operator: RR/AS

Inst : GC/MS Ins

Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:54:07 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.73 | 96 | 293670 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 0.00 | 117 | 0 | 0.00 | ug/L | -20.12 |
| 58) 1,4-Dichlorobenzene-d4 | 0.00 | 152 | 0 | 0.00 | ug/L | -24.35 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|------|
| 16) Dibromofluoromethane | 13.28 | 113 | 85050 | 26.69 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 106.76% | |
| 19) 1,2-Dichloroethane-d4 | 14.25 | 65 | 80179 | 24.83 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 99.32% | |
| 28) Toluene-d8 | 17.42 | 98 | 341762 | 25.77 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 103.08% | |
| 45) 4-Bromofluorobenzene | 0.00 | 95 | 0 | 0.00 | ug/L | |
| Spiked Amount | 25.000 | | Recovery | = | 0.00% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | d | |

(#) = qualifier out of range (m) = manual integration

7519_01.D 01_12_07.M

Thu Jan 25 14:18:06 2007

Data File : C:\HPCHEM\1\DATA\0102207\7519_01.D

Vial: 5

Acq On : 22 Jan 2007 12:27 pm

Operator: RR/AS

Sample : 7519_01 0.10 ml

Inst : GC/MS Ins

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 23 14:57 2007

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:54:07 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|--------|------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 2244882 | 391.29 | ug/L | 100 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 9.06 | 43 | 70745 | N.D. | | |
| 69) Carbon disulfide | 9.06 | 76 | 248583 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

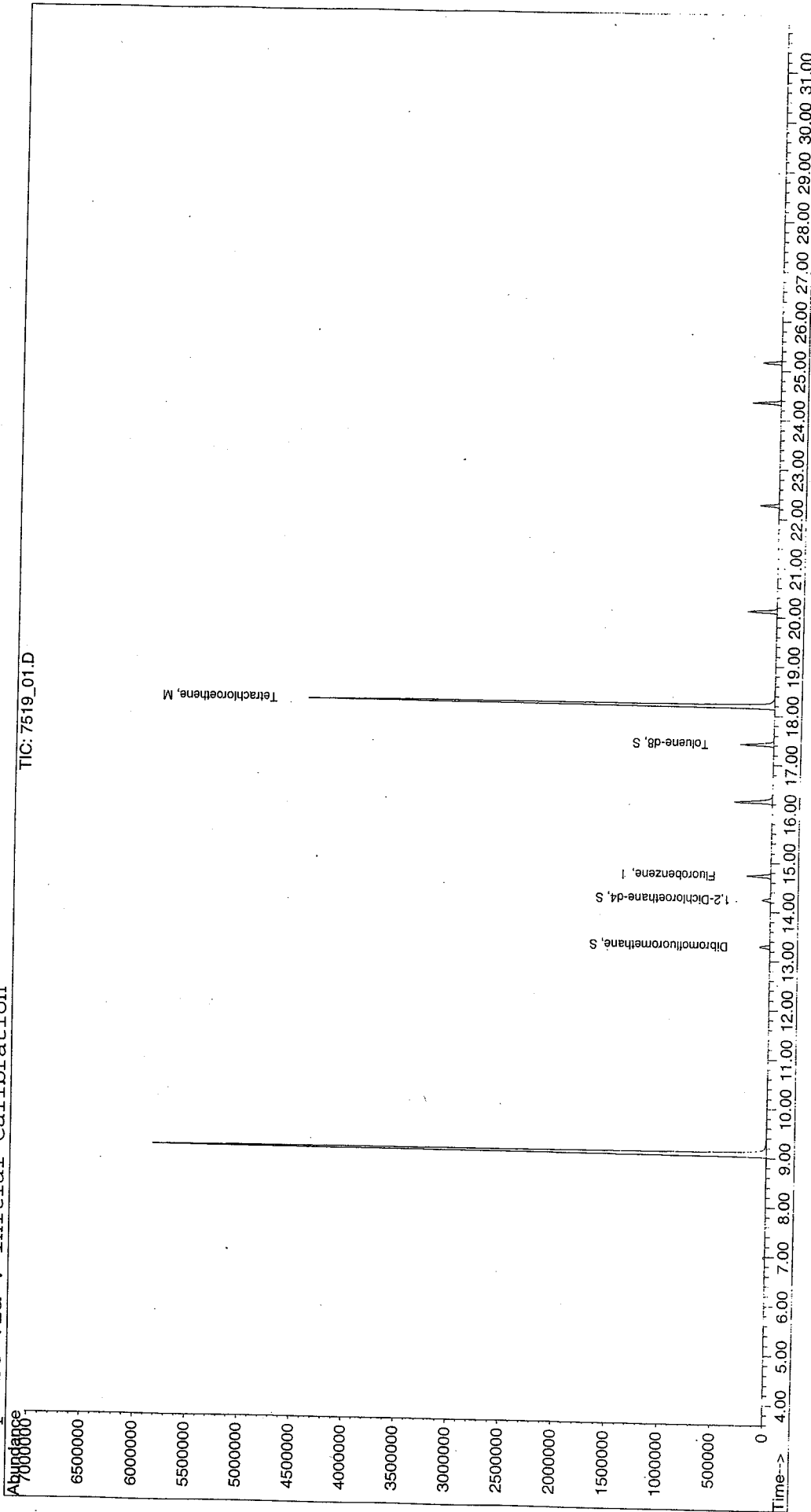
7519_01.D 01_12_07.M Thu Jan 25 14:18:09 2007

Page 2

Quantitation Report

Data File : C:\HPCHEM\1\DATA\0102207\7519_01.D
Acq On : 22 Jan 2007 12:27 pm
Sample : 7519_01 0.10 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 23 14:57 2007
Vial: 5
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: 01_12_07.RES

Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:54:07 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\0102207\7519_02.D
 Acq On : 22 Jan 2007 1:06 pm
 Sample : 7519_02 0.10 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 23 14:59 2007

Vial: 6
 Operator: RR/AS
 Inst : GC/MS Ins
 Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Mon Jan 15 14:54:07 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.73 | 96 | 295750 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 20.12 | 117 | 261159 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.36 | 152 | 130228 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|--------|-------|------|--------------------|
| 16) Dibromofluoromethane | 13.29 | 113 | 85491 | 26.63 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 106.52% |
| 19) 1,2-Dichloroethane-d4 | 14.25 | 65 | 81692 | 25.12 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 100.48% |
| 28) Toluene-d8 | 17.42 | 98 | 348016 | 26.06 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 104.24% |
| 45) 4-Bromofluorobenzene | 22.27 | 95 | 92442 | 20.73 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | | | | |
| | | | | | | Recovery = 82.92% |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | d | |

(#) = qualifier out of range (m) = manual integration
 7519_02.D 01_12_07.M Thu Jan 25 14:18:17 2007

Data File : C:\HPCHEM\1\DATA\0102207\7519_02.D
Acq On : 22 Jan 2007 1:06 pm
Sample : 7519_02 0.10 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 23 14:59 2007

Vial: 6
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:54:07 2007
Response via : Initial Calibration
DataAcq Meth : RUN

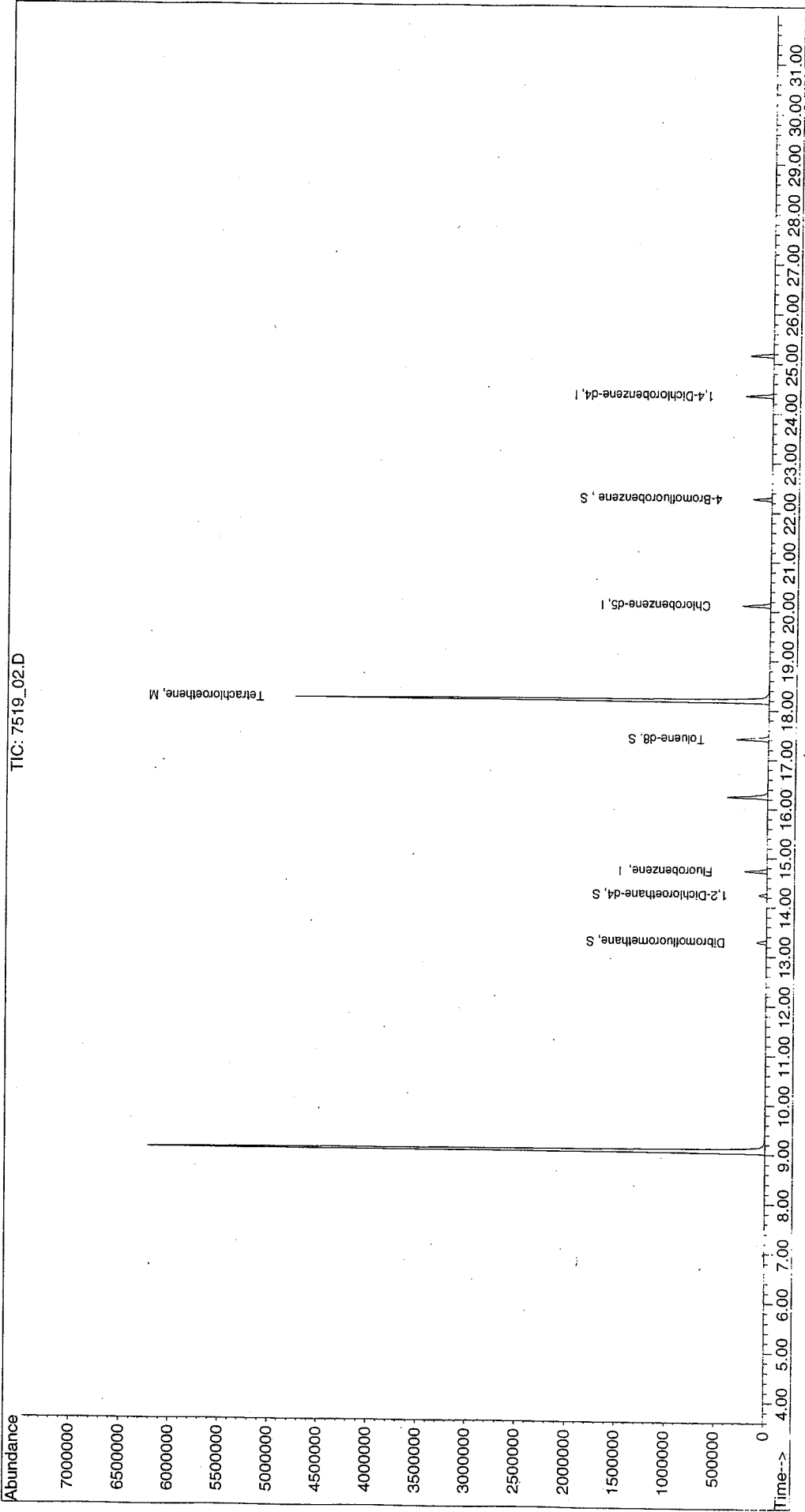
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|--------|------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 2439524 | 422.23 | ug/L | 99 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | d | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | d | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | d | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration
7519_02.D 01_12_07.M Thu Jan 25 14:18:19 2007

Quantitation Report

Data File : C:\HPCHEM\1\DATA\0102207\7519_02.D
Acq On : 22 Jan 2007 1:06 pm
Sample : 7519_02 0.10 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 23 14:59 2007
Vial: 6
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: 01_12_07.REB

Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:54:07 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\0102207\7519_03.D
 Acq On : 22 Jan 2007 1:44 pm
 Sample : 7519_03 0.10 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Jan 24 11:11 2007

Vial: 7
 Operator: RR/AS
 Inst : GC/MS Ins
 Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Mon Jan 15 14:54:07 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN

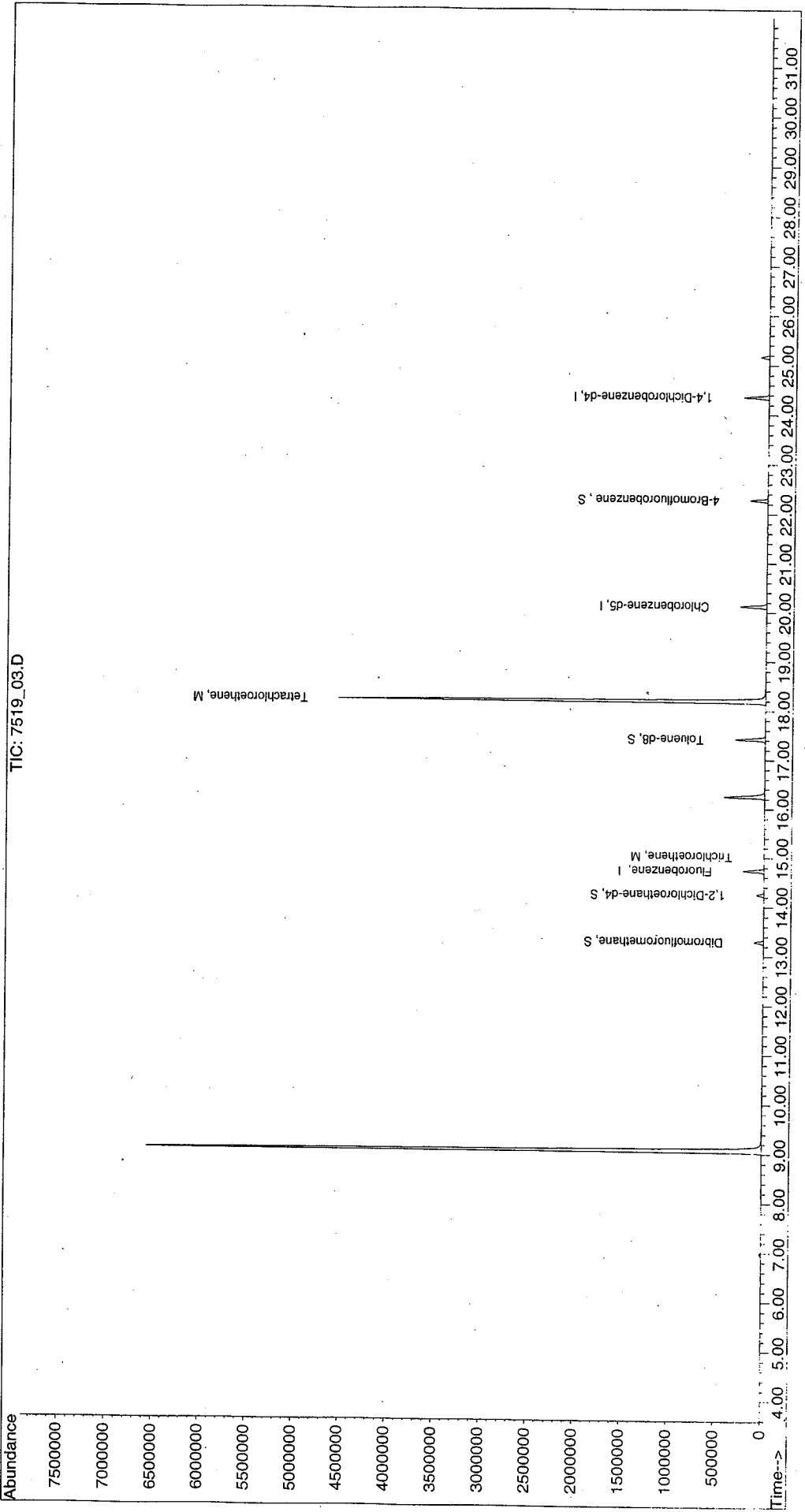
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|--------|------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 2326948 | 402.24 | ug/L | 99 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | d | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | d | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | d | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration
 7519_03.D 01_12_07.M Thu Jan 25 14:18:30 2007

Quantitation Report

Data File : C:\HPCHEM\1\DATA\0102207\7519_03.D
Acq On : 22 Jan 2007 1:44 pm
Sample : 7519_03 0.10 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Jan 24 11:11 2007
Vial: 7
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: 01_12_07.RES

Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:54:07 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\0102207\7519_3.D

Vial: 9

Acq On : 22 Jan 2007 10:07 pm

Operator: RR/AS

Sample : 7519-3 0.020ml

Inst : GC/MS Ins

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Jan 24 11:13 2007

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:54:07 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|--------------------|------|------|----------|------|-------|----------|
|--------------------|------|------|----------|------|-------|----------|

| | | | | | | |
|----------------------------|-------|-----|--------|-------|------|------|
| 1) Fluorobenzene | 14.73 | 96 | 295454 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 20.12 | 117 | 245383 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.35 | 152 | 117674 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|------|
| 16) Dibromofluoromethane | 13.29 | 113 | 77548 | 24.18 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 96.72% | |
| 19) 1,2-Dichloroethane-d4 | 14.24 | 65 | 74269 | 22.86 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 91.44% | |
| 28) Toluene-d8 | 17.42 | 98 | 335295 | 25.13 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 100.52% | |
| 45) 4-Bromofluorobenzene | 22.28 | 95 | 81836 | 19.53 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 78.12% | |

Target Compounds

| | | | | | Qvalue |
|-------------------------------|------|-----|---|------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | d |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | |

(#) = qualifier out of range (m) = manual integration

7519_3.D 01_12_07.M

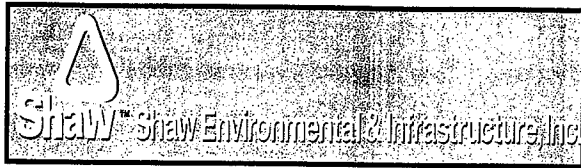
Thu Jan 25 14:18:59 2007

MEEP Results with non-heated samples
Using Henry's law

| Run Date | Sample | Liquid Volume (mL) | Headspace Volume (mL) | Room temp. (°C) | Room temp. (K) | Obs | | | Obs Measured Ethane (ppmV) | Obs Measured Ethene (ppmV) | Obs Measured Propane (ppmV) | dilution | Calculated Aq. Concentrations | | | Calculated umolar Conc. | | |
|-----------|-------------|--------------------|-----------------------|-----------------|----------------|----------------|---------------|----------------|----------------------------|----------------------------|-----------------------------|----------|-------------------------------|---------------|----------------|-------------------------|-------------|--------------|
| | | | | | | Methane (ppmV) | Ethane (ppmV) | Propane (ppmV) | | | | | Methane (ug/L) | Ethane (ug/L) | Propane (ug/L) | Methane (uM) | Ethane (uM) | Propane (uM) |
| 1/29/2007 | blank | 32 | 10 | 21.67 | 294.7 | 32.374 | 0 | 0 | 0 | 0 | 1.0 | 8.09 | 0.00 | 0.00 | 0.51 | 0.00 | 0.00 | |
| | 100ppmv | 32 | 10 | 21.67 | 294.7 | 112.436 | 97.921 | 97.61 | 95.408 | 0 | 1.0 | 26.09 | 43.85 | 50.65 | 1.76 | 1.46 | 1.81 | |
| | 7519-1 | 3.75 | 1.25 | 21.67 | 294.7 | 74.427 | 4.428 | 0 | 0 | 0 | 1.0 | 19.82 | 2.10 | 0.00 | 1.23 | 0.07 | 0.00 | |
| | 7519-2 | 3.75 | 1.25 | 21.67 | 294.7 | 78.804 | 4.502 | 0 | 0 | 0 | 1.0 | 20.77 | 2.13 | 0.00 | 1.30 | 0.07 | 0.00 | |
| | 7519-3 | 3.75 | 1.25 | 21.67 | 294.7 | 85.485 | 4.131 | 0 | 0 | 0 | 1.0 | 22.54 | 1.96 | 0.00 | 1.41 | 0.07 | 0.00 | |
| | 7519-4 | 3.75 | 1.25 | 21.67 | 294.7 | 82.104 | 4.217 | 0 | 0 | 0 | 1.0 | 21.64 | 2.00 | 0.00 | 1.35 | 0.07 | 0.00 | |
| | 7519-5 | 3.75 | 1.25 | 21.67 | 294.7 | 85.04 | 4.35 | 0 | 0 | 0 | 1.0 | 25.05 | 2.08 | 0.00 | 1.57 | 0.07 | 0.00 | |
| | 7519-6 | 3.75 | 1.25 | 21.67 | 294.7 | 85.248 | 4.511 | 0 | 0 | 0 | 1.0 | 22.47 | 2.14 | 0.00 | 1.40 | 0.07 | 0.00 | |
| | 7519-7 | 3.75 | 1.25 | 21.67 | 294.7 | 83.3 | 4.386 | 0 | 0 | 0 | 1.0 | 21.96 | 2.08 | 0.00 | 1.37 | 0.07 | 0.00 | |
| | 7519-8 | 3.75 | 1.25 | 21.67 | 294.7 | 39.381 | 5.082 | 0 | 0 | 0 | 1.0 | 10.38 | 2.41 | 0.00 | 0.65 | 0.08 | 0.00 | |
| | 7519-9 | 3.75 | 1.25 | 21.67 | 294.7 | 60.906 | 4.447 | 0 | 0 | 0 | 1.0 | 16.06 | 2.11 | 0.00 | 1.00 | 0.07 | 0.00 | |
| | 7519-10 | 3.75 | 1.25 | 21.67 | 294.7 | 60.222 | 9.003 | 0 | 0 | 0 | 1.0 | 15.88 | 4.26 | 0.00 | 0.99 | 0.14 | 0.00 | |
| | blank | 32 | 10 | 21.67 | 294.7 | 9.154 | 0 | 0 | 0 | 0 | 1.0 | 2.29 | 0.00 | 0.00 | 0.14 | 0.00 | 0.00 | |
| | 50ppmv | 32 | 10 | 21.67 | 294.7 | 51.205 | 52.166 | 50.58 | 51.733 | 0 | 1.0 | 12.79 | 23.36 | 26.24 | 0.80 | 0.78 | 0.94 | |
| | 7519-11 | 3.75 | 1.25 | 21.67 | 294.7 | 68.538 | 8.052 | 0 | 0 | 0 | 1.0 | 18.07 | 3.81 | 0.00 | 1.13 | 0.13 | 0.00 | |
| | 7519-11 | 3.75 | 1.25 | 21.67 | 294.7 | 70.623 | 12.135 | 0 | 1.8 | 0 | 1.0 | 18.62 | 5.75 | 0.00 | 1.16 | 0.19 | 0.00 | |
| | 7519-12 MS | 3.75 | 1.25 | 21.67 | 294.7 | 170.238 | 105.873 | 97.222 | 95.157 | 0 | 1.0 | 44.88 | 50.14 | 52.79 | 2.80 | 1.67 | 1.89 | |
| | 7519-12 MSD | 3.75 | 1.25 | 21.67 | 294.7 | 171.865 | 104.782 | 97 | 96.545 | 0 | 1.0 | 45.33 | 49.63 | 52.67 | 2.83 | 1.65 | 1.88 | |
| | blank | 32 | 10 | 21.67 | 294.7 | 12.793 | 0 | 0 | 0 | 0 | 1.0 | 3.20 | 0.00 | 0.00 | 0.20 | 0.00 | 0.00 | |

MEE- henry's constant adjusted to temperature

Propane - henry's constant at 25 C



17 Princess Rd
Lawrenceville, New Jersey 08648
Tel: 609/895-5370
Fax: 609/895-1858

Volatile Organic Compound Data Summary Package

Prepared for
Gorham Textron

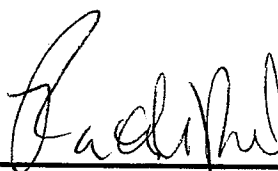
Lab ID
7543

Project Number 101960 02000000

Samples Received
8-Feb-07

Reported
28-Feb-07

NJDEP Certified Lab 11001

 2/28/07

Randi K Rothmel, PhD Date
Laboratory Director

1.0 Chain of Custody

Lawrenceville, NJ 08648
609-895-5340/ 609-895-1858



Shaw Environmental and Infrastructure Inc.

7543

Project Contact: Charles Schaefer, Chuck Condee
(Name & phone #)

Send Report To: Chuck Condee
Phone/Fax Number:
Address:
City/State:

CHAIN OF CUSTODY

Ref. Document #

Page 1 of 1

Project Number/Cost code: 101960 / 02

Project Name / Location: Gorham Textron

Send bill to:

Shipment date:
Lab Destination:
Lab Contact Name / ph. #:

| Lab No. | Sample ID Number | Sample Description | Collection Information | | # of containers | Container Type | Preservative | | | | | VOC | MER | Antions | VFA | Analyses Requested | Turn Around Time Requested |
|---------|------------------|--------------------|------------------------|-------|-----------------|----------------|--------------|-----|------|------------------|--------------------------------|-----|-----|---------|-----|--------------------|----------------------------|
| | | | Date | Time | | | G/C | HCL | NaOH | HNO ₃ | H ₂ SO ₄ | | | | | | |
| 1 | Killed A | " " | 02/08/07 | 12:00 | Aq | 1 | 10mL | X | | | | | | X | | | |
| 2 | Killed B | " " | | | Aq | 1 | 3.75mL | X | | | | | | X | | | |
| 3 | Killed C | " " | | | Aq | 1 | 10mL | X | | | | | | X | | | |
| 4 | Live A | " " | | | Aq | 1 | 3.75mL | X | | | | | | X | | | |
| 5 | Live B | " " | | | Aq | 1 | 3.75mL | X | | | | | | X | | | |
| 6 | Live C | " " | | | Aq | 1 | 10mL | X | | | | | | X | | | |
| 7 | Bioaug 1 LA | " " | | | Aq | 1 | 10mL | X | | | | | | X | | | |
| 8 | Bioaug 1 LB | " " | | | Aq | 1 | 3.75mL | X | | | | | | X | | | |
| 9 | Bioaug 1 LC | " " | | | Aq | 1 | 3.75mL | X | | | | | | X | | | |
| 10 | Bioaug 2 EA | " " | | | Aq | 1 | 3.75mL | X | | | | | | X | | | |
| 11 | Bioaug 2 EB | " " | | | Aq | 1 | 10mL | X | | | | | | X | | | |
| 12 | Bioaug 2 EC | " " | | | Aq | 1 | 3.75mL | X | | | | | | X | | | |
| 13 | Killed D | " " | | | Aq | 1 | 3.75mL | X | | | | | | X | | | |
| 14 | Live D | " " | | | Aq | 1 | 15mL | X | | | | | | X | | | |
| 15 | Bioaug 1 LD | " " | | | Aq | 1 | 15mL | X | | | | | | X | | | |
| 16 | Bioaug 2 ED | " " | | | Aq | 1 | 15mL | X | | | | | | X | | | |

Special Instructions: MEE bottles hold 5 mL with 3.75 mL sample inside.

Known Waste Stream Code:
RCRA PCB/dioxin PAH/oil
QC/Data Package Level Required: I II III IV

Flammable
Corrosive
Reactive
GIS EDD
Preliminary data

G/C Codes
C = Composite
G = Grab

QC Package Codes

Relinquished By: *Mary Higum*
Date: 2/8/07
Time: 1:00pm

Received By: *[Signature]*
Date: 02/08/07
Time: 1:46

Level II = data summary + basic QC
Level III = New Jersey QC reduced deliverable
Level IV = Full deliverable CLP package
Cooler temperature upon arrival at Lab:

Shaw E&I Analytical and Treatability Laboratories Internal Chain of Custody

Pg _____ of _____

Lab ID **7543**

Client:

Date Received:

| Sample ID | Parameter | Bottle Type | Preservative | Date/Time Removed | Relinquishing Custodian Initials | Receiving Analyst Initials | Date/Time Returned | Receiving Custodian Initials | Relinquishing Analyst Initials |
|-----------|-----------|-------------|--------------|-------------------|----------------------------------|----------------------------|--------------------|------------------------------|--------------------------------|
| 7543-1 | VOCs | 1-10-1 | | 02-08-07 9:00 | AM | AM | 02-08-07 17:00 | PM | AM |
| 2 | | | | | | | | | |
| 3 | | | | | | | | | |
| 4 | | | | | | | | | |
| 5 | | | | | | | | | |
| 6 | | | | | | | | | |
| 7 | | | | | | | | | |
| 8 | | | | | | | | | |
| 9 | | | | | | | | | |
| 10 | | | | | | | | | |
| 11 | | | | | | | | | |
| 7543-12 | VOCs | 1-10-1 | | 02-08-07 9:00 | AM | AM | 02-08-07 17:00 | PM | AM |
| 7543-1 | MEERs | 1-10-1 | | 2/12/07 9:00 | AM | AM | 2/12/07 17:30 | PM | AM |
| 2 | | | | | | | | | |
| 3 | | | | | | | | | |
| 4 | | | | | | | | | |
| 5 | | | | | | | | | |
| 6 | | | | | | | | | |
| 7 | | | | | | | | | |
| 8 | | | | | | | | | |
| 9 | | | | | | | | | |
| 10 | | | | | | | | | |
| 11 | | | | | | | | | |
| 7543-12 | MEERs | 1-10-1 | | 2/12/07 17:00 | AM | AM | 2/14/07 16:00 | PM | AM |
| 7543-13 | MEERs | 1-10-1 | | 2/12/07 17:00 | AM | AM | 2/14/07 16:00 | PM | AM |
| 14 | | | | | | | | | |
| 15 | | | | | | | | | |
| 7543-16 | MEERs | 1-10-1 | | 2/14/07 17:00 | AM | AM | 2/14/07 16:00 | PM | AM |
| 7543-17 | VFA | 1-10-1 | | 2/14/07 17:00 | AM | AM | 2/14/07 16:00 | PM | AM |
| 18 | | | | | | | | | |
| 19 | | | | | | | | | |
| 7543-15 | VFA | 1-10-1 | | 2/14/07 17:00 | AM | AM | 2/14/07 16:00 | PM | AM |

2.0. Sample Summaries

Sample summaries are enclosed

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-1

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_1 0.010 ml
 Sample wt/vol: 0.010 (g/ml) ML Lab File ID: 7543_1.D
 Level: (low/med) LOW Date Received: 02/07/07
 % Moisture: not dec. _____ Date Analyzed: 02/08/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 5000 | U |
| 74-87-3 | Chloromethane | | 5000 | U |
| 75-01-4 | Vinyl chloride | | 5000 | U |
| 74-83-9 | Bromomethane | | 5000 | U |
| 75-00-3 | Chloroethane | | 5000 | U |
| 75-69-4 | Trichlorofluoromethane | | 5000 | U |
| 75-35-4 | 1,1-Dichloroethene | | 5000 | U |
| 75-09-2 | Methylene chloride | | 5000 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 5000 | U |
| 75-34-3 | 1,1-Dichloroethane | | 5000 | U |
| 594-20-7 | 2,2-Dichloropropane | | 5000 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 5000 | U |
| 74-97-5 | Bromochloromethane | | 5000 | U |
| 67-66-3 | Chloroform | | 5000 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 5000 | U |
| 56-23-5 | Carbon tetrachloride | | 5000 | U |
| 563-58-6 | 1,1-Dichloropropene | | 5000 | U |
| 71-43-2 | Benzene | | 5000 | U |
| 107-06-2 | 1,2-Dichloroethane | | 5000 | U |
| 79-01-6 | Trichloroethene | | 5000 | U |
| 78-87-5 | 1,2-Dichloropropane | | 5000 | U |
| 74-95-3 | Dibromomethane | | 5000 | U |
| 75-27-4 | Bromodichloromethane | | 5000 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 5000 | U |
| 108-88-3 | Toluene | | 5000 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 5000 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 5000 | U |
| 127-18-4 | Tetrachloroethene | | 26000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 5000 | U |
| 124-48-1 | Dibromochloromethane | | 5000 | U |
| 108-90-7 | Chlorobenzene | | 5000 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 5000 | U |
| 100-41-4 | Ethylbenzene | | 5000 | U |
| 1330-20-7 | Xylene (para & meta) | | 5000 | U |
| 95-47-6 | Xylene (Ortho) | | 5000 | U |
| 100-42-5 | Styrene | | 5000 | U |
| 75-25-2 | Bromoform | | 5000 | U |
| 98-82-8 | Isopropylbenzene | | 5000 | U |
| 108-86-1 | Bromobenzene | | 5000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-1

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_1 0.010 ml
 Sample wt/vol: 0.01^o (g/ml) ML Lab File ID: 7543_1.D
 Level: (low/med) LOW Date Received: 02/07/07
 % Moisture: not dec. _____ Date Analyzed: 02/08/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

CAS NO. - COMPOUND (ug/L or ug/Kg) UG/L Q

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5000 | U | U |
| 96-18-4 | 1,2,3-Trichloropropane | 5000 | U | U |
| 103-65-1 | n-Propylbenzene | 5000 | U | U |
| 95-49-8 | 2-Chlorotoluene | 5000 | U | U |
| 106-43-4 | 4-Chlorotoluene | 5000 | U | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | 5000 | U | U |
| 98-06-6 | tert-Butylbenzene | 5000 | U | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | 5000 | U | U |
| 135-98-8 | sec-Butylbenzene | 5000 | U | U |
| 541-73-1 | 1,3-Dichlorobenzene | 5000 | U | U |
| 99-87-6 | 4-Isopropyltoluene | 5000 | U | U |
| 106-46-7 | 1,4-Dichlorobenzene | 5000 | U | U |
| 95-50-1 | 1,2-Dichlorobenzene | 5000 | U | U |
| 104-51-8 | n-Butylbenzene | 5000 | U | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | 5000 | U | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | 5000 | U | U |
| 87-68-3 | Hexachlorobutadiene | 5000 | U | U |
| 91-20-3 | Naphthalene | 5000 | U | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | 5000 | U | U |
| 1634-04-4 | MTBE | 5000 | U | U |
| 67-64-1 | Acetone | 10000 | U | U |
| 75-15-0 | Carbon disulfide | 5000 | U | U |
| 78-93-3 | 2-Butanone | 10000 | U | U |
| 109-99-9 | Tetrahydrofuran | 10000 | U | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 10000 | U | U |
| 591-78-6 | 2-Hexanone | 10000 | U | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | 10000 | U | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-2

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_2 0.010 ml
 Sample wt/vol: 0.010 (g/ml) ML Lab File ID: 7543_2.D
 Level: (low/med) LOW Date Received: 02/07/07
 % Moisture: not dec. _____ Date Analyzed: 02/08/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | 5000 | | U |
| 74-87-3 | Chloromethane | 5000 | | U |
| 75-01-4 | Vinyl chloride | 5000 | | U |
| 74-83-9 | Bromomethane | 5000 | | U |
| 75-00-3 | Chloroethane | 5000 | | U |
| 75-69-4 | Trichlorofluoromethane | 5000 | | U |
| 75-35-4 | 1,1-Dichloroethene | 5000 | | U |
| 75-09-2 | Methylene chloride | 5000 | | U |
| 156-60-5 | trans-1,2-Dichloroethene | 5000 | | U |
| 75-34-3 | 1,1-Dichloroethane | 5000 | | U |
| 594-20-7 | 2,2-Dichloropropane | 5000 | | U |
| 156-59-2 | cis-1,2-Dichloroethene | 5000 | | U |
| 74-97-5 | Bromochloromethane | 5000 | | U |
| 67-66-3 | Chloroform | 5000 | | U |
| 71-55-6 | 1,1,1-Trichloroethane | 5000 | | U |
| 56-23-5 | Carbon tetrachloride | 5000 | | U |
| 563-58-6 | 1,1-Dichloropropene | 5000 | | U |
| 71-43-2 | Benzene | 5000 | | U |
| 107-06-2 | 1,2-Dichloroethane | 5000 | | U |
| 79-01-6 | Trichloroethene | 5000 | | U |
| 78-87-5 | 1,2-Dichloropropane | 5000 | | U |
| 74-95-3 | Dibromomethane | 5000 | | U |
| 75-27-4 | Bromodichloromethane | 5000 | | U |
| 10061-02-5 | cis-1,3-Dichloropropene | 5000 | | U |
| 108-88-3 | Toluene | 5000 | | U |
| 10061-01-6 | trans-1,3-Dichloropropene | 5000 | | U |
| 79-00-5 | 1,1,2-Trichloroethane | 5000 | | U |
| 127-18-4 | Tetrachloroethene | 24000 | | D |
| 142-28-9 | 1,3-Dichloropropane | 5000 | | U |
| 124-48-1 | Dibromochloromethane | 5000 | | U |
| 108-90-7 | Chlorobenzene | 5000 | | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 5000 | | U |
| 100-41-4 | Ethylbenzene | 5000 | | U |
| 1330-20-7 | Xylene (para & meta) | 5000 | | U |
| 95-47-6 | Xylene (Ortho) | 5000 | | U |
| 100-42-5 | Styrene | 5000 | | U |
| 75-25-2 | Bromoform | 5000 | | U |
| 98-82-8 | Isopropylbenzene | 5000 | | U |
| 108-86-1 | Bromobenzene | 5000 | | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-2

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_2 0.010 ml
 Sample wt/vol: 0.010 (g/ml) ML Lab File ID: 7543_2.D
 Level: (low/med) LOW Date Received: 02/07/07
 % Moisture: not dec. _____ Date Analyzed: 02/08/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5000 | U | |
| 96-18-4 | 1,2,3-Trichloropropane | 5000 | U | |
| 103-65-1 | n-Propylbenzene | 5000 | U | |
| 95-49-8 | 2-Chlorotoluene | 5000 | U | |
| 106-43-4 | 4-Chlorotoluene | 5000 | U | |
| 108-67-8 | 1,3,5-Trimethylbenzene | 5000 | U | |
| 98-06-6 | tert-Butylbenzene | 5000 | U | |
| 95-63-6 | 1,2,4-Trimethylbenzene | 5000 | U | |
| 135-98-8 | sec-Butylbenzene | 5000 | U | |
| 541-73-1 | 1,3-Dichlorobenzene | 5000 | U | |
| 99-87-6 | 4-Isopropyltoluene | 5000 | U | |
| 106-46-7 | 1,4-Dichlorobenzene | 5000 | U | |
| 95-50-1 | 1,2-Dichlorobenzene | 5000 | U | |
| 104-51-8 | n-Butylbenzene | 5000 | U | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | 5000 | U | |
| 120-82-1 | 1,2,4-Trichlorobenzene | 5000 | U | |
| 87-68-3 | Hexachlorobutadiene | 5000 | U | |
| 91-20-3 | Naphthalene | 5000 | U | |
| 87-61-6 | 1,2,3-Trichlorobenzene | 5000 | U | |
| 1634-04-4 | MTBE | 5000 | U | |
| 67-64-1 | Acetone | 10000 | U | |
| 75-15-0 | Carbon disulfide | 5000 | U | |
| 78-93-3 | 2-Butanone | 10000 | U | |
| 109-99-9 | Tetrahydrofuran | 10000 | U | |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 10000 | U | |
| 591-78-6 | 2-Hexanone | 10000 | U | |
| 110-75-8 | 2-Chloroethyl vinyl ether | 10000 | U | |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-3

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_3 0.010 ml
 Sample wt/vol: 0.010 (g/ml) ML Lab File ID: 7543_3.D
 Level: (low/med) LOW Date Received: 02/07/07
 % Moisture: not dec. _____ Date Analyzed: 02/08/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | 5000 | | U |
| 74-87-3 | Chloromethane | 5000 | | U |
| 75-01-4 | Vinyl chloride | 5000 | | U |
| 74-83-9 | Bromomethane | 5000 | | U |
| 75-00-3 | Chloroethane | 5000 | | U |
| 75-69-4 | Trichlorofluoromethane | 5000 | | U |
| 75-35-4 | 1,1-Dichloroethene | 5000 | | U |
| 75-09-2 | Methylene chloride | 5000 | | U |
| 156-60-5 | trans-1,2-Dichloroethene | 5000 | | U |
| 75-34-3 | 1,1-Dichloroethane | 5000 | | U |
| 594-20-7 | 2,2-Dichloropropane | 5000 | | U |
| 156-59-2 | cis-1,2-Dichloroethene | 5000 | | U |
| 74-97-5 | Bromochloromethane | 5000 | | U |
| 67-66-3 | Chloroform | 5000 | | U |
| 71-55-6 | 1,1,1-Trichloroethane | 5000 | | U |
| 56-23-5 | Carbon tetrachloride | 5000 | | U |
| 563-58-6 | 1,1-Dichloropropene | 5000 | | U |
| 71-43-2 | Benzene | 5000 | | U |
| 107-06-2 | 1,2-Dichloroethane | 5000 | | U |
| 79-01-6 | Trichloroethene | 5000 | | U |
| 78-87-5 | 1,2-Dichloropropane | 5000 | | U |
| 74-95-3 | Dibromomethane | 5000 | | U |
| 75-27-4 | Bromodichloromethane | 5000 | | U |
| 10061-02-5 | cis-1,3-Dichloropropene | 5000 | | U |
| 108-88-3 | Toluene | 5000 | | U |
| 10061-01-6 | trans-1,3-Dichloropropene | 5000 | | U |
| 79-00-5 | 1,1,2-Trichloroethane | 5000 | | U |
| 127-18-4 | Tetrachloroethene | 21000 | | D |
| 142-28-9 | 1,3-Dichloropropane | 5000 | | U |
| 124-48-1 | Dibromochloromethane | 5000 | | U |
| 108-90-7 | Chlorobenzene | 5000 | | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 5000 | | U |
| 100-41-4 | Ethylbenzene | 5000 | | U |
| 1330-20-7 | Xylene (para & meta) | 5000 | | U |
| 95-47-6 | Xylene (Ortho) | 5000 | | U |
| 100-42-5 | Styrene | 5000 | | U |
| 75-25-2 | Bromoform | 5000 | | U |
| 98-82-8 | Isopropylbenzene | 5000 | | U |
| 108-86-1 | Bromobenzene | 5000 | | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-3

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_3 0.010 ml
 Sample wt/vol: 0.010 (g/ml) ML Lab File ID: 7543_3.D
 Level: (low/med) LOW Date Received: 02/07/07
 % Moisture: not dec. _____ Date Analyzed: 02/08/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5000 | U | |
| 96-18-4 | 1,2,3-Trichloropropane | 5000 | U | |
| 103-65-1 | n-Propylbenzene | 5000 | U | |
| 95-49-8 | 2-Chlorotoluene | 5000 | U | |
| 106-43-4 | 4-Chlorotoluene | 5000 | U | |
| 108-67-8 | 1,3,5-Trimethylbenzene | 5000 | U | |
| 98-06-6 | tert-Butylbenzene | 5000 | U | |
| 95-63-6 | 1,2,4-Trimethylbenzene | 5000 | U | |
| 135-98-8 | sec-Butylbenzene | 5000 | U | |
| 541-73-1 | 1,3-Dichlorobenzene | 5000 | U | |
| 99-87-6 | 4-Isopropyltoluene | 5000 | U | |
| 106-46-7 | 1,4-Dichlorobenzene | 5000 | U | |
| 95-50-1 | 1,2-Dichlorobenzene | 5000 | U | |
| 104-51-8 | n-Butylbenzene | 5000 | U | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | 5000 | U | |
| 120-82-1 | 1,2,4-Trichlorobenzene | 5000 | U | |
| 87-68-3 | Hexachlorobutadiene | 5000 | U | |
| 91-20-3 | Naphthalene | 5000 | U | |
| 87-61-6 | 1,2,3-Trichlorobenzene | 5000 | U | |
| 1634-04-4 | MTBE | 5000 | U | |
| 67-64-1 | Acetone | 10000 | U | |
| 75-15-0 | Carbon disulfide | 5000 | U | |
| 78-93-3 | 2-Butanone | 10000 | U | |
| 109-99-9 | Tetrahydrofuran | 10000 | U | |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 10000 | U | |
| 591-78-6 | 2-Hexanone | 10000 | U | |
| 110-75-8 | 2-Chloroethyl vinyl ether | 10000 | U | |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-4

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA.
 Matrix: (soil/water) WATER Lab Sample ID: 7543_4 0.010 ml
 Sample wt/vol: 0.01 (g/ml) ML Lab File ID: 7543_4.D
 Level: (low/med) LOW Date Received: 02/07/07
 % Moisture: not dec. _____ Date Analyzed: 02/08/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

CAS NO. - COMPOUND (ug/L or ug/Kg) UG/L Q

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 5000 | U |
| 74-87-3 | Chloromethane | | 5000 | U |
| 75-01-4 | Vinyl chloride | | 5000 | U |
| 74-83-9 | Bromomethane | | 5000 | U |
| 75-00-3 | Chloroethane | | 5000 | U |
| 75-69-4 | Trichlorofluoromethane | | 5000 | U |
| 75-35-4 | 1,1-Dichloroethene | | 5000 | U |
| 75-09-2 | Methylene chloride | | 5000 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 5000 | U |
| 75-34-3 | 1,1-Dichloroethane | | 5000 | U |
| 594-20-7 | 2,2-Dichloropropane | | 5000 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 5000 | U |
| 74-97-5 | Bromochloromethane | | 5000 | U |
| 67-66-3 | Chloroform | | 5000 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 5000 | U |
| 56-23-5 | Carbon tetrachloride | | 5000 | U |
| 563-58-6 | 1,1-Dichloropropene | | 5000 | U |
| 71-43-2 | Benzene | | 5000 | U |
| 107-06-2 | 1,2-Dichloroethane | | 5000 | U |
| 79-01-6 | Trichloroethene | | 5000 | U |
| 78-87-5 | 1,2-Dichloropropane | | 5000 | U |
| 74-95-3 | Dibromomethane | | 5000 | U |
| 75-27-4 | Bromodichloromethane | | 5000 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 5000 | U |
| 108-88-3 | Toluene | | 5000 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 5000 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 5000 | U |
| 127-18-4 | Tetrachloroethene | | 19000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 5000 | U |
| 124-48-1 | Dibromochloromethane | | 5000 | U |
| 108-90-7 | Chlorobenzene | | 5000 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 5000 | U |
| 100-41-4 | Ethylbenzene | | 5000 | U |
| 1330-20-7 | Xylene (para & meta) | | 5000 | U |
| 95-47-6 | Xylene (Ortho) | | 5000 | U |
| 100-42-5 | Styrene | | 5000 | U |
| 75-25-2 | Bromoform | | 5000 | U |
| 98-82-8 | Isopropylbenzene | | 5000 | U |
| 108-86-1 | Bromobenzene | | 5000 | U |

.1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-4

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_4 0.010 ml
 Sample wt/vol: 0.01 (g/ml) ML Lab File ID: 7543_4.D
 Level: (low/med) LOW Date Received: 02/07/07
 % Moisture: not dec. _____ Date Analyzed: 02/08/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5000 | U | |
| 96-18-4 | 1,2,3-Trichloropropane | 5000 | U | |
| 103-65-1 | n-Propylbenzene | 5000 | U | |
| 95-49-8 | 2-Chlorotoluene | 5000 | U | |
| 106-43-4 | 4-Chlorotoluene | 5000 | U | |
| 108-67-8 | 1,3,5-Trimethylbenzene | 5000 | U | |
| 98-06-6 | tert-Butylbenzene | 5000 | U | |
| 95-63-6 | 1,2,4-Trimethylbenzene | 5000 | U | |
| 135-98-8 | sec-Butylbenzene | 5000 | U | |
| 541-73-1 | 1,3-Dichlorobenzene | 5000 | U | |
| 99-87-6 | 4-Isopropyltoluene | 5000 | U | |
| 106-46-7 | 1,4-Dichlorobenzene | 5000 | U | |
| 95-50-1 | 1,2-Dichlorobenzene | 5000 | U | |
| 104-51-8 | n-Butylbenzene | 5000 | U | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | 5000 | U | |
| 120-82-1 | 1,2,4-Trichlorobenzene | 5000 | U | |
| 87-68-3 | Hexachlorobutadiene | 5000 | U | |
| 91-20-3 | Naphthalene | 5000 | U | |
| 87-61-6 | 1,2,3-Trichlorobenzene | 5000 | U | |
| 1634-04-4 | MTBE | 5000 | U | |
| 67-64-1 | Acetone | 10000 | U | |
| 75-15-0 | Carbon disulfide | 5000 | U | |
| 78-93-3 | 2-Butanone | 10000 | U | |
| 109-99-9 | Tetrahydrofuran | 10000 | U | |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 10000 | U | |
| 591-78-6 | 2-Hexanone | 10000 | U | |
| 110-75-8 | 2-Chloroethyl vinyl ether | 10000 | U | |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-5

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_5 0.010 ml
 Sample wt/vol: 0.0 / 0 (g/ml) ML Lab File ID: 7543_5.D
 Level: (low/med) LOW Date Received: 02/07/07
 % Moisture: not dec. _____ Date Analyzed: 02/08/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | 5000 | | U |
| 74-87-3 | Chloromethane | 5000 | | U |
| 75-01-4 | Vinyl chloride | 5000 | | U |
| 74-83-9 | Bromomethane | 5000 | | U |
| 75-00-3 | Chloroethane | 5000 | | U |
| 75-69-4 | Trichlorofluoromethane | 5000 | | U |
| 75-35-4 | 1,1-Dichloroethene | 5000 | | U |
| 75-09-2 | Methylene chloride | 5000 | | U |
| 156-60-5 | trans-1,2-Dichloroethene | 5000 | | U |
| 75-34-3 | 1,1-Dichloroethane | 5000 | | U |
| 594-20-7 | 2,2-Dichloropropane | 5000 | | U |
| 156-59-2 | cis-1,2-Dichloroethene | 5000 | | U |
| 74-97-5 | Bromochloromethane | 5000 | | U |
| 67-66-3 | Chloroform | 5000 | | U |
| 71-55-6 | 1,1,1-Trichloroethane | 5000 | | U |
| 56-23-5 | Carbon tetrachloride | 5000 | | U |
| 563-58-6 | 1,1-Dichloropropene | 5000 | | U |
| 71-43-2 | Benzene | 5000 | | U |
| 107-06-2 | 1,2-Dichloroethane | 5000 | | U |
| 79-01-6 | Trichloroethene | 5000 | | U |
| 78-87-5 | 1,2-Dichloropropane | 5000 | | U |
| 74-95-3 | Dibromomethane | 5000 | | U |
| 75-27-4 | Bromodichloromethane | 5000 | | U |
| 10061-02-5 | cis-1,3-Dichloropropene | 5000 | | U |
| 108-88-3 | Toluene | 5000 | | U |
| 10061-01-6 | trans-1,3-Dichloropropene | 5000 | | U |
| 79-00-5 | 1,1,2-Trichloroethane | 5000 | | U |
| 127-18-4 | Tetrachloroethene | 22000 | | D |
| 142-28-9 | 1,3-Dichloropropane | 5000 | | U |
| 124-48-1 | Dibromochloromethane | 5000 | | U |
| 108-90-7 | Chlorobenzene | 5000 | | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 5000 | | U |
| 100-41-4 | Ethylbenzene | 5000 | | U |
| 1330-20-7 | Xylene (para & meta) | 5000 | | U |
| 95-47-6 | Xylene (Ortho) | 5000 | | U |
| 100-42-5 | Styrene | 5000 | | U |
| 75-25-2 | Bromoform | 5000 | | U |
| 98-82-8 | Isopropylbenzene | 5000 | | U |
| 108-86-1 | Bromobenzene | 5000 | | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-5

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_5 0.010 ml
 Sample wt/vol: 0.010 (g/ml) ML Lab File ID: 7543_5.D
 Level: (low/med) LOW Date Received: 02/07/07
 % Moisture: not dec. _____ Date Analyzed: 02/08/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|-------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 5000 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 5000 | U |
| 103-65-1 | n-Propylbenzene | | 5000 | U |
| 95-49-8 | 2-Chlorotoluene | | 5000 | U |
| 106-43-4 | 4-Chlorotoluene | | 5000 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 5000 | U |
| 98-06-6 | tert-Butylbenzene | | 5000 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 5000 | U |
| 135-98-8 | sec-Butylbenzene | | 5000 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 5000 | U |
| 99-87-6 | 4-Isopropyltoluene | | 5000 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 5000 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 5000 | U |
| 104-51-8 | n-Butylbenzene | | 5000 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 5000 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 5000 | U |
| 87-68-3 | Hexachlorobutadiene | | 5000 | U |
| 91-20-3 | Naphthalene | | 5000 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 5000 | U |
| 1634-04-4 | MTBE | | 5000 | U |
| 67-64-1 | Acetone | | 10000 | U |
| 75-15-0 | Carbon disulfide | | 5000 | U |
| 78-93-3 | 2-Butanone | | 10000 | U |
| 109-99-9 | Tetrahydrofuran | | 10000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 10000 | U |
| 591-78-6 | 2-Hexanone | | 10000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 10000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-6

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA.
 Matrix: (soil/water) WATER Lab Sample ID: 7543_6 0.010 ml
 Sample wt/vol: 0.01 (g/ml) ML Lab File ID: 7543_6.D
 Level: (low/med) LOW Date Received: 02/07/07
 % Moisture: not dec. Date Analyzed: 02/09/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 5000 | U |
| 74-87-3 | Chloromethane | | 5000 | U |
| 75-01-4 | Vinyl chloride | | 5000 | U |
| 74-83-9 | Bromomethane | | 5000 | U |
| 75-00-3 | Chloroethane | | 5000 | U |
| 75-69-4 | Trichlorofluoromethane | | 5000 | U |
| 75-35-4 | 1,1-Dichloroethene | | 5000 | U |
| 75-09-2 | Methylene chloride | | 5000 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 5000 | U |
| 75-34-3 | 1,1-Dichloroethane | | 5000 | U |
| 594-20-7 | 2,2-Dichloropropane | | 5000 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 5000 | U |
| 74-97-5 | Bromochloromethane | | 5000 | U |
| 67-66-3 | Chloroform | | 5000 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 5000 | U |
| 56-23-5 | Carbon tetrachloride | | 5000 | U |
| 563-58-6 | 1,1-Dichloropropene | | 5000 | U |
| 71-43-2 | Benzene | | 5000 | U |
| 107-06-2 | 1,2-Dichloroethane | | 5000 | U |
| 79-01-6 | Trichloroethene | | 5000 | U |
| 78-87-5 | 1,2-Dichloropropane | | 5000 | U |
| 74-95-3 | Dibromomethane | | 5000 | U |
| 75-27-4 | Bromodichloromethane | | 5000 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 5000 | U |
| 108-88-3 | Toluene | | 5000 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 5000 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 5000 | U |
| 127-18-4 | Tetrachloroethene | | 17000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 5000 | U |
| 124-48-1 | Dibromochloromethane | | 5000 | U |
| 108-90-7 | Chlorobenzene | | 5000 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 5000 | U |
| 100-41-4 | Ethylbenzene | | 5000 | U |
| 1330-20-7 | Xylene (para & meta) | | 5000 | U |
| 95-47-6 | Xylene (Ortho) | | 5000 | U |
| 100-42-5 | Styrene | | 5000 | U |
| 75-25-2 | Bromoform | | 5000 | U |
| 98-82-8 | Isopropylbenzene | | 5000 | U |
| 108-86-1 | Bromobenzene | | 5000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-6

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_6 0.010 ml
 Sample wt/vol: 0.01 (g/ml) ML Lab File ID: 7543_6.D
 Level: (low/med) LOW Date Received: 02/07/07
 % Moisture: not dec. _____ Date Analyzed: 02/09/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5000 | U | |
| 96-18-4 | 1,2,3-Trichloropropane | 5000 | U | |
| 103-65-1 | n-Propylbenzene | 5000 | U | |
| 95-49-8 | 2-Chlorotoluene | 5000 | U | |
| 106-43-4 | 4-Chlorotoluene | 5000 | U | |
| 108-67-8 | 1,3,5-Trimethylbenzene | 5000 | U | |
| 98-06-6 | tert-Butylbenzene | 5000 | U | |
| 95-63-6 | 1,2,4-Trimethylbenzene | 5000 | U | |
| 135-98-8 | sec-Butylbenzene | 5000 | U | |
| 541-73-1 | 1,3-Dichlorobenzene | 5000 | U | |
| 99-87-6 | 4-Isopropyltoluene | 5000 | U | |
| 106-46-7 | 1,4-Dichlorobenzene | 5000 | U | |
| 95-50-1 | 1,2-Dichlorobenzene | 5000 | U | |
| 104-51-8 | n-Butylbenzene | 5000 | U | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | 5000 | U | |
| 120-82-1 | 1,2,4-Trichlorobenzene | 5000 | U | |
| 87-68-3 | Hexachlorobutadiene | 5000 | U | |
| 91-20-3 | Naphthalene | 5000 | U | |
| 87-61-6 | 1,2,3-Trichlorobenzene | 5000 | U | |
| 1634-04-4 | MTBE | 5000 | U | |
| 67-64-1 | Acetone | 10000 | U | |
| 75-15-0 | Carbon disulfide | 5000 | U | |
| 78-93-3 | 2-Butanone | 10000 | U | |
| 109-99-9 | Tetrahydrofuran | 10000 | U | |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 10000 | U | |
| 591-78-6 | 2-Hexanone | 10000 | U | |
| 110-75-8 | 2-Chloroethyl vinyl ether | 10000 | U | |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-7

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_7 0.010 ml
 Sample wt/vol: 0.01 (g/ml) ML Lab File ID: 7543_7.D
 Level: (low/med) LOW Date Received: 02/07/07
 % Moisture: not dec. _____ Date Analyzed: 02/09/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | 5000 | | U |
| 74-87-3 | Chloromethane | 5000 | | U |
| 75-01-4 | Vinyl chloride | 5000 | | U |
| 74-83-9 | Bromomethane | 5000 | | U |
| 75-00-3 | Chloroethane | 5000 | | U |
| 75-69-4 | Trichlorofluoromethane | 5000 | | U |
| 75-35-4 | 1,1-Dichloroethene | 5000 | | U |
| 75-09-2 | Methylene chloride | 5000 | | U |
| 156-60-5 | trans-1,2-Dichloroethene | 5000 | | U |
| 75-34-3 | 1,1-Dichloroethane | 5000 | | U |
| 594-20-7 | 2,2-Dichloropropane | 5000 | | U |
| 156-59-2 | cis-1,2-Dichloroethene | 5000 | | U |
| 74-97-5 | Bromochloromethane | 5000 | | U |
| 67-66-3 | Chloroform | 5000 | | U |
| 71-55-6 | 1,1,1-Trichloroethane | 5000 | | U |
| 56-23-5 | Carbon tetrachloride | 5000 | | U |
| 563-58-6 | 1,1-Dichloropropene | 5000 | | U |
| 71-43-2 | Benzene | 5000 | | U |
| 107-06-2 | 1,2-Dichloroethane | 5000 | | U |
| 79-01-6 | Trichloroethene | 5000 | | U |
| 78-87-5 | 1,2-Dichloropropane | 5000 | | U |
| 74-95-3 | Dibromomethane | 5000 | | U |
| 75-27-4 | Bromodichloromethane | 5000 | | U |
| 10061-02-5 | cis-1,3-Dichloropropene | 5000 | | U |
| 108-88-3 | Toluene | 5000 | | U |
| 10061-01-6 | trans-1,3-Dichloropropene | 5000 | | U |
| 79-00-5 | 1,1,2-Trichloroethane | 5000 | | U |
| 127-18-4 | Tetrachloroethene | 21000 | | D |
| 142-28-9 | 1,3-Dichloropropane | 5000 | | U |
| 124-48-1 | Dibromochloromethane | 5000 | | U |
| 108-90-7 | Chlorobenzene | 5000 | | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 5000 | | U |
| 100-41-4 | Ethylbenzene | 5000 | | U |
| 1330-20-7 | Xylene (para & meta) | 5000 | | U |
| 95-47-6 | Xylene (Ortho) | 5000 | | U |
| 100-42-5 | Styrene | 5000 | | U |
| 75-25-2 | Bromoform | 5000 | | U |
| 98-82-8 | Isopropylbenzene | 5000 | | U |
| 108-86-1 | Bromobenzene | 5000 | | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-7

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_7 0.010 ml
 Sample wt/vol: 0.01 (g/ml) ML Lab File ID: 7543_7.D
 Level: (low/med) LOW Date Received: 02/07/07
 % Moisture: not dec. _____ Date Analyzed: 02/09/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 5000 | U | |
| 96-18-4 | 1,2,3-Trichloropropane | 5000 | U | |
| 103-65-1 | n-Propylbenzene | 5000 | U | |
| 95-49-8 | 2-Chlorotoluene | 5000 | U | |
| 106-43-4 | 4-Chlorotoluene | 5000 | U | |
| 108-67-8 | 1,3,5-Trimethylbenzene | 5000 | U | |
| 98-06-6 | tert-Butylbenzene | 5000 | U | |
| 95-63-6 | 1,2,4-Trimethylbenzene | 5000 | U | |
| 135-98-8 | sec-Butylbenzene | 5000 | U | |
| 541-73-1 | 1,3-Dichlorobenzene | 5000 | U | |
| 99-87-6 | 4-Isopropyltoluene | 5000 | U | |
| 106-46-7 | 1,4-Dichlorobenzene | 5000 | U | |
| 95-50-1 | 1,2-Dichlorobenzene | 5000 | U | |
| 104-51-8 | n-Butylbenzene | 5000 | U | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | 5000 | U | |
| 120-82-1 | 1,2,4-Trichlorobenzene | 5000 | U | |
| 87-68-3 | Hexachlorobutadiene | 5000 | U | |
| 91-20-3 | Naphthalene | 5000 | U | |
| 87-61-6 | 1,2,3-Trichlorobenzene | 5000 | U | |
| 1634-04-4 | MTBE | 5000 | U | |
| 67-64-1 | Acetone | 10000 | U | |
| 75-15-0 | Carbon disulfide | 5000 | U | |
| 78-93-3 | 2-Butanone | 10000 | U | |
| 109-99-9 | Tetrahydrofuran | 10000 | U | |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | 10000 | U | |
| 591-78-6 | 2-Hexanone | 10000 | U | |
| 110-75-8 | 2-Chloroethyl vinyl ether | 10000 | U | |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-08

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_08 0.010 ml
 Sample wt/vol: 0.010 (g/ml) ML Lab File ID: 7543_08.D
 Level: (low/med) LOW Date Received: 02/08/07
 % Moisture: not dec. _____ Date Analyzed: 02/14/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | 5000 | | U |
| 74-87-3 | Chloromethane | 5000 | | U |
| 75-01-4 | Vinyl chloride | 5000 | | U |
| 74-83-9 | Bromomethane | 5000 | | U |
| 75-00-3 | Chloroethane | 5000 | | U |
| 75-69-4 | Trichlorofluoromethane | 5000 | | U |
| 75-35-4 | 1,1-Dichloroethene | 5000 | | U |
| 75-09-2 | Methylene chloride | 5000 | | U |
| 156-60-5 | trans-1,2-Dichloroethene | 5000 | | U |
| 75-34-3 | 1,1-Dichloroethane | 5000 | | U |
| 594-20-7 | 2,2-Dichloropropane | 5000 | | U |
| 156-59-2 | cis-1,2-Dichloroethene | 5000 | | U |
| 74-97-5 | Bromochloromethane | 5000 | | U |
| 67-66-3 | Chloroform | 5000 | | U |
| 71-55-6 | 1,1,1-Trichloroethane | 5000 | | U |
| 56-23-5 | Carbon tetrachloride | 5000 | | U |
| 563-58-6 | 1,1-Dichloropropene | 5000 | | U |
| 71-43-2 | Benzene | 5000 | | U |
| 107-06-2 | 1,2-Dichloroethane | 5000 | | U |
| 79-01-6 | Trichloroethene | 5000 | | U |
| 78-87-5 | 1,2-Dichloropropane | 5000 | | U |
| 74-95-3 | Dibromomethane | 5000 | | U |
| 75-27-4 | Bromodichloromethane | 5000 | | U |
| 10061-02-5 | cis-1,3-Dichloropropene | 5000 | | U |
| 108-88-3 | Toluene | 5000 | | U |
| 10061-01-6 | trans-1,3-Dichloropropene | 5000 | | U |
| 79-00-5 | 1,1,2-Trichloroethane | 5000 | | U |
| 127-18-4 | Tetrachloroethene | 26000 | | D |
| 142-28-9 | 1,3-Dichloropropane | 5000 | | U |
| 124-48-1 | Dibromochloromethane | 5000 | | U |
| 108-90-7 | Chlorobenzene | 5000 | | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 5000 | | U |
| 100-41-4 | Ethylbenzene | 5000 | | U |
| 1330-20-7 | Xylene (para & meta) | 5000 | | U |
| 95-47-6 | Xylene (Ortho) | 5000 | | U |
| 100-42-5 | Styrene | 5000 | | U |
| 75-25-2 | Bromoform | 5000 | | U |
| 98-82-8 | Isopropylbenzene | 5000 | | U |
| 108-86-1 | Bromobenzene | 5000 | | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-08

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_08 0.010 ml
 Sample wt/vol: 0.010 (g/ml) ML Lab File ID: 7543_08.D
 Level: (low/med) LOW Date Received: 02/08/07
 % Moisture: not dec. _____ Date Analyzed: 02/14/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|-------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 5000 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 5000 | U |
| 103-65-1 | n-Propylbenzene | | 5000 | U |
| 95-49-8 | 2-Chlorotoluene | | 5000 | U |
| 106-43-4 | 4-Chlorotoluene | | 5000 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 5000 | U |
| 98-06-6 | tert-Butylbenzene | | 5000 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 5000 | U |
| 135-98-8 | sec-Butylbenzene | | 5000 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 5000 | U |
| 99-87-6 | 4-Isopropyltoluene | | 5000 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 5000 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 5000 | U |
| 104-51-8 | n-Butylbenzene | | 5000 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 5000 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 5000 | U |
| 87-68-3 | Hexachlorobutadiene | | 5000 | U |
| 91-20-3 | Naphthalene | | 5000 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 5000 | U |
| 1634-04-4 | MTBE | | 5000 | U |
| 67-64-1 | Acetone | | 10000 | U |
| 75-15-0 | Carbon disulfide | | 5000 | U |
| 78-93-3 | 2-Butanone | | 10000 | U |
| 109-99-9 | Tetrahydrofuran | | 10000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 10000 | U |
| 591-78-6 | 2-Hexanone | | 10000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 10000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-09

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_09 0.010 ml
 Sample wt/vol: 0.0|0 (g/ml) ML Lab File ID: 7543_09.D
 Level: (low/med) LOW Date Received: 02/08/07
 % Moisture: not dec. _____ Date Analyzed: 02/14/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 5000 | U |
| 74-87-3 | Chloromethane | | 5000 | U |
| 75-01-4 | Vinyl chloride | | 5000 | U |
| 74-83-9 | Bromomethane | | 5000 | U |
| 75-00-3 | Chloroethane | | 5000 | U |
| 75-69-4 | Trichlorofluoromethane | | 5000 | U |
| 75-35-4 | 1,1-Dichloroethene | | 5000 | U |
| 75-09-2 | Methylene chloride | | 5000 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 5000 | U |
| 75-34-3 | 1,1-Dichloroethane | | 5000 | U |
| 594-20-7 | 2,2-Dichloropropane | | 5000 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 5000 | U |
| 74-97-5 | Bromochloromethane | | 5000 | U |
| 67-66-3 | Chloroform | | 5000 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 5000 | U |
| 56-23-5 | Carbon tetrachloride | | 5000 | U |
| 563-58-6 | 1,1-Dichloropropene | | 5000 | U |
| 71-43-2 | Benzene | | 5000 | U |
| 107-06-2 | 1,2-Dichloroethane | | 5000 | U |
| 79-01-6 | Trichloroethene | | 5000 | U |
| 78-87-5 | 1,2-Dichloropropane | | 5000 | U |
| 74-95-3 | Dibromomethane | | 5000 | U |
| 75-27-4 | Bromodichloromethane | | 5000 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 5000 | U |
| 108-88-3 | Toluene | | 5000 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 5000 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 5000 | U |
| 127-18-4 | Tetrachloroethene | | 21000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 5000 | U |
| 124-48-1 | Dibromochloromethane | | 5000 | U |
| 108-90-7 | Chlorobenzene | | 5000 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 5000 | U |
| 100-41-4 | Ethylbenzene | | 5000 | U |
| 1330-20-7 | Xylene (para & meta) | | 5000 | U |
| 95-47-6 | Xylene (Ortho) | | 5000 | U |
| 100-42-5 | Styrene | | 5000 | U |
| 75-25-2 | Bromoform | | 5000 | U |
| 98-82-8 | Isopropylbenzene | | 5000 | U |
| 108-86-1 | Bromobenzene | | 5000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-09

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_09 0.010 ml
 Sample wt/vol: 0.010 (g/ml) ML Lab File ID: 7543_09.D
 Level: (low/med) LOW Date Received: 02/08/07
 % Moisture: not dec. _____ Date Analyzed: 02/14/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|-------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 5000 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 5000 | U |
| 103-65-1 | n-Propylbenzene | | 5000 | U |
| 95-49-8 | 2-Chlorotoluene | | 5000 | U |
| 106-43-4 | 4-Chlorotoluene | | 5000 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 5000 | U |
| 98-06-6 | tert-Butylbenzene | | 5000 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 5000 | U |
| 135-98-8 | sec-Butylbenzene | | 5000 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 5000 | U |
| 99-87-6 | 4-Isopropyltoluene | | 5000 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 5000 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 5000 | U |
| 104-51-8 | n-Butylbenzene | | 5000 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 5000 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 5000 | U |
| 87-68-3 | Hexachlorobutadiene | | 5000 | U |
| 91-20-3 | Naphthalene | | 5000 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 5000 | U |
| 1634-04-4 | MTBE | | 5000 | U |
| 67-64-1 | Acetone | | 10000 | U |
| 75-15-0 | Carbon disulfide | | 5000 | U |
| 78-93-3 | 2-Butanone | | 10000 | U |
| 109-99-9 | Tetrahydrofuran | | 10000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 10000 | U |
| 591-78-6 | 2-Hexanone | | 10000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 10000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-10

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_10 0.010 ml
 Sample wt/vol: 0.010 (g/ml) ML Lab File ID: 7543_10.D
 Level: (low/med) LOW Date Received: 02/08/07
 % Moisture: not dec. _____ Date Analyzed: 02/14/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|-------|---|
| 75-71-8 | Dichlorodifluoromethane | | 5000 | U |
| 74-87-3 | Chloromethane | | 5000 | U |
| 75-01-4 | Vinyl chloride | | 5000 | U |
| 74-83-9 | Bromomethane | | 5000 | U |
| 75-00-3 | Chloroethane | | 5000 | U |
| 75-69-4 | Trichlorofluoromethane | | 5000 | U |
| 75-35-4 | 1,1-Dichloroethene | | 5000 | U |
| 75-09-2 | Methylene chloride | | 5000 | U |
| 156-60-5 | trans-1,2-Dichloroethene | | 5000 | U |
| 75-34-3 | 1,1-Dichloroethane | | 5000 | U |
| 594-20-7 | 2,2-Dichloropropane | | 5000 | U |
| 156-59-2 | cis-1,2-Dichloroethene | | 5000 | U |
| 74-97-5 | Bromochloromethane | | 5000 | U |
| 67-66-3 | Chloroform | | 5000 | U |
| 71-55-6 | 1,1,1-Trichloroethane | | 5000 | U |
| 56-23-5 | Carbon tetrachloride | | 5000 | U |
| 563-58-6 | 1,1-Dichloropropene | | 5000 | U |
| 71-43-2 | Benzene | | 5000 | U |
| 107-06-2 | 1,2-Dichloroethane | | 5000 | U |
| 79-01-6 | Trichloroethene | | 5000 | U |
| 78-87-5 | 1,2-Dichloropropane | | 5000 | U |
| 74-95-3 | Dibromomethane | | 5000 | U |
| 75-27-4 | Bromodichloromethane | | 5000 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | | 5000 | U |
| 108-88-3 | Toluene | | 5000 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | | 5000 | U |
| 79-00-5 | 1,1,2-Trichloroethane | | 5000 | U |
| 127-18-4 | Tetrachloroethene | | 14000 | D |
| 142-28-9 | 1,3-Dichloropropane | | 5000 | U |
| 124-48-1 | Dibromochloromethane | | 5000 | U |
| 108-90-7 | Chlorobenzene | | 5000 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | | 5000 | U |
| 100-41-4 | Ethylbenzene | | 5000 | U |
| 1330-20-7 | Xylene (para & meta) | | 5000 | U |
| 95-47-6 | Xylene (Ortho) | | 5000 | U |
| 100-42-5 | Styrene | | 5000 | U |
| 75-25-2 | Bromoform | | 5000 | U |
| 98-82-8 | Isopropylbenzene | | 5000 | U |
| 108-86-1 | Bromobenzene | | 5000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-10

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_10 0.010 ml
 Sample wt/vol: 0.010 (g/ml) ML Lab File ID: 7543_10.D
 Level: (low/med) LOW Date Received: 02/08/07
 % Moisture: not dec. _____ Date Analyzed: 02/14/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|-------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 5000 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 5000 | U |
| 103-65-1 | n-Propylbenzene | | 5000 | U |
| 95-49-8 | 2-Chlorotoluene | | 5000 | U |
| 106-43-4 | 4-Chlorotoluene | | 5000 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 5000 | U |
| 98-06-6 | tert-Butylbenzene | | 5000 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 5000 | U |
| 135-98-8 | sec-Butylbenzene | | 5000 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 5000 | U |
| 99-87-6 | 4-Isopropyltoluene | | 5000 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 5000 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 5000 | U |
| 104-51-8 | n-Butylbenzene | | 5000 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 5000 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 5000 | U |
| 87-68-3 | Hexachlorobutadiene | | 5000 | U |
| 91-20-3 | Naphthalene | | 5000 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 5000 | U |
| 1634-04-4 | MTBE | | 5000 | U |
| 67-64-1 | Acetone | | 10000 | U |
| 75-15-0 | Carbon disulfide | | 5000 | U |
| 78-93-3 | 2-Butanone | | 10000 | U |
| 109-99-9 | Tetrahydrofuran | | 10000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 10000 | U |
| 591-78-6 | 2-Hexanone | | 10000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 10000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-11

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_11 0.010 ml
 Sample wt/vol: 0.010 (g/ml) ML Lab File ID: 7543_11.D
 Level: (low/med) LOW Date Received: 02/08/07
 % Moisture: not dec. _____ Date Analyzed: 02/14/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|------------|---------------------------|-----------------|------|---|
| 75-71-8 | Dichlorodifluoromethane | 5000 | U | |
| 74-87-3 | Chloromethane | 5000 | U | |
| 75-01-4 | Vinyl chloride | 5000 | U | |
| 74-83-9 | Bromomethane | 5000 | U | |
| 75-00-3 | Chloroethane | 5000 | U | |
| 75-69-4 | Trichlorofluoromethane | 5000 | U | |
| 75-35-4 | 1,1-Dichloroethene | 5000 | U | |
| 75-09-2 | Methylene chloride | 5000 | U | |
| 156-60-5 | trans-1,2-Dichloroethene | 5000 | U | |
| 75-34-3 | 1,1-Dichloroethane | 5000 | U | |
| 594-20-7 | 2,2-Dichloropropane | 5000 | U | |
| 156-59-2 | cis-1,2-Dichloroethene | 5000 | U | |
| 74-97-5 | Bromochloromethane | 5000 | U | |
| 67-66-3 | Chloroform | 5000 | U | |
| 71-55-6 | 1,1,1-Trichloroethane | 5000 | U | |
| 56-23-5 | Carbon tetrachloride | 5000 | U | |
| 563-58-6 | 1,1-Dichloropropene | 5000 | U | |
| 71-43-2 | Benzene | 5000 | U | |
| 107-06-2 | 1,2-Dichloroethane | 5000 | U | |
| 79-01-6 | Trichloroethene | 5000 | U | |
| 78-87-5 | 1,2-Dichloropropane | 5000 | U | |
| 74-95-3 | Dibromomethane | 5000 | U | |
| 75-27-4 | Bromodichloromethane | 5000 | U | |
| 10061-02-5 | cis-1,3-Dichloropropene | 5000 | U | |
| 108-88-3 | Toluene | 5000 | U | |
| 10061-01-6 | trans-1,3-Dichloropropene | 5000 | U | |
| 79-00-5 | 1,1,2-Trichloroethane | 5000 | U | |
| 127-18-4 | Tetrachloroethene | 24000 | D | |
| 142-28-9 | 1,3-Dichloropropane | 5000 | U | |
| 124-48-1 | Dibromochloromethane | 5000 | U | |
| 108-90-7 | Chlorobenzene | 5000 | U | |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 5000 | U | |
| 100-41-4 | Ethylbenzene | 5000 | U | |
| 1330-20-7 | Xylene (para & meta) | 5000 | U | |
| 95-47-6 | Xylene (Ortho) | 5000 | U | |
| 100-42-5 | Styrene | 5000 | U | |
| 75-25-2 | Bromoform | 5000 | U | |
| 98-82-8 | Isopropylbenzene | 5000 | U | |
| 108-86-1 | Bromobenzene | 5000 | U | |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-11

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_11 0.010 ml
 Sample wt/vol: 0.010 (g/ml) ML Lab File ID: 7543_11.D
 Level: (low/med) LOW Date Received: 02/08/07
 % Moisture: not dec. _____ Date Analyzed: 02/14/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|-------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 5000 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 5000 | U |
| 103-65-1 | n-Propylbenzene | | 5000 | U |
| 95-49-8 | 2-Chlorotoluene | | 5000 | U |
| 106-43-4 | 4-Chlorotoluene | | 5000 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 5000 | U |
| 98-06-6 | tert-Butylbenzene | | 5000 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 5000 | U |
| 135-98-8 | sec-Butylbenzene | | 5000 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 5000 | U |
| 99-87-6 | 4-Isopropyltoluene | | 5000 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 5000 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 5000 | U |
| 104-51-8 | n-Butylbenzene | | 5000 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 5000 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 5000 | U |
| 87-68-3 | Hexachlorobutadiene | | 5000 | U |
| 91-20-3 | Naphthalene | | 5000 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 5000 | U |
| 1634-04-4 | MTBE | | 5000 | U |
| 67-64-1 | Acetone | | 10000 | U |
| 75-15-0 | Carbon disulfide | | 5000 | U |
| 78-93-3 | 2-Butanone | | 10000 | U |
| 109-99-9 | Tetrahydrofuran | | 10000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 10000 | U |
| 591-78-6 | 2-Hexanone | | 10000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 10000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-12

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_12 0.010 ml
 Sample wt/vol: 0.01⁰ (g/ml) ML Lab File ID: 7543_12.D
 Level: (low/med) LOW Date Received: 02/08/07
 % Moisture: not dec. _____ Date Analyzed: 02/14/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

CAS NO. COMPOUND (ug/L or ug/Kg) UG/L Q

| | | | |
|------------|---------------------------|------|----|
| 75-71-8 | Dichlorodifluoromethane | 5000 | U |
| 74-87-3 | Chloromethane | 5000 | U |
| 75-01-4 | Vinyl chloride | 5000 | U |
| 74-83-9 | Bromomethane | 5000 | U |
| 75-00-3 | Chloroethane | 5000 | U |
| 75-69-4 | Trichlorofluoromethane | 5000 | U |
| 75-35-4 | 1,1-Dichloroethene | 5000 | U |
| 75-09-2 | Methylene chloride | 5000 | U |
| 156-60-5 | trans-1,2-Dichloroethene | 5000 | U |
| 75-34-3 | 1,1-Dichloroethane | 5000 | U |
| 594-20-7 | 2,2-Dichloropropane | 5000 | U |
| 156-59-2 | cis-1,2-Dichloroethene | 5000 | U |
| 74-97-5 | Bromochloromethane | 5000 | U |
| 67-66-3 | Chloroform | 5000 | U |
| 71-55-6 | 1,1,1-Trichloroethane | 5000 | U |
| 56-23-5 | Carbon tetrachloride | 5000 | U |
| 563-58-6 | 1,1-Dichloropropene | 5000 | U |
| 71-43-2 | Benzene | 5000 | U |
| 107-06-2 | 1,2-Dichloroethane | 5000 | U |
| 79-01-6 | Trichloroethene | 5000 | U |
| 78-87-5 | 1,2-Dichloropropane | 5000 | U |
| 74-95-3 | Dibromomethane | 5000 | U |
| 75-27-4 | Bromodichloromethane | 5000 | U |
| 10061-02-5 | cis-1,3-Dichloropropene | 5000 | U |
| 108-88-3 | Toluene | 5000 | U |
| 10061-01-6 | trans-1,3-Dichloropropene | 5000 | U |
| 79-00-5 | 1,1,2-Trichloroethane | 5000 | U |
| 127-18-4 | Tetrachloroethene | 2100 | JD |
| 142-28-9 | 1,3-Dichloropropane | 5000 | U |
| 124-48-1 | Dibromochloromethane | 5000 | U |
| 108-90-7 | Chlorobenzene | 5000 | U |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 5000 | U |
| 100-41-4 | Ethylbenzene | 5000 | U |
| 1330-20-7 | Xylene (para & meta) | 5000 | U |
| 95-47-6 | Xylene (Ortho) | 5000 | U |
| 100-42-5 | Styrene | 5000 | U |
| 75-25-2 | Bromoform | 5000 | U |
| 98-82-8 | Isopropylbenzene | 5000 | U |
| 108-86-1 | Bromobenzene | 5000 | U |

1A
VOLATILE ORGANICS ANALYSIS DATA SHEET

SAMPLE:

7543-12

Lab Name: ATL Analyzed: Tony Soto
 N.J. Dep # 11001 Cal Date: 01/12/07 Gc/Ms: 3 Client: NA
 Matrix: (soil/water) WATER Lab Sample ID: 7543_12 0.010 ml
 Sample wt/vol: 0.010 (g/ml) ML Lab File ID: 7543_12.D
 Level: (low/med) LOW Date Received: 02/08/07
 % Moisture: not dec. _____ Date Analyzed: 02/14/07
 GC Column: RTX 502 ID: 0.25 (mm) Dilution Factor: 1000.0
 Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)

CONCENTRATION UNITS:

| CAS NO. | COMPOUND | (ug/L or ug/Kg) | UG/L | Q |
|-----------|-----------------------------|-----------------|-------|---|
| 79-34-5 | 1,1,2,2-Tetrachloroethane | | 5000 | U |
| 96-18-4 | 1,2,3-Trichloropropane | | 5000 | U |
| 103-65-1 | n-Propylbenzene | | 5000 | U |
| 95-49-8 | 2-Chlorotoluene | | 5000 | U |
| 106-43-4 | 4-Chlorotoluene | | 5000 | U |
| 108-67-8 | 1,3,5-Trimethylbenzene | | 5000 | U |
| 98-06-6 | tert-Butylbenzene | | 5000 | U |
| 95-63-6 | 1,2,4-Trimethylbenzene | | 5000 | U |
| 135-98-8 | sec-Butylbenzene | | 5000 | U |
| 541-73-1 | 1,3-Dichlorobenzene | | 5000 | U |
| 99-87-6 | 4-Isopropyltoluene | | 5000 | U |
| 106-46-7 | 1,4-Dichlorobenzene | | 5000 | U |
| 95-50-1 | 1,2-Dichlorobenzene | | 5000 | U |
| 104-51-8 | n-Butylbenzene | | 5000 | U |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | | 5000 | U |
| 120-82-1 | 1,2,4-Trichlorobenzene | | 5000 | U |
| 87-68-3 | Hexachlorobutadiene | | 5000 | U |
| 91-20-3 | Naphthalene | | 5000 | U |
| 87-61-6 | 1,2,3-Trichlorobenzene | | 5000 | U |
| 1634-04-4 | MTBE | | 5000 | U |
| 67-64-1 | Acetone | | 10000 | U |
| 75-15-0 | Carbon disulfide | | 5000 | U |
| 78-93-3 | 2-Butanone | | 10000 | U |
| 109-99-9 | Tetrahydrofuran | | 10000 | U |
| 108-10-1 | 4-Methyl-2-pentanone (MIBK) | | 10000 | U |
| 591-78-6 | 2-Hexanone | | 10000 | U |
| 110-75-8 | 2-Chloroethyl vinyl ether | | 10000 | U |

3.0 Raw Data and Chromatograms

Raw data and Chromatograms are attached.

Data File : C:\HPCHEM\1\DATA\020807\7543_1.D

Acq On : 8 Feb 2007 7:50 pm

Sample : 7543_1 0.010 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Feb 13 15:36 2007

Vial: 6

Operator: RR/AS

Inst : GC/MS Ins

Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:51:14 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.73 | 96 | 237928 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 20.12 | 117 | 201829 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.35 | 152 | 104901 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|------|
| 16) Dibromofluoromethane | 13.29 | 113 | 67536 | 26.15 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 104.60% | |
| 19) 1,2-Dichloroethane-d4 | 14.24 | 65 | 62081 | 23.73 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 94.92% | |
| 28) Toluene-d8 | 17.42 | 98 | 264441 | 24.61 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 98.44% | |
| 45) 4-Bromofluorobenzene | 22.27 | 95 | 71243 | 20.67 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 82.68% | |

Target Compounds

| | | | | | Qvalue |
|-------------------------------|------|-----|---|------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | |

(#) = qualifier out of range (m) = manual integration

7543_1.D 01_12_07.M

Wed Feb 14 16:08:58 2007

Page 1

Acq On : 8 Feb 2007 7:50 pm

Operator: RR/AS

Sample : 7543_1 0.010 ml

Inst : GC/MS Ins

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Feb 13 15:36 2007

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:51:14 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 122197 | 26.29 | ug/L | 99 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 20.14 | 112 | 5779 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | d | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | d | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

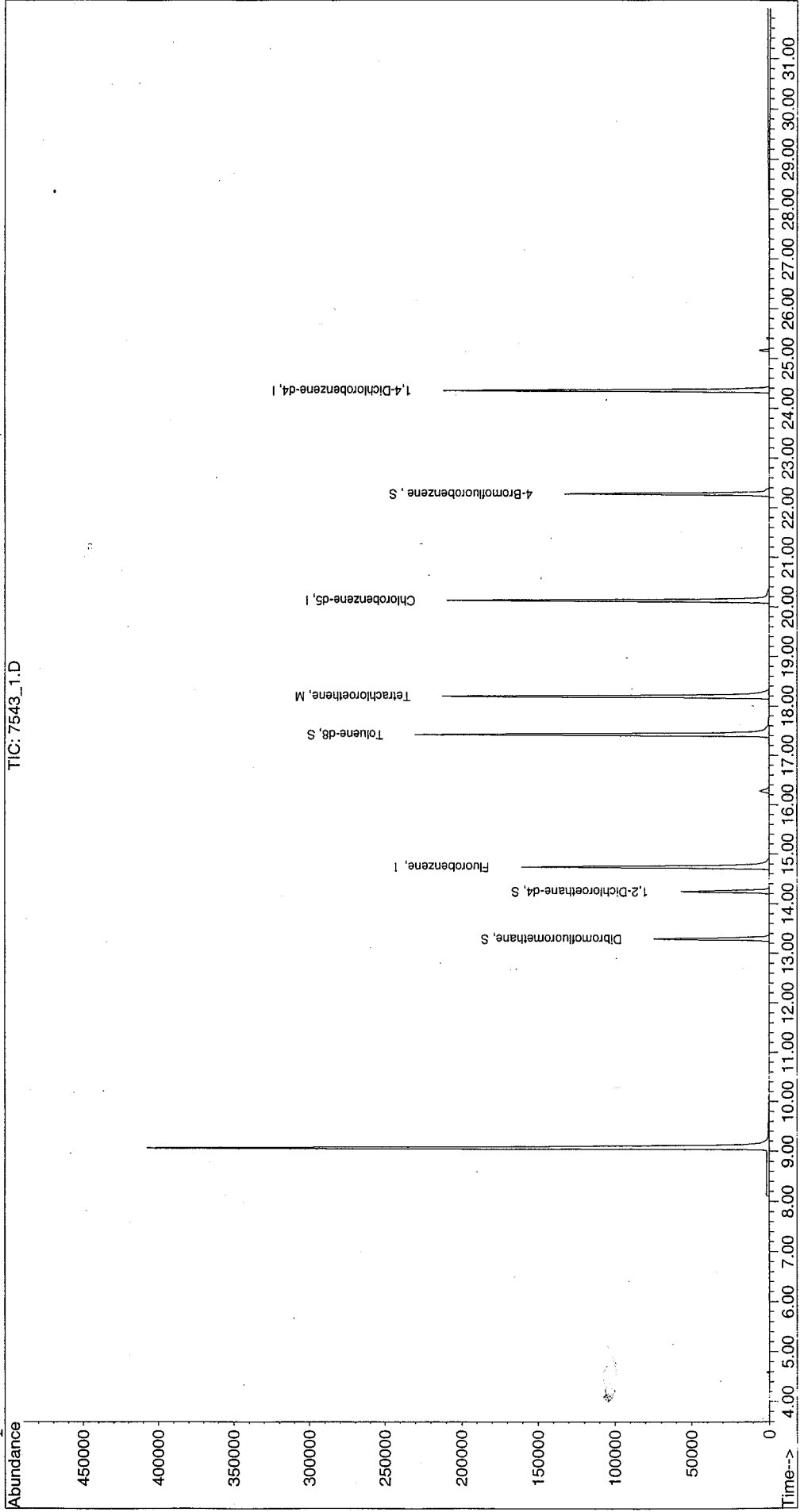
7543_1.D 01_12_07.M Wed Feb 14 16:08:59 2007

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Quantitation Report

Data File : C:\HPCHEM\1\DATA\020807\7543_1.D Vial: 6
Acq On : 8 Feb 2007 7:50 pm Operator: RR/AS
Sample : 7543_1 0.010 ml Inst : GC/MS Ins
Misc : Multiplr: 1.00
MS Integration Params: ODD.P
Quant Time: Feb 13 15:36 2007 Quant Results File: 01_12_07.RES

Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:54:07 2007
Response via : Initial Calibration



Acq On : 8 Feb 2007 9:44 pm

Operator: RR/AS

Sample : 7543_2 0.010 ml

Inst : GC/MS Ins

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Feb 13 15:37 2007

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:51:14 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|--------------------|------|------|----------|------|-------|----------|
|--------------------|------|------|----------|------|-------|----------|

| | | | | | | |
|----------------------------|-------|-----|--------|-------|------|------|
| 1) Fluorobenzene | 14.73 | 96 | 300813 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 20.12 | 117 | 244768 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.35 | 152 | 132175 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|--------|------|
| 16) Dibromofluoromethane | 13.29 | 113 | 74554 | 22.84 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 91.36% | |
| 19) 1,2-Dichloroethane-d4 | 14.24 | 65 | 67997 | 20.56 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 82.24% | |
| 28) Toluene-d8 | 17.42 | 98 | 314405 | 23.14 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 92.56% | |
| 45) 4-Bromofluorobenzene | 22.27 | 95 | 90857 | 21.74 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 86.96% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | d | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#)=qualifier out of range (m)=manual integration

7543_2.D 01_12_07.M

Wed Feb 14 16:09:32 2007

Data File : C:\HPCHEM\1\DATA\020807\7543_2.D

Vial: 9

Acq On : 8 Feb 2007 9:44 pm

Operator: RR/AS

Sample : 7543_2 0.010 ml

Inst : GC/MS Ins

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Feb 13 15:37 2007

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:51:14 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 143105 | 24.35 | ug/L | 99 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 20.15 | 112 | 7011 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | d | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | d | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7543_2.D 01_12_07.M

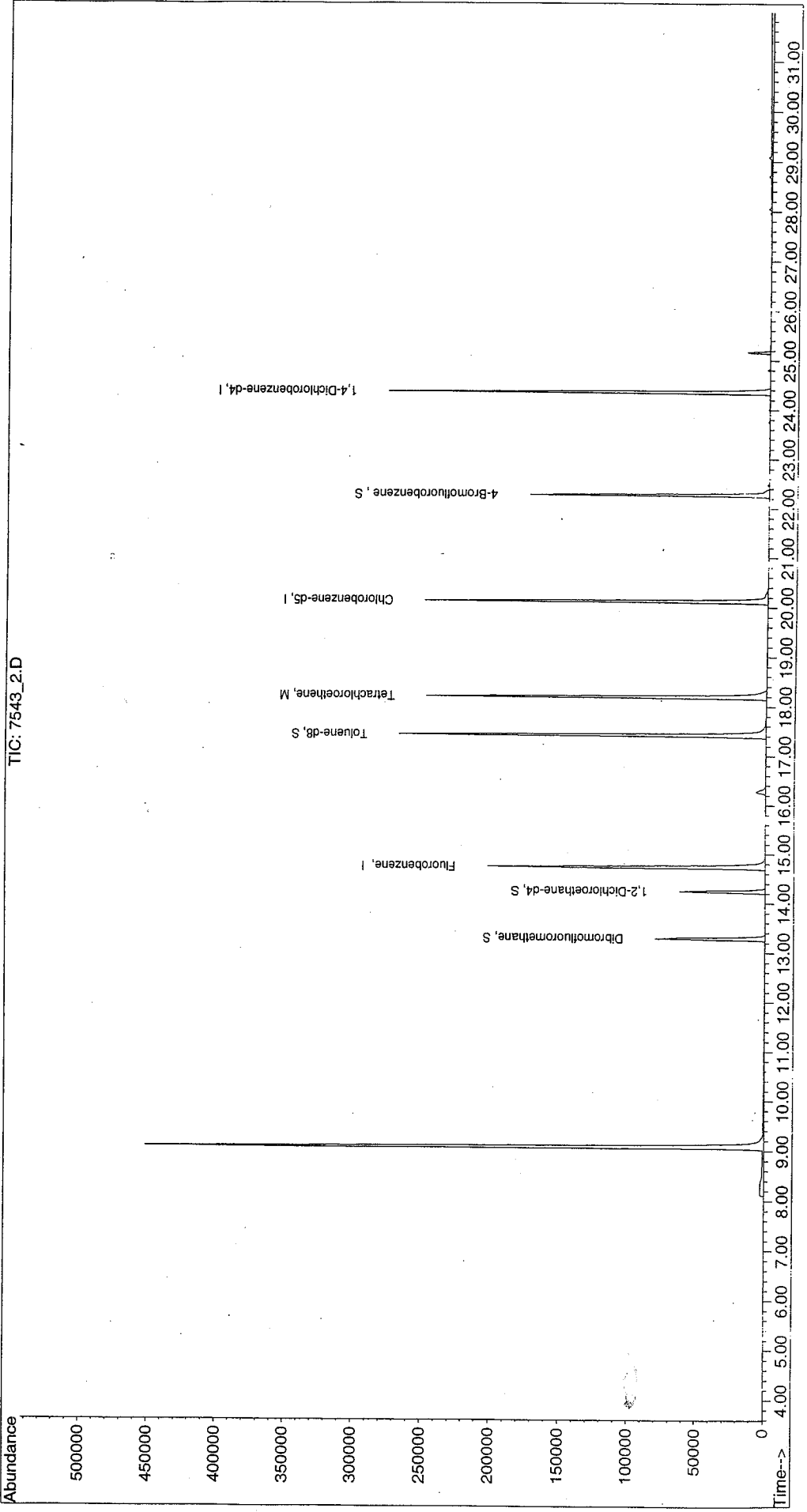
Wed Feb 14 16:09:34 2007

Page 2

Quantification Report

Data File : C:\HPCHEM\1\DATA\020807\7543_2.D
Acq On : 8 Feb 2007 9:44 pm
Sample : 7543_2 0.010 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Feb 13 15:37 2007
Vial: 9
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: 01_12_07.RES

Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:54:07 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\020807\7543_3.D
 Acq On : 8 Feb 2007 10:22 pm
 Sample : 7543_3 0.010 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Feb 13 15:37 2007

Vial: 10
 Operator: RR/AS
 Inst : GC/MS Ins
 Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Mon Jan 15 14:51:14 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|--------|------|----------|-------|--------|----------|
| 1) Fluorobenzene | 14.73 | 96 | 285444 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 20.12 | 117 | 235106 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.35 | 152 | 120866 | 25.00 | ug/L | 0.00 |
| System Monitoring Compounds | | | | | | |
| 16) Dibromofluoromethane | 13.28 | 113 | 73008 | 23.57 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 94.28% | |
| 19) 1,2-Dichloroethane-d4 | 14.24 | 65 | 67010 | 21.35 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 85.40% | |
| 28) Toluene-d8 | 17.41 | 98 | 303498 | 23.54 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 94.16% | |
| 45) 4-Bromofluorobenzene | 22.27 | 95 | 85072 | 21.19 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 84.76% | |
| Target Compounds | | | | | | |
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | Qvalue |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration
 7543_3.D 01_12_07.M Wed Feb 14 16:09:41 2007

Data File : C:\HPCHEM\1\DATA\020807\7543_3.D

Acq On : 8 Feb 2007 10:22 pm

Sample : 7543_3 0.010 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Feb 13 15:37 2007

Vial: 10

Operator: RR/AS

Inst : GC/MS Ins

Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:51:14 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 117455 | 21.06 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 20.15 | 112 | 5829 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | d | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | d | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

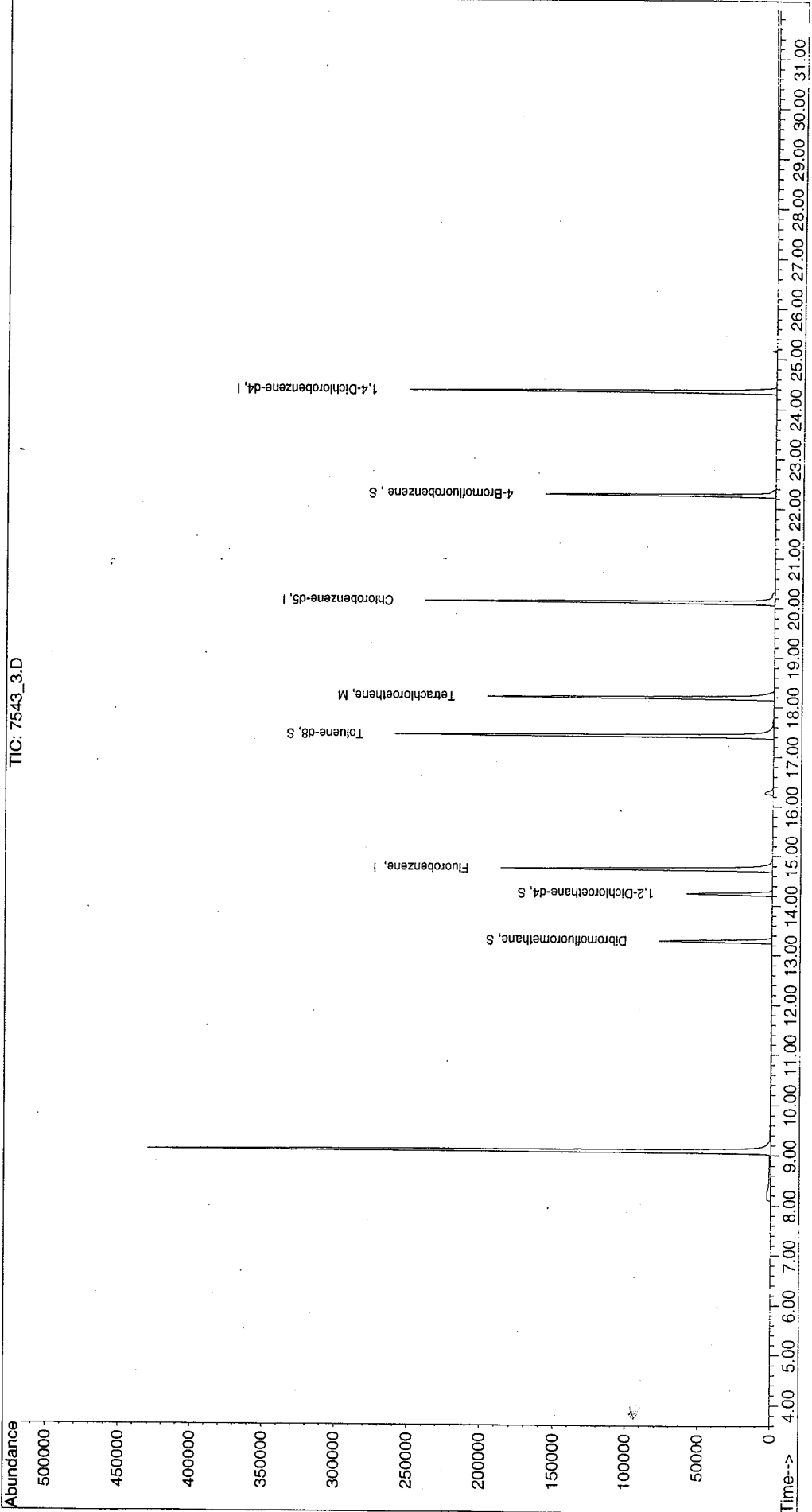
7543_3.D 01_12_07.M

Wed Feb 14 16:09:43 2007

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Quantitation report

Data File : C:\HPCHEM\1\DATA\020807\7543_3.D
Acq On : 8 Feb 2007 10:22 pm
Sample : 7543_3 0.010 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Feb 13 15:37 2007
Quant Results File: 01_12_07.RES
Vial: 10
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:54:07 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\020807\7543_4.D
 Acq On : 8 Feb 2007 11:00 pm
 Sample : 7543_4 0.010 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Feb 13 15:38 2007

Vial: 11
 Operator: RR/AS
 Inst : GC/MS Ins
 Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Mon Jan 15 14:51:14 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.73 | 96 | 268258 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 20.12 | 117 | 221481 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.35 | 152 | 115372 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|--------|------|
| 16) Dibromofluoromethane | 13.28 | 113 | 68632 | 23.57 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 94.28% | |
| 19) 1,2-Dichloroethane-d4 | 14.24 | 65 | 63402 | 21.49 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 85.96% | |
| 28) Toluene-d8 | 17.42 | 98 | 279734 | 23.09 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 92.36% | |
| 45) 4-Bromofluorobenzene | 22.27 | 95 | 79864 | 21.12 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 84.48% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\020807\7543_4.D
Acq On : 8 Feb 2007 11:00 pm
Sample : 7543_4 0.010 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Feb 13 15:38 2007

Vial: 11
Operator: RR/AS
Inst : GC/MS Ins.
Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:51:14 2007
Response via : Initial Calibration
DataAcq Meth : RUN

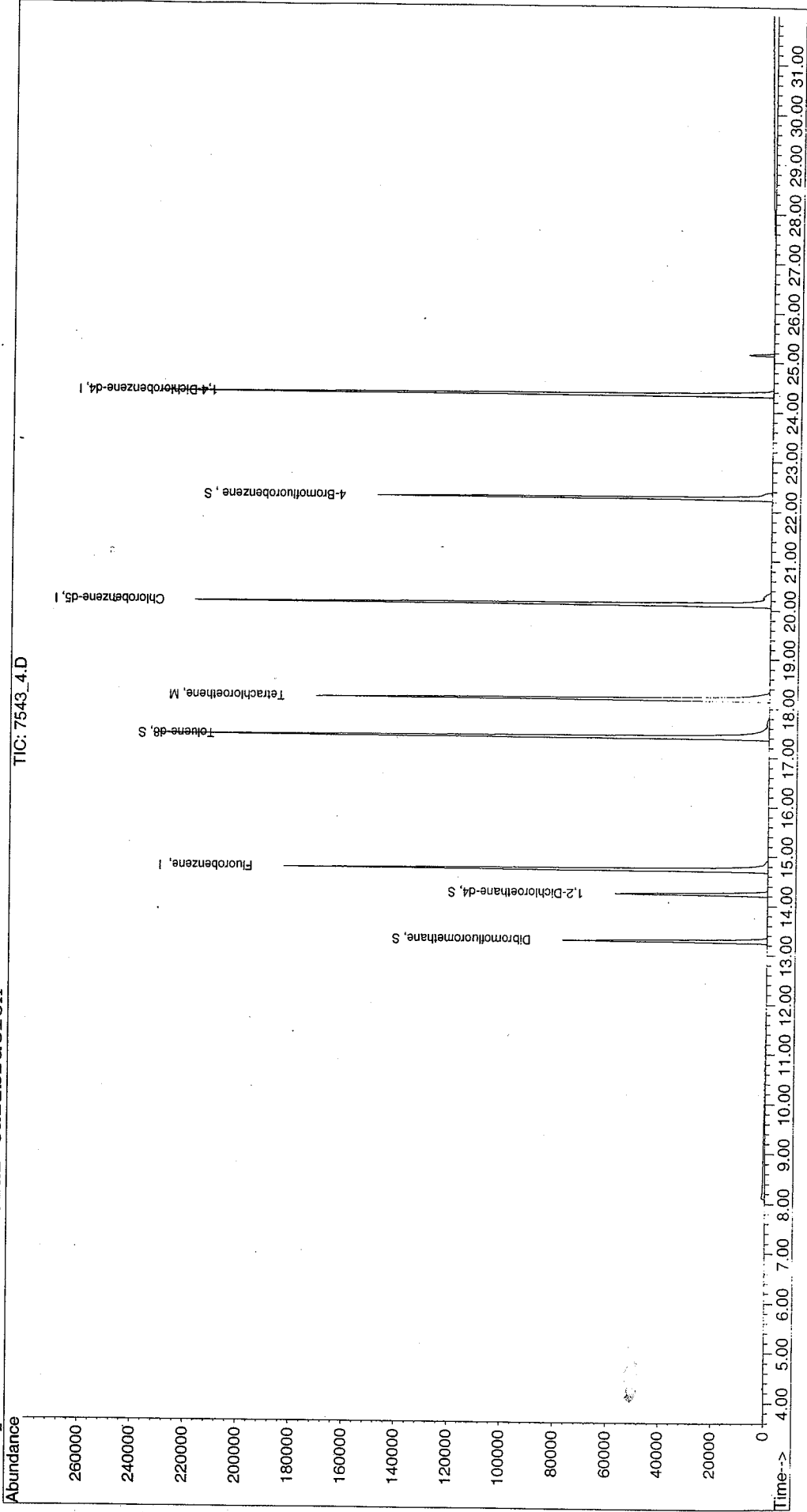
| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|--------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 99171 | 18.92 | ug/L | 99 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | | N.D. | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | | N.D. d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | | N.D. | |
| 37) Chlorobenzene | 20.15 | 112 | 5079 | | N.D. | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | | N.D. | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | | N.D. | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | | N.D. | |
| 42) Styrene | 0.00 | 104 | 0 | | N.D. | |
| 43) Bromoform | 0.00 | 173 | 0 | | N.D. | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 46) Bromobenzene | 0.00 | 156 | 0 | | N.D. | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | | N.D. | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | | N.D. | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | | N.D. | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | | N.D. | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | | N.D. | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | | N.D. | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | | N.D. | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | | N.D. | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | | N.D. | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | | N.D. | |
| 65) Naphthalene | 0.00 | 128 | 0 | | N.D. | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | | N.D. | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | | N.D. | |
| 68) Acetone | 0.00 | 43 | 0 | | N.D. | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | | N.D. | |
| 70) 2-Butanone | 0.00 | 43 | 0 | | N.D. | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | | N.D. | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | | N.D. | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | | N.D. | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration
7543_4.D 01_12_07.M Wed Feb 14 16:09:52 2007

Quantitation Report

Data File : C:\HPCHEM\1\DATA\020807\7543_4.D
Acq On : 8 Feb 2007 11:00 pm
Sample : 7543_4 0.010 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Feb 13 15:38 2007
Vial: 11
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: 01_12_07.RES

Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:54:07 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\020807\7543_5.D

Acq On : 8 Feb 2007 11:38 pm

Sample : 7543_5 0.010 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Feb 13 15:39 2007

Vial: 12

Operator: RR/AS

Inst : GC/MS Ins

Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:51:14 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.73 | 96 | 249935 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 20.12 | 117 | 206996 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.35 | 152 | 107016 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|--------|------|
| 16) Dibromofluoromethane | 13.28 | 113 | 65332 | 24.09 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 96.36% | |
| 19) 1,2-Dichloroethane-d4 | 14.24 | 65 | 60188 | 21.90 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 87.60% | |
| 28) Toluene-d8 | 17.41 | 98 | 263120 | 23.31 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 93.24% | |
| 45) 4-Bromofluorobenzene | 22.27 | 95 | 74654 | 21.12 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 84.48% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|--------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

7543_5.D 01_12_07.M

Wed Feb 14 16:09:59 2007

Data File : C:\HPCHEM\1\DATA\020807\7543_5.D

Vial: 12

Acq On : 8 Feb 2007 11:38 pm

Operator: RR/AS

Sample : 7543_5 0.010 ml

Inst : GC/MS Ins

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Feb 13 15:39 2007

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:51:14 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 106530 | 21.82 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7543_5.D 01_12_07.M

Wed Feb 14 16:10:00 2007

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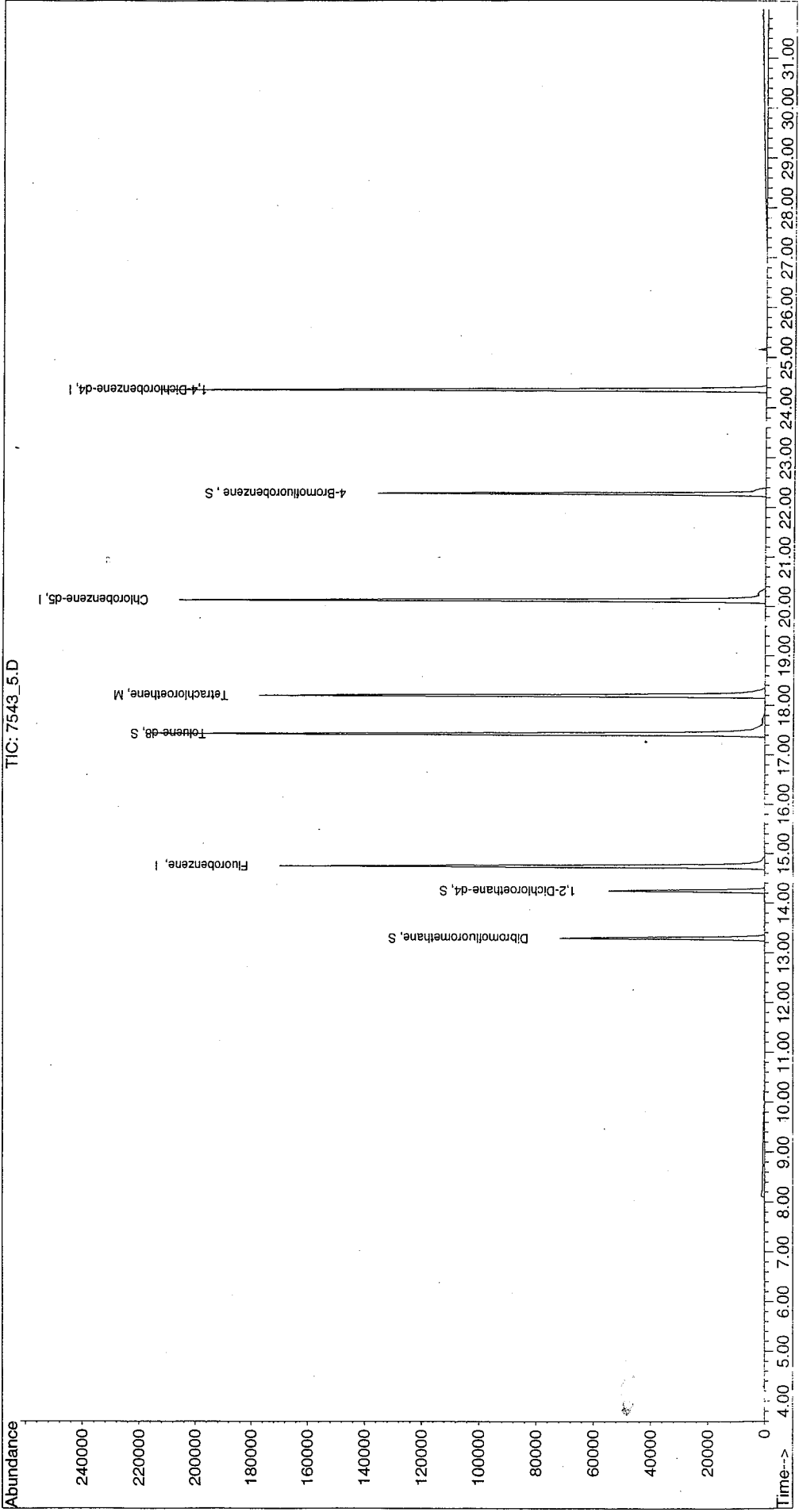
Quantitation report

Data File : C:\HPCHEM\1\DATA\020807\7543_5.D
Acq On : 8 Feb 2007 11:38 pm
Sample : 7543_5 0.010 ml
Misc :
Vial: 12
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00

MS Integration Params: ODD.P
Quant Time: Feb 13 15:39 2007

Quant Results File: 01_12_07.RES

Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:54:07 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\020807\7543_6.D
 Acq On : 9 Feb 2007 12:16 am
 Sample : 7543_6 0.010 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Feb 9 11:53 2007

Vial: 13
 Operator: RR/AS
 Inst : GC/MS Ins
 Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Mon Jan 15 14:51:14 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.73 | 96 | 269400 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 20.12 | 117 | 220734 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.35 | 152 | 112609 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|--------|-------|
| 16) Dibromofluoromethane | 13.28 | 113 | 69173 | 23.66 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 94.64% | |
| 19) 1,2-Dichloroethane-d4 | 14.23 | 65 | 63511 | 21.44 | ug/L | -0.01 |
| Spiked Amount | 25.000 | | Recovery | = | 85.76% | |
| 28) Toluene-d8 | 17.42 | 98 | 282073 | 23.18 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 92.72% | |
| 45) 4-Bromofluorobenzene | 22.27 | 95 | 79504 | 21.09 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 84.36% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\020807\7543_6.D

Vial: 13

Acq On : 9 Feb 2007 12:16 am

Operator: RR/AS

Sample : 7543_6 0.010 ml

Inst : GC/MS Ins

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Feb 9 11:53 2007

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:51:14 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 87227 | 16.57 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 20.15 | 112 | 5025 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7543_6.D 01_12_07.M

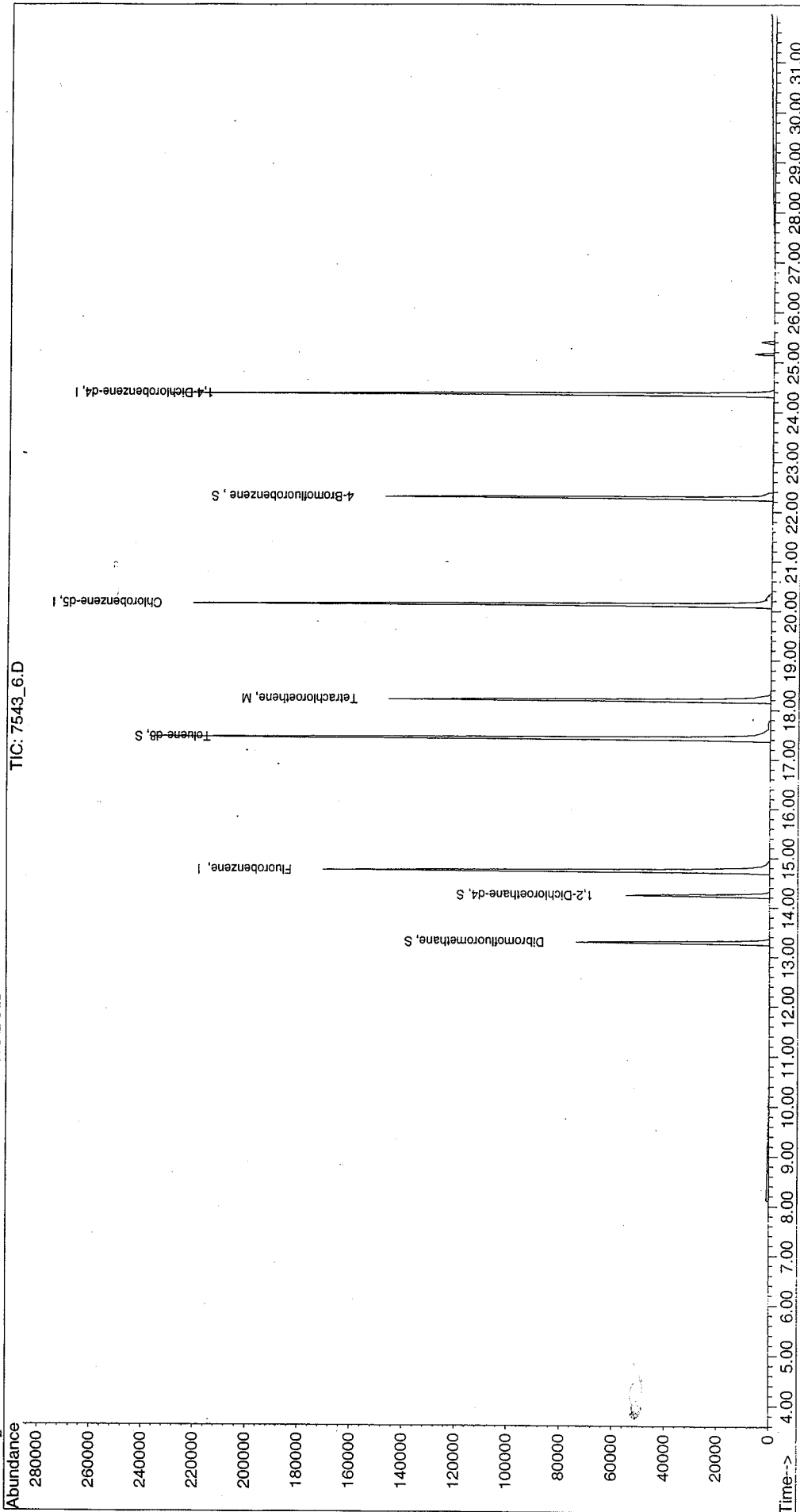
Wed Feb 14 16:10:09 2007

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QUALIFICATION REPORT

Data File : C:\HPCHEM\1\DATA\020807\7543_6.D
Acq On : 9 Feb 2007 12:16 am
Sample : 7543_6 0.010 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Feb 9 11:53 2007
Vial: 13
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: 01_12_07.RES

Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:54:07 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\020807\7543_7.D
 Acq On : 9 Feb 2007 12:54 am
 Sample : 7543_7 0.010 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Feb 13 15:40 2007

Vial: 14
 Operator: RR/AS
 Inst : GC/MS Ins
 Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Mon Jan 15 14:51:14 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.73 | 96 | 259154 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 20.12 | 117 | 211853 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.35 | 152 | 105514 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|--------|-------|
| 16) Dibromofluoromethane | 13.29 | 113 | 65950 | 23.45 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 93.80% | |
| 19) 1,2-Dichloroethane-d4 | 14.23 | 65 | 61117 | 21.45 | ug/L | -0.01 |
| Spiked Amount | 25.000 | | Recovery | = | 85.80% | |
| 28) Toluene-d8 | 17.42 | 98 | 269553 | 23.03 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 92.12% | |
| 45) 4-Bromofluorobenzene | 22.27 | 95 | 75101 | 20.76 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 83.04% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | | N.D. | |
| 3) Chloromethane | 0.00 | 50 | 0 | | N.D. | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | | N.D. | |
| 5) Bromomethane | 0.00 | 94 | 0 | | N.D. | |
| 6) Chloroethane | 0.00 | 64 | 0 | | N.D. | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | | N.D. | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 9) Methylene chloride | 0.00 | 84 | 0 | | N.D. | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | | N.D. | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | | N.D. | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | | N.D. | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | | N.D. | |
| 15) Chloroform | 0.00 | 83 | 0 | | N.D. | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | | N.D. | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | | N.D. | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 21) Benzene | 0.00 | 78 | 0 | | N.D. | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | | N.D. | |
| 23) Trichloroethene | 0.00 | 95 | 0 | | N.D. | d |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | | N.D. | |
| 25) Dibromomethane | 0.00 | 93 | 0 | | N.D. | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | | N.D. | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 29) Toluene | 0.00 | 92 | 0 | | N.D. | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | | N.D. | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | | N.D. | |

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\020807\7543_7.D

Acq On : 9 Feb 2007 12:54 am

Sample : 7543_7 0.010 ml

Misc :

MS Integration Params: ODD.P

Quant Time: Feb 13 15:40 2007

Vial: 14

Operator: RR/AS

Inst : GC/MS Ins

Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:51:14 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 18.19 | 166 | 104518 | 20.64 | ug/L | 97 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

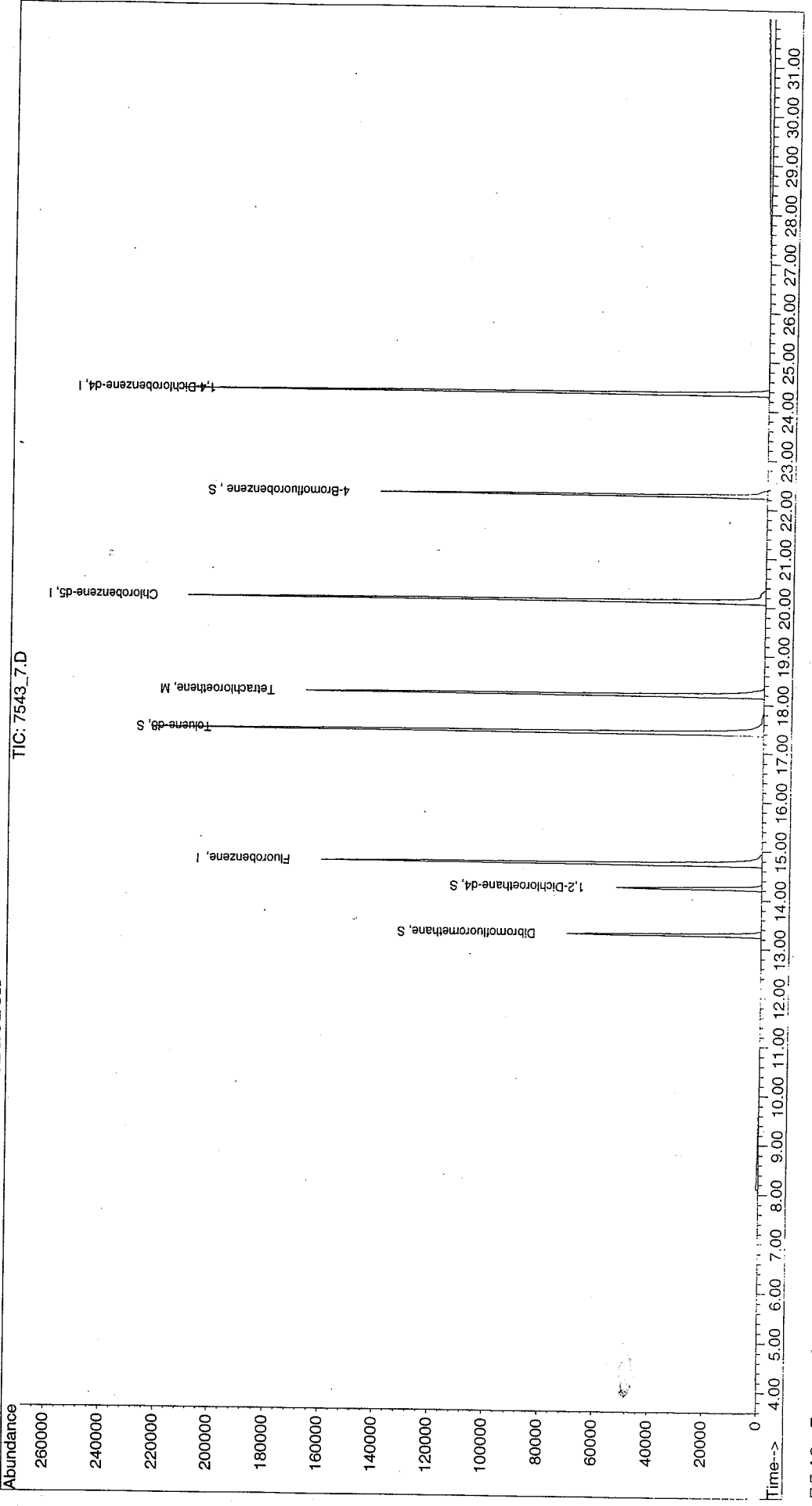
7543_7.D 01_12_07.M

Wed Feb 14 16:10:18 2007

Page 2

Quantitation Report

Data File : C:\HPCHEM\1\DATA\020807\7543_7.D
Acq On : 9 Feb 2007 12:54 am
Sample : 7543_7 0.010 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Feb 13 15:40 2007
Vial: 14
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: 01_12_07.RES
Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:54:07 2007
Response via : Initial Calibration



Acq On : 14 Feb 2007 12:16 pm

Operator: RR/AS

Sample : 7543_08 0.010 ml

Inst : GC/MS Ins

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Feb 14 15:18 2007

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:54:07 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.73 | 96 | 230994 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 20.12 | 117 | 194202 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.35 | 152 | 103007 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|------|
| 16) Dibromofluoromethane | 13.28 | 113 | 70281 | 28.03 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 112.12% | |
| 19) 1,2-Dichloroethane-d4 | 14.24 | 65 | 66953 | 26.36 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 105.44% | |
| 28) Toluene-d8 | 17.42 | 98 | 253908 | 24.34 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 97.36% | |
| 45) 4-Bromofluorobenzene | 22.27 | 95 | 66969 | 20.19 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 80.76% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene. | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#)= qualifier out of range (m) = manual integration

7543_08.D 01_12_07.M

Mon Feb 26 14:25:02 2007

Page 1

Data File : C:\HPCHEM\1\DATA\0201407\7543_08.D

Vial: 7

Acq On : 14 Feb 2007 12:16 pm

Operator: RR/AS

Sample : 7543_08 0.010 ml

Inst : GC/MS Ins

Misc :

Multiplr: 1.00

MS Integration Params: ODD:P

Quant Time: Feb 14 15:18 2007

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:54:07 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 18.18 | 166 | 118934 | 26.36 | ug/L | 99 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

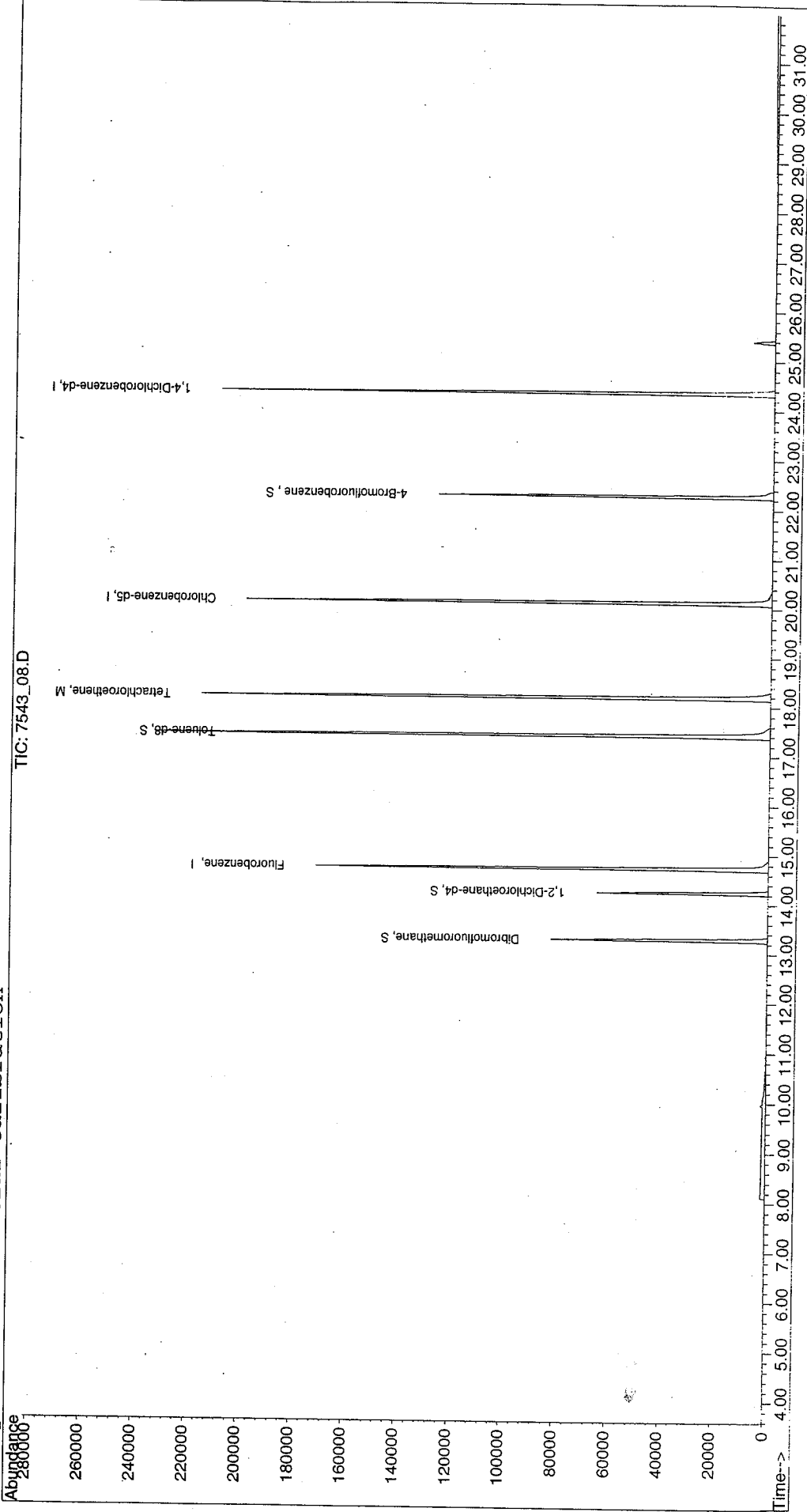
7543_08.D 01_12_07.M Mon Feb 26 14:25:04 2007

Page 2

Quantitation Report

Data File : C:\HPCHEM\1\DATA\0201407\7543_08.D
Acq On : 14 Feb 2007 12:16 pm
Sample : 7543_08 0.010 ml
Misc :
MS Integration Params: ODD.P
Quant Time: Feb 14 15:18 2007
Vial: 7
Operator: RR/AS
Inst : GC/MS Ins
Multiplr: 1.00
Quant Results File: 01_12_07.RES

Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
Title : EPA Method 8260A
Last Update : Mon Jan 15 14:54:07 2007
Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\0201407\7543_09.D
 Acq On : 14 Feb 2007 12:54 pm
 Sample : 7543_09 0.010 ml
 Misc :
 MS Integration Params: ODD.P
 Quant Time: Feb 14 15:19 2007

Vial: 8
 Operator: RR/AS
 Inst : GC/MS Ins
 Multiplr: 1.00

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)
 Title : EPA Method 8260A
 Last Update : Mon Jan 15 14:54:07 2007
 Response via : Initial Calibration
 DataAcq Meth : RUN

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------|-------|------|----------|-------|-------|----------|
| 1) Fluorobenzene | 14.73 | 96 | 237154 | 25.00 | ug/L | 0.00 |
| 36) Chlorobenzene-d5 | 20.11 | 117 | 198219 | 25.00 | ug/L | 0.00 |
| 58) 1,4-Dichlorobenzene-d4 | 24.35 | 152 | 104340 | 25.00 | ug/L | 0.00 |

System Monitoring Compounds

| | | | | | | |
|---------------------------|--------|-----|----------|-------|---------|------|
| 16) Dibromofluoromethane | 13.29 | 113 | 71674 | 27.85 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 111.40% | |
| 19) 1,2-Dichloroethane-d4 | 14.24 | 65 | 67292 | 25.80 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 103.20% | |
| 28) Toluene-d8 | 17.41 | 98 | 262935 | 24.55 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 98.20% | |
| 45) 4-Bromofluorobenzene | 22.27 | 95 | 67181 | 19.85 | ug/L | 0.00 |
| Spiked Amount | 25.000 | | Recovery | = | 79.40% | |

Target Compounds

| Target Compounds | R.T. | QIon | Response | Conc | Units | Qvalue |
|-------------------------------|------|------|----------|------|-------|--------|
| 2) Dichlorodifluoromethane | 0.00 | 85 | 0 | N.D. | | |
| 3) Chloromethane | 0.00 | 50 | 0 | N.D. | | |
| 4) Vinyl chloride | 0.00 | 62 | 0 | N.D. | | |
| 5) Bromomethane | 0.00 | 94 | 0 | N.D. | | |
| 6) Chloroethane | 0.00 | 64 | 0 | N.D. | | |
| 7) Trichlorofluoromethane | 0.00 | 101 | 0 | N.D. | | |
| 8) 1,1-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 9) Methylene chloride | 0.00 | 84 | 0 | N.D. | | |
| 10) trans-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 11) 1,1-Dichloroethane | 0.00 | 63 | 0 | N.D. | | |
| 12) 2,2-Dichloropropane | 0.00 | 77 | 0 | N.D. | | |
| 13) cis-1,2-Dichloroethene | 0.00 | 96 | 0 | N.D. | | |
| 14) Bromochloromethane | 0.00 | 128 | 0 | N.D. | | |
| 15) Chloroform | 0.00 | 83 | 0 | N.D. | | |
| 17) 1,1,1-Trichloroethane | 0.00 | 97 | 0 | N.D. | | |
| 18) Carbon tetrachloride | 0.00 | 117 | 0 | N.D. | | |
| 20) 1,1-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 21) Benzene | 0.00 | 78 | 0 | N.D. | | |
| 22) 1,2-Dichloroethane | 0.00 | 62 | 0 | N.D. | | |
| 23) Trichloroethene | 0.00 | 95 | 0 | N.D. | d | |
| 24) 1,2-Dichloropropane | 0.00 | 63 | 0 | N.D. | | |
| 25) Dibromomethane | 0.00 | 93 | 0 | N.D. | | |
| 26) Bromodichloromethane | 0.00 | 83 | 0 | N.D. | | |
| 27) cis-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 29) Toluene | 0.00 | 92 | 0 | N.D. | | |
| 30) trans-1,3-Dichloropropene | 0.00 | 75 | 0 | N.D. | | |
| 31) 1,1,2-Trichloroethane | 0.00 | 83 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7543_09.D 01_12_07.M Mon Feb 26 14:25:13 2007

Data File : C:\HPCHEM\1\DATA\0201407\7543_09.D

Vial: 8

Acq On : 14 Feb 2007 12:54 pm

Operator: RR/AS

Sample : 7543_09 0.010 ml

Inst : GC/MS Ins

Misc :

Multiplr: 1.00

MS Integration Params: ODD.P

Quant Time: Feb 14 15:19 2007

Quant Results File: 01_12_07.RES

Quant Method : C:\HPCHEM\1\METHODS\01_12_07.M (RTE Integrator)

Title : EPA Method 8260A

Last Update : Mon Jan 15 14:54:07 2007

Response via : Initial Calibration

DataAcq Meth : RUN

| Compound | R.T. | QIon | Response | Conc | Unit | Qvalue |
|---------------------------------|-------|------|----------|-------|------|--------|
| 32) Tetrachloroethene | 18.18 | 166 | 98299 | 21.22 | ug/L | 98 |
| 33) 1,3-Dichloropropane | 0.00 | 76 | 0 | N.D. | | |
| 34) Dibromochloromethane | 0.00 | 129 | 0 | N.D. | d | |
| 35) 1,2-Dibromomethane | 0.00 | 107 | 0 | N.D. | | |
| 37) Chlorobenzene | 0.00 | 112 | 0 | N.D. | | |
| 38) 1,1,1,2-Tetrachloroethane | 0.00 | 131 | 0 | N.D. | | |
| 39) Ethylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 40) Xylene (para & meta) | 0.00 | 106 | 0 | N.D. | | |
| 41) Xylene (Ortho) | 0.00 | 106 | 0 | N.D. | | |
| 42) Styrene | 0.00 | 104 | 0 | N.D. | | |
| 43) Bromoform | 0.00 | 173 | 0 | N.D. | | |
| 44) Isopropylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 46) Bromobenzene | 0.00 | 156 | 0 | N.D. | | |
| 47) 1,1,2,2-Tetrachloroethane | 0.00 | 83 | 0 | N.D. | | |
| 48) 1,2,3-Trichloropropane | 0.00 | 110 | 0 | N.D. | | |
| 49) n-Propylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 50) 2-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 51) 4-Chlorotoluene | 0.00 | 91 | 0 | N.D. | | |
| 52) 1,3,5-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 53) tert-Butylbenzene | 0.00 | 119 | 0 | N.D. | | |
| 54) 1,2,4-Trimethylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 55) sec-Butylbenzene | 0.00 | 105 | 0 | N.D. | | |
| 56) 1,3-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 57) 4-Isopropyltoluene | 0.00 | 119 | 0 | N.D. | | |
| 59) 1,4-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 60) 1,2-Dichlorobenzene | 0.00 | 146 | 0 | N.D. | | |
| 61) n-Butylbenzene | 0.00 | 91 | 0 | N.D. | | |
| 62) 1,2-Dibromo-3-chloropropan | 0.00 | 75 | 0 | N.D. | | |
| 63) 1,2,4-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 64) Hexachlorobutadiene | 0.00 | 225 | 0 | N.D. | | |
| 65) Naphthalene | 0.00 | 128 | 0 | N.D. | | |
| 66) 1,2,3-Trichlorobenzene | 0.00 | 180 | 0 | N.D. | | |
| 67) Methyl-tert butyl ether | 0.00 | 73 | 0 | N.D. | | |
| 68) Acetone | 0.00 | 43 | 0 | N.D. | | |
| 69) Carbon disulfide | 0.00 | 76 | 0 | N.D. | | |
| 70) 2-Butanone | 0.00 | 43 | 0 | N.D. | | |
| 71) Tetrahydrofuran | 0.00 | 42 | 0 | N.D. | | |
| 72) 4-Methyl-2-pentanone (MIBK) | 0.00 | 58 | 0 | N.D. | | |
| 73) 2-Hexanone | 0.00 | 43 | 0 | N.D. | | |
| 74) 2-Chloroethyl vinyl ether | 0.00 | 63 | 0 | N.D. | | |

(#) = qualifier out of range (m) = manual integration

7543_09.D 01_12_07.M

Mon Feb 26 14:25:15 2007

Page 2

COVER PAGE

OrderID: Y1443 **ProjectID:** Textron
CustomerName: Shaw E & I

LAB SAMPLE NO.
Y1443-01
Y1443-02

CLIENT SAMPLE NO
MW-101D
SB-219A(29-29)

I certify that the data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hard copy data package has been authorized by the laboratory manager or his designee, as verified by the following signature.

Signature: Mildred V Reyes Name: Mildred V Reyes
Date: 2/21/07 Title: QA/QC

Y1443

17 Princess Rd
Lawrenceville, NJ 08648
609-895-5370/ 609-895-1858
Shaw
Shaw Environmental and Infrastructure Inc.

CHAIN OF CUSTODY

Ref. Document #

Page 1 of 1

Project Number/Cost code: 101960 / 0902000
Project Name / Location: Textron /
Purchase Order #: 87235

Project Contact: _____
(Name & phone #)

Send Report To: Rand: Potune
Phone/Fax Number: _____
Address: 45 Home
City/State: _____

Shipment Date: _____
Waybill/Airbill Number: _____
Lab Destination: CharTech
Lab Contact Name / ph. #: C

| Analyses Requested | Turn Around Time Requested |
|---|----------------------------|
| Priority Metals (13) | X |
| Sb, As, Be, Cd, Cr, Cu, Pb, Hg, Ni, Se, V, Tl, Zn | X |
| SVOCs 8230 | X |

| Collection Information | Preservative | | | | |
|------------------------|--------------------------------|------------------|------|-----|-----|
| | H ₂ SO ₄ | HNO ₃ | NaOH | HCL | Ice |
| Date: 12/16/06 | | X | | | |
| Time: 1235 | | | | | |
| G/C: 6 | | | | | |
| Date: 12/16/06 | | | | | |
| Time: 130 | | | | | |
| G/C: 5 | | | | | |

| Lab No. | Sample ID Number | Sample Description | Matrix | # of containers | Container type |
|---------|------------------|--------------------|--------|-----------------|----------------|
| | MW-101D | ↓ | A | 1 | 120 |
| | SB-219A 27-29 | | S | 1 | 230 |

Special Instructions:
Analyze SVOCs eventhough out of hold

Known Waste Stream Circle:
RCRA PCB/dioxin PAH/coil RAD Corrosive Flammable Reactive G = Grab

QC Package Level Required:
I II III IV NJ EDD GIS EDD Preliminary data

QC Package Codes:
Level I = data summary
Level II = data summary + basic QC
Level III = New Jersey QC reduced deliverable
Level IV = Full deliverable CLP package

Received By:
Date: 2/5/07
Time: 1330

Relinquished By:
Date: 2-6-07
Time: 11:15

Received By:
Date: 2-6-07
Time: 11:15

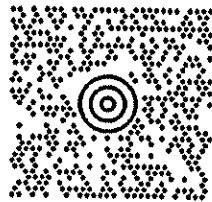

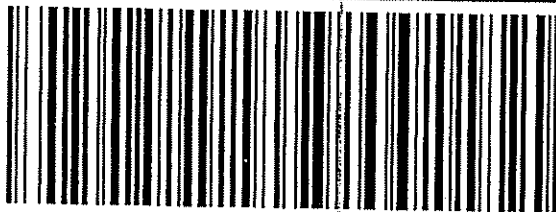
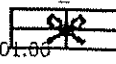
Relinquished By:
Date: 2-6-07
Time: 11:15

Cooler temperature upon arrival at Lab:
5°C

UPS

D. Jackson

https://www.campusship.ups.com/ship/create?ActionOriginPair=print__PrinterPage&PO... 2/5/2007

| | | |
|--|--|---|
| MARION HORVATH 609.895.5340 LAWRENCEVILLE 15386E 17 PRINCESS RD LAWRENCEVILLE NJ 08648 | 55 LBS DWT:24,13,15 | 1 OF 1 |
| SHIP TO: CHEMTECH 908 789 8900 284 SHEFFIELD ST. MOUNTAINSIDE NJ 07092-2319 | | |
|  | NJ 078 9-61  | |
| UPS GROUND TRACKING #: 1Z 153 86E 03 9683 1822 | | |
|  | | |
| BILLING: F/C BILL RECEIVER | |  |
| Project Number or Department Number: 00501501414425.5701.00 Sender's Name: randi rothmel | | TM |
| CS 9.0.19.0 WAPES0 60.1A 10/2006 | | |

JL
2-6-07
11:15

000000

Laboratory Certification

| State | License No. |
|----------------|-------------|
| New Jersey | 20012 |
| New York | 11376 |
| Arizona | AZ0653 |
| Connecticut | PH-0649 |
| Florida | E87935 |
| Kansas | E-10355 |
| Maryland | 296 |
| Massachusetts | M-NJ503 |
| Maine | NJ0503 |
| North Carolina | 630 |
| Oklahoma | 9705 |
| Pennsylvania | 68-548 |
| Rhode Island | LAO00259 |

QA Control Code: A2070148

DATA REPORTING QUALIFIERS- ORGANIC

For reporting results, the following " Results Qualifiers" are used:

- Value If the result is a value greater than or equal to the detection limit, report the value
- U** Indicates the compound was analyzed for but was not detected. Report the minimum detection limit for the sample with the U, i.e. "10 U". This is not necessarily the instrument detection limit attainable for this particular sample based on any concentration or dilution that may have been required.
- J** Indicates an estimated value. This flag is used:
- (1) When estimating a concentration for a tentatively identified compound (library search hits, where a 1:1 response is assumed.)
 - (2) When the mass spectral data indicated the identification, however the result was less than the specified detection limit greater than zero. If the detection limit was 10ug/L and a concentration of 3 ug/L was calculated report as 3 J. This is flag is used when similar situation arise on any organic parameter i.e. Pest, PCB and others.
- B** Indicates the analyte was found in the blank as well as the sample report as "12 B".
- E** Indicates the analyte 's concentration exceeds the calibrated range of the instrument for that specific analysis.
- D** This flag identifies all compounds identified in an analysis at a secondary dilution factor.
- P** This flag is used for Pesticide/PCB target analyte when there is >25% difference for detected concentrations between the two GC columns. The lower of the two values is reported on Form 1 and flagged with a "P".
- N** This flag indicates presumptive evidence of a compound. This is only used for tentatively identified compounds (TICs), where the identification is based on a mass spectral library search. It applies to all TIC results. For generic characterization of a TIC, such as chlorinated hydrocarbon, the flag is not used.

DATA REPORTING QUALIFIERS- INORGANIC

For reporting results, the following "Results Qualifiers" are used:

- J** If the reported value was obtained from a reading that was less than the Contract Required Detection Limit (CRDL), but greater than or equal to the Instrument Detection Limit (IDL).
- U** If the analyte was analyzed for, but not detected.
- E** The reported value is estimated because of the presence of interference
- M** Duplicate injection precision not met.
- N** Spiked sample recovery not within control limits.
- S** The reported value was determined by the Method of Standard Addition (MSA).
- W** Post-digestion spike for Furnace AA analysis is out of control limits (85-115%), while absorbance is less than 50% of spike absorbance.
- *** Duplicate analysis not within control limits.
- +** Correlation coefficient for the MSA is less than 0.995.
- ***** Entering "S", "W" or "+" is mutually exclusive. NO combination of these qualifiers can appear in the same field for an analyte.
- D** The reported value is from a secondary analysis with a dilution factor. The original analysis exceeded the calibration range.
- M** Method qualifiers
"P" for ICP instrument
"A" for Flame AA
"PM" for ICP when Microwave Digestion is used
"AM" for flame AA when Microwave Digestion is used
"FM" for furnace AA when Microwave Digestion is used
"CV" for Manual Cold Vapor AA
"AV" for automated Cold Vapor AA
"CA" for MIDI-Distillation Spectrophotometric
"AS" for Semi-Automated Spectrophotometric
"C" for Manual Spectrophotometric
"T" for Titrimetric
"NR" for analyte not required to be analyzed
- OR** Indicates the analyte's concentration exceeds the calibrated range of the instrument for that specific analysis.

APPENDIX A

QA REVIEW GENERAL DOCUMENTATION

Project #: 71664

Completed

For thorough review, the report must have the following:

GENERAL:

Are all original paperwork present (chain of custody, record of communication, airbill, sample management lab chronicle, login page) ✓
Check chain-of-custody for proper relinquish/return of samples ✓
Is the chain of custody signed and complete ✓
Check internal chain-of-custody for proper relinquish/return of samples /sample extracts ✓
Collect information for each project id from server. Were all requirements followed ✓

COVER PAGE:

Do numbers of samples correspond to the number of samples in the Chain of Custody and on login page ✓
Do lab numbers and client Ids on cover page agree with the Chain of Custody ✓

CHAIN OF CUSTODY:

Do requested analyses on Chain of Custody agree with form I results ✓
Do requested analyses on Chain of Custody agree with the log-in page ✓
Were the correct method log-in for analysis according to the Analytical Request and Chain of Custody ✓
Were the samples received within hold time ✓
Were any problems found with the samples at arrival recorded in the Sample Management Laboratory Chronicle ✓

ANALYTICAL:

Was method requirement followed? ✓
Was client requirement followed? ✓
Does the case narrative summarize all QC failure? ✓
All runlogs reviewed for manual integration requirements ✓

1st Level QA Review Signature: P. G. Pennington Date: 02/21/07

2nd Level QA Review Signature: M. J. Deo V. Reyes Date: 2/21/07

Report of Analysis

| | | | |
|---------------------------|-----------------------|------------------------|-------------------|
| Client: | Shaw E & I | Date Collected: | 12/16/2006 |
| Project: | Textron | Date Received: | 2/6/2007 |
| Client Sample ID: | MW-101D | SDG No.: | Y1443 |
| Lab Sample ID: | Y1443-01 | Matrix: | WATER |
| Analytical Method: | 8270 | % Moisture: | 100 |
| Sample Wt/Wol: | 980.0 mL | Extract Vol: | 1000 uL |

| | | | | |
|-------------------|-----------------|-----------------------|----------------------|----------------------------|
| File ID | Dilution | Date Extracted | Date Analyzed | Analytical Batch ID |
| BE038475.D | 1 | 2/8/2007 | 2/9/2007 | BE020907 |

| CAS Number | Parameter | Conc. | Qualifier | RL | MDL | Units |
|----------------|-----------------------------|-------|-----------|----|-------|-------|
| TARGETS | | | | | | |
| 108-95-2 | Phenol | 1.3 | U | 10 | 1.3 | ug/L |
| 111-44-4 | bis(2-Chloroethyl)ether | 1.5 | U | 10 | 1.5 | ug/L |
| 95-57-8 | 2-Chlorophenol | 1.2 | U | 10 | 1.2 | ug/L |
| 95-50-1 | 1,2-Dichlorobenzene | 1.2 | U | 10 | 1.2 | ug/L |
| 541-73-1 | 1,3-Dichlorobenzene | 1.2 | U | 10 | 1.2 | ug/L |
| 106-46-7 | 1,4-Dichlorobenzene | 1.2 | U | 10 | 1.2 | ug/L |
| 95-48-7 | 2-Methylphenol | 1.5 | U | 10 | 1.5 | ug/L |
| 108-60-1 | 2,2-oxybis(1-Chloropropane) | 1.2 | U | 10 | 1.2 | ug/L |
| 106-44-5 | 3+4-Methylphenols | 1.3 | U | 10 | 1.3 | ug/L |
| 621-64-7 | N-Nitroso-di-n-propylamine | 1.4 | U | 10 | 1.4 | ug/L |
| 67-72-1 | Hexachloroethane | 1.2 | U | 10 | 1.2 | ug/L |
| 98-95-3 | Nitrobenzene | 1.6 | U | 10 | 1.6 | ug/L |
| 78-59-1 | Isophorone | 1.3 | U | 10 | 1.3 | ug/L |
| 88-75-5 | 2-Nitrophenol | 1.4 | U | 10 | 1.4 | ug/L |
| 105-67-9 | 2,4-Dimethylphenol | 1.2 | U | 10 | 1.2 | ug/L |
| 111-91-1 | bis(2-Chloroethoxy)methane | 1.4 | U | 10 | 1.4 | ug/L |
| 120-83-2 | 2,4-Dichlorophenol | 1.5 | U | 10 | 1.5 | ug/L |
| 120-82-1 | 1,2,4-Trichlorobenzene | 1.4 | U | 10 | 1.4 | ug/L |
| 91-20-3 | Naphthalene | 1.4 | U | 10 | 1.4 | ug/L |
| 106-47-8 | 4-Chloroaniline | 0.880 | U | 10 | 0.880 | ug/L |
| 87-68-3 | Hexachlorobutadiene | 1.4 | U | 10 | 1.4 | ug/L |
| 59-50-7 | 4-Chloro-3-methylphenol | 1.4 | U | 10 | 1.4 | ug/L |
| 91-57-6 | 2-Methylnaphthalene | 1.1 | U | 10 | 1.1 | ug/L |
| 77-47-4 | Hexachlorocyclopentadiene | 1.2 | U | 10 | 1.2 | ug/L |
| 88-06-2 | 2,4,6-Trichlorophenol | 1.2 | U | 10 | 1.2 | ug/L |
| 95-95-4 | 2,4,5-Trichlorophenol | 1.2 | U | 10 | 1.2 | ug/L |
| 91-58-7 | 2-Chloronaphthalene | 1.4 | U | 10 | 1.4 | ug/L |
| 88-74-4 | 2-Nitroaniline | 1.1 | U | 10 | 1.1 | ug/L |
| 131-11-3 | Dimethylphthalate | 1.3 | U | 10 | 1.3 | ug/L |
| 208-96-8 | Acenaphthylene | 1.3 | U | 10 | 1.3 | ug/L |
| 606-20-2 | 2,6-Dinitrotoluene | 1.3 | U | 10 | 1.3 | ug/L |

U = Not Detected

RL = Reporting Limit

MDL = Method Detection Limit

E = Value Exceeds Calibration Range

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

Report of Analysis

| | | | |
|---------------------------|------------|------------------------|------------|
| Client: | Shaw E & I | Date Collected: | 12/16/2006 |
| Project: | Textron | Date Received: | 2/6/2007 |
| Client Sample ID: | MW-101D | SDG No.: | Y1443 |
| Lab Sample ID: | Y1443-01 | Matrix: | WATER |
| Analytical Method: | 8270 | % Moisture: | 100 |
| Sample Wt/Wol: | 980.0 mL | Extract Vol: | 1000 uL |

| | | | | |
|----------------|-----------------|-----------------------|----------------------|----------------------------|
| File ID | Dilution | Date Extracted | Date Analyzed | Analytical Batch ID |
| BE038475.D | 1 | 2/8/2007 | 2/9/2007 | BE020907 |

| CAS Number | Parameter | Conc. | Qualifier | RL | MDL | Units |
|----------------|----------------------------|-------|-----------|----|-------|-------|
| TARGETS | | | | | | |
| 99-09-2 | 3-Nitroaniline | 1.0 | U | 10 | 1.0 | ug/L |
| 83-32-9 | Acenaphthene | 1.4 | U | 10 | 1.4 | ug/L |
| 51-28-5 | 2,4-Dinitrophenol | 3.6 | U | 10 | 3.6 | ug/L |
| 100-02-7 | 4-Nitrophenol | 3.2 | U | 10 | 3.2 | ug/L |
| 132-64-9 | Dibenzofuran | 1.3 | U | 10 | 1.3 | ug/L |
| 121-14-2 | 2,4-Dinitrotoluene | 1.2 | U | 10 | 1.2 | ug/L |
| 84-66-2 | Diethylphthalate | 1.4 | U | 10 | 1.4 | ug/L |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 1.4 | U | 10 | 1.4 | ug/L |
| 86-73-7 | Fluorene | 1.4 | U | 10 | 1.4 | ug/L |
| 100-01-6 | 4-Nitroaniline | 1.1 | U | 10 | 1.1 | ug/L |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 1.6 | U | 10 | 1.6 | ug/L |
| 86-30-6 | N-Nitrosodiphenylamine | 1.3 | U | 10 | 1.3 | ug/L |
| 101-55-3 | 4-Bromophenyl-phenylether | 1.5 | U | 10 | 1.5 | ug/L |
| 118-74-1 | Hexachlorobenzene | 1.2 | U | 10 | 1.2 | ug/L |
| 87-86-5 | Pentachlorophenol | 1.6 | U | 10 | 1.6 | ug/L |
| 85-01-8 | Phenanthrene | 1.4 | U | 10 | 1.4 | ug/L |
| 120-12-7 | Anthracene | 1.4 | U | 10 | 1.4 | ug/L |
| 86-74-8 | Carbazole | 1.3 | U | 10 | 1.3 | ug/L |
| 84-74-2 | Di-n-butylphthalate | 1.3 | U | 10 | 1.3 | ug/L |
| 206-44-0 | Fluoranthene | 1.2 | U | 10 | 1.2 | ug/L |
| 129-00-0 | Pyrene | 1.5 | U | 10 | 1.5 | ug/L |
| 85-68-7 | Butylbenzylphthalate | 1.5 | U | 10 | 1.5 | ug/L |
| 91-94-1 | 3,3-Dichlorobenzidine | 1.1 | U | 10 | 1.1 | ug/L |
| 56-55-3 | Benzo(a)anthracene | 1.1 | U | 10 | 1.1 | ug/L |
| 218-01-9 | Chrysene | 1.7 | U | 10 | 1.7 | ug/L |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | 1.7 | J | 10 | 1.6 | ug/L |
| 117-84-0 | Di-n-octyl phthalate | 1.3 | U | 10 | 1.3 | ug/L |
| 205-99-2 | Benzo(b)fluoranthene | 0.760 | U | 10 | 0.760 | ug/L |
| 207-08-9 | Benzo(k)fluoranthene | 1.9 | U | 10 | 1.9 | ug/L |
| 50-32-8 | Benzo(a)pyrene | 1.2 | U | 10 | 1.2 | ug/L |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 0.840 | U | 10 | 0.840 | ug/L |

U = Not Detected

RL = Reporting Limit

MDL = Method Detection Limit

E = Value Exceeds Calibration Range

J = Estimated Value

B = Analyte Found In Associated Method Blank

N = Presumptive Evidence of a Compound

Report of Analysis

| | | | |
|---------------------------|------------|------------------------|------------|
| Client: | Shaw E & I | Date Collected: | 12/16/2006 |
| Project: | Textron | Date Received: | 2/6/2007 |
| Client Sample ID: | MW-101D | SDG No.: | Y1443 |
| Lab Sample ID: | Y1443-01 | Matrix: | WATER |
| Analytical Method: | 8270 | % Moisture: | 100 |
| Sample Wt/Wol: | 980.0 mL | Extract Vol: | 1000 uL |

| | | | | |
|----------------|-----------------|-----------------------|----------------------|----------------------------|
| File ID | Dilution | Date Extracted | Date Analyzed | Analytical Batch ID |
| BE038475.D | 1 | 2/8/2007 | 2/9/2007 | BE020907 |

| CAS Number | Parameter | Conc. | Qualifier | RL | MDL | Units |
|---------------------------------------|------------------------|--------|-----------|----------|-------|---------|
| TARGETS | | | | | | |
| 53-70-3 | Dibenz(a,h)anthracene | 0.880 | U | 10 | 0.880 | ug/L |
| 191-24-2 | Benzo(g,h,i)perylene | 1.1 | U | 10 | 1.1 | ug/L |
| SURROGATES | | | | | | |
| 367-12-4 | 2-Fluorophenol | 83.21 | 55 % | 21 - 100 | | SPK: 15 |
| 13127-88-3 | Phenol-d5 | 87.79 | 59 % | 10 - 94 | | SPK: 15 |
| 4165-60-0 | Nitrobenzene-d5 | 63.65 | 64 % | 35 - 114 | | SPK: 10 |
| 321-60-8 | 2-Fluorobiphenyl | 65.85 | 66 % | 43 - 116 | | SPK: 10 |
| 118-79-6 | 2,4,6-Tribromophenol | 112.57 | 75 % | 10 - 123 | | SPK: 15 |
| 1718-51-0 | Terphenyl-d14 | 73.31 | 73 % | 33 - 141 | | SPK: 10 |
| INTERNAL STANDARDS | | | | | | |
| 3855-82-1 | 1,4-Dichlorobenzene-d4 | 78240 | 4.28 | | | |
| 1146-65-2 | Naphthalene-d8 | 272632 | 5.44 | | | |
| 15067-26-2 | Acenaphthene-d10 | 130124 | 7.14 | | | |
| 1517-22-2 | Phenanthrene-d10 | 207421 | 8.61 | | | |
| 1719-03-5 | Chrysene-d12 | 160862 | 11.21 | | | |
| 1520-96-3 | Perylene-d12 | 92027 | 12.85 | | | |
| TENTITIVE IDENTIFIED COMPOUNDS | | | | | | |
| 127-18-4 | Tetrachloroethylene | 260 | J | 2.70 | | ug/L |
| | ACP3.00 | 46 | A | 3.00 | | ug/L |

U = Not Detected
 RL = Reporting Limit
 MDL = Method Detection Limit
 E = Value Exceeds Calibration Range

J = Estimated Value
 B = Analyte Found In Associated Method Blank
 N = Presumptive Evidence of a Compound

Report of Analysis

| | | | |
|---------------------------|-------------------------------|------------------------|------------|
| Client: | Shaw E & I | Date Collected: | 12/16/2006 |
| Project: | Textron | Date Received: | 2/6/2007 |
| Client Sample ID: | SB-219A ²⁷ (29-29) | SDG No.: | Y1443 |
| Lab Sample ID: | Y1443-02 | Matrix: | SOIL |
| Analytical Method: | 8270 | % Moisture: | 15 |
| Sample Wt/Wol: | 30.1 g | Extract Vol: | 1000 uL |

| | | | | |
|----------------|-----------------|-----------------------|----------------------|----------------------------|
| File ID | Dilution | Date Extracted | Date Analyzed | Analytical Batch ID |
| BB036131.D | 1 | 2/8/2007 | 2/8/2007 | BB012907 |

| CAS Number | Parameter | Conc. | Qualifier | RL | MDL | Units |
|----------------|-----------------------------|-------|-----------|-----|-----|-------|
| TARGETS | | | | | | |
| 108-95-2 | Phenol | 240 | J | 390 | 59 | ug/Kg |
| 111-44-4 | bis(2-Chloroethyl)ether | 61 | U | 390 | 61 | ug/Kg |
| 95-57-8 | 2-Chlorophenol | 62 | U | 390 | 62 | ug/Kg |
| 95-50-1 | 1,2-Dichlorobenzene | 58 | U | 390 | 58 | ug/Kg |
| 541-73-1 | 1,3-Dichlorobenzene | 61 | U | 390 | 61 | ug/Kg |
| 106-46-7 | 1,4-Dichlorobenzene | 68 | U | 390 | 68 | ug/Kg |
| 95-48-7 | 2-Methylphenol | 65 | U | 390 | 65 | ug/Kg |
| 108-60-1 | 2,2-oxybis(1-Chloropropane) | 63 | U | 390 | 63 | ug/Kg |
| 106-44-5 | 3+4-Methylphenols | 61 | U | 390 | 61 | ug/Kg |
| 621-64-7 | N-Nitroso-di-n-propylamine | 64 | U | 390 | 64 | ug/Kg |
| 67-72-1 | Hexachloroethane | 66 | U | 390 | 66 | ug/Kg |
| 98-95-3 | Nitrobenzene | 85 | U | 390 | 85 | ug/Kg |
| 78-59-1 | Isophorone | 58 | U | 390 | 58 | ug/Kg |
| 88-75-5 | 2-Nitrophenol | 60 | U | 390 | 60 | ug/Kg |
| 105-67-9 | 2,4-Dimethylphenol | 62 | U | 390 | 62 | ug/Kg |
| 111-91-1 | bis(2-Chloroethoxy)methane | 64 | U | 390 | 64 | ug/Kg |
| 120-83-2 | 2,4-Dichlorophenol | 72 | U | 390 | 72 | ug/Kg |
| 120-82-1 | 1,2,4-Trichlorobenzene | 66 | U | 390 | 66 | ug/Kg |
| 91-20-3 | Naphthalene | 66 | U | 390 | 66 | ug/Kg |
| 106-47-8 | 4-Chloroaniline | 46 | U | 390 | 46 | ug/Kg |
| 87-68-3 | Hexachlorobutadiene | 60 | U | 390 | 60 | ug/Kg |
| 59-50-7 | 4-Chloro-3-methylphenol | 54 | U | 390 | 54 | ug/Kg |
| 91-57-6 | 2-Methylnaphthalene | 65 | U | 390 | 65 | ug/Kg |
| 77-47-4 | Hexachlorocyclopentadiene | 62 | U | 390 | 62 | ug/Kg |
| 88-06-2 | 2,4,6-Trichlorophenol | 57 | U | 390 | 57 | ug/Kg |
| 95-95-4 | 2,4,5-Trichlorophenol | 59 | U | 970 | 59 | ug/Kg |
| 91-58-7 | 2-Chloronaphthalene | 64 | U | 390 | 64 | ug/Kg |
| 88-74-4 | 2-Nitroaniline | 49 | U | 970 | 49 | ug/Kg |
| 131-11-3 | Dimethylphthalate | 62 | U | 390 | 62 | ug/Kg |
| 208-96-8 | Acenaphthylene | 63 | U | 390 | 63 | ug/Kg |
| 606-20-2 | 2,6-Dinitrotoluene | 55 | U | 390 | 55 | ug/Kg |

U = Not Detected
 RL = Reporting Limit
 MDL = Method Detection Limit
 E = Value Exceeds Calibration Range

J = Estimated Value
 B = Analyte Found In Associated Method Blank
 N = Presumptive Evidence of a Compound

Report of Analysis

| | | | |
|---------------------------|--------------------------------|------------------------|------------|
| Client: | Shaw E & I | Date Collected: | 12/16/2006 |
| Project: | Textron | Date Received: | 2/6/2007 |
| Client Sample ID: | SB-219A ^{2,3} (29-29) | SDG No.: | Y1443 |
| Lab Sample ID: | Y1443-02 | Matrix: | SOIL |
| Analytical Method: | 8270 | % Moisture: | 15 |
| Sample Wt/Wol: | 30.1 g | Extract Vol: | 1000 uL |

| | | | | |
|----------------|-----------------|-----------------------|----------------------|----------------------------|
| File ID | Dilution | Date Extracted | Date Analyzed | Analytical Batch ID |
| BB036131.D | 1 | 2/8/2007 | 2/8/2007 | BB012907 |

| CAS Number | Parameter | Conc. | Qualifier | RL | MDL | Units |
|----------------|----------------------------|-------|-----------|-----|-----|-------|
| TARGETS | | | | | | |
| 99-09-2 | 3-Nitroaniline | 51 | U | 970 | 51 | ug/Kg |
| 83-32-9 | Acenaphthene | 69 | U | 390 | 69 | ug/Kg |
| 51-28-5 | 2,4-Dinitrophenol | 330 | U | 970 | 330 | ug/Kg |
| 100-02-7 | 4-Nitrophenol | 48 | U | 970 | 48 | ug/Kg |
| 132-64-9 | Dibenzofuran | 64 | U | 390 | 64 | ug/Kg |
| 121-14-2 | 2,4-Dinitrotoluene | 57 | U | 390 | 57 | ug/Kg |
| 84-66-2 | Diethylphthalate | 67 | U | 390 | 67 | ug/Kg |
| 7005-72-3 | 4-Chlorophenyl-phenylether | 61 | U | 390 | 61 | ug/Kg |
| 86-73-7 | Fluorene | 65 | U | 390 | 65 | ug/Kg |
| 100-01-6 | 4-Nitroaniline | 66 | U | 970 | 66 | ug/Kg |
| 534-52-1 | 4,6-Dinitro-2-methylphenol | 75 | U | 970 | 75 | ug/Kg |
| 86-30-6 | N-Nitrosodiphenylamine | 64 | U | 390 | 64 | ug/Kg |
| 101-55-3 | 4-Bromophenyl-phenylether | 58 | U | 390 | 58 | ug/Kg |
| 118-74-1 | Hexachlorobenzene | 62 | U | 390 | 62 | ug/Kg |
| 87-86-5 | Pentachlorophenol | 90 | U | 970 | 90 | ug/Kg |
| 85-01-8 | Phenanthrene | 62 | U | 390 | 62 | ug/Kg |
| 120-12-7 | Anthracene | 59 | U | 390 | 59 | ug/Kg |
| 86-74-8 | Carbazole | 59 | U | 390 | 59 | ug/Kg |
| 84-74-2 | Di-n-butylphthalate | 59 | U | 390 | 59 | ug/Kg |
| 206-44-0 | Fluoranthene | 58 | U | 390 | 58 | ug/Kg |
| 129-00-0 | Pyrene | 69 | U | 390 | 69 | ug/Kg |
| 85-68-7 | Butylbenzylphthalate | 63 | U | 390 | 63 | ug/Kg |
| 91-94-1 | 3,3-Dichlorobenzidine | 66 | U | 390 | 66 | ug/Kg |
| 56-55-3 | Benzo(a)anthracene | 54 | U | 390 | 54 | ug/Kg |
| 218-01-9 | Chrysene | 70 | U | 390 | 70 | ug/Kg |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | 130 | J | 390 | 75 | ug/Kg |
| 117-84-0 | Di-n-octyl phthalate | 66 | U | 390 | 66 | ug/Kg |
| 205-99-2 | Benzo(b)fluoranthene | 43 | U | 390 | 43 | ug/Kg |
| 207-08-9 | Benzo(k)fluoranthene | 85 | U | 390 | 85 | ug/Kg |
| 50-32-8 | Benzo(a)pyrene | 62 | U | 390 | 62 | ug/Kg |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 49 | U | 390 | 49 | ug/Kg |

U = Not Detected
 RL = Reporting Limit
 MDL = Method Detection Limit
 E = Value Exceeds Calibration Range

J = Estimated Value
 B = Analyte Found In Associated Method Blank
 N = Presumptive Evidence of a Compound

Report of Analysis

| | | | |
|---------------------------|----------------|------------------------|------------|
| Client: | Shaw E & I | Date Collected: | 12/16/2006 |
| Project: | Textron | Date Received: | 2/6/2007 |
| Client Sample ID: | SB-219A(29-29) | SDG No.: | Y1443 |
| Lab Sample ID: | Y1443-02 | Matrix: | SOIL |
| Analytical Method: | 8270 | % Moisture: | 15 |
| Sample Wt/Wol: | 30.1 g | Extract Vol: | 1000 uL |

| | | | | |
|----------------|-----------------|-----------------------|----------------------|----------------------------|
| File ID | Dilution | Date Extracted | Date Analyzed | Analytical Batch ID |
| BB036131.D | 1 | 2/8/2007 | 2/8/2007 | BB012907 |

| CAS Number | Parameter | Conc. | Qualifier | RL | MDL | Units |
|---------------------------------------|------------------------------------|---------|-----------|----------|-----|---------|
| TARGETS | | | | | | |
| 53-70-3 | Dibenz(a,h)anthracene | 49 | U | 390 | 49 | ug/Kg |
| 191-24-2 | Benzo(g,h,i)perylene | 64 | U | 390 | 64 | ug/Kg |
| SURROGATES | | | | | | |
| 367-12-4 | 2-Fluorophenol | 98.48 | 66 % | 25 - 121 | | SPK: 15 |
| 13127-88-3 | Phenol-d5 | 104.55 | 70 % | 24 - 113 | | SPK: 15 |
| 4165-60-0 | Nitrobenzene-d5 | 74.62 | 75 % | 23 - 120 | | SPK: 10 |
| 321-60-8 | 2-Fluorobiphenyl | 69.49 | 69 % | 30 - 116 | | SPK: 10 |
| 118-79-6 | 2,4,6-Tribromophenol | 101.47 | 68 % | 19 - 122 | | SPK: 15 |
| 1718-51-0 | Terphenyl-d14 | 91.79 | 92 % | 18 - 137 | | SPK: 10 |
| INTERNAL STANDARDS | | | | | | |
| 3855-82-1 | 1,4-Dichlorobenzene-d4 | 268737 | 6.44 | | | |
| 1146-65-2 | Naphthalene-d8 | 1141981 | 8.74 | | | |
| 15067-26-2 | Acenaphthene-d10 | 596432 | 12.20 | | | |
| 1517-22-2 | Phenanthrene-d10 | 968614 | 15.19 | | | |
| 1719-03-5 | Chrysene-d12 | 617230 | 20.52 | | | |
| 1520-96-3 | Perylene-d12 | 697419 | 23.76 | | | |
| TENTITIVE IDENTIFIED COMPOUNDS | | | | | | |
| | ACP3.97 | 5800 | AB | 3.97 | | ug/Kg |
| 67-72-1 | Ethane, hexachloro- | 600 | J | 7.32 | | ug/Kg |
| 88104-31-8 | 2- Chloropropionic acid, octadecyl | 150 | J | 15.81 | | ug/Kg |
| 5454-48-8 | Bromoacetic acid, hexadecyl ester | 250 | J | 17.47 | | ug/Kg |
| 54833-48-6 | Heptadecane, 2,6,10,15-tetramethy | 150 | J | 17.53 | | ug/Kg |
| 629-78-7 | Heptadecane | 540 | J | 18.28 | | ug/Kg |
| 638-67-5 | Tricosane | 1500 | J | 19.00 | | ug/Kg |
| 646-31-1 | Tetracosane | 2600 | J | 19.69 | | ug/Kg |
| 593-45-3 | Octadecane | 2900 | J | 20.35 | | ug/Kg |
| 13187-99-0 | 2-Bromo dodecane | 1400 | J | 20.99 | | ug/Kg |
| 55282-13-8 | Octadecane, 5,14-dibutyl- | 410 | J | 21.63 | | ug/Kg |
| 7683-64-9 | Squalene | 380 | J | 22.51 | | ug/Kg |

U = Not Detected
 RL = Reporting Limit
 MDL = Method Detection Limit
 E = Value Exceeds Calibration Range

J = Estimated Value
 B = Analyte Found in Associated Method Blank
 N = Presumptive Evidence of a Compound

Hit Summary Report

SDG No.: Y1443

Order ID: Y1443

Client: Shaw E & I

Project ID: Textron

Test: SVOC-TCL BNA -20

| Sample ID | Client ID | Matrix | Parameter | Concentration | C | RDL | MDL | Units |
|-------------------------|-----------|--------|----------------------------|---------------|---|-----|-----|-------|
| Client ID: | MW-101D | | | | | | | |
| Y1443-01 | MW-101D | WATER | bis(2-Ethylhexyl)phthalate | 1.7 | J | 10 | 1.6 | ug/L |
| Y1443-01 | MW-101D | WATER | Tetrachloroethylene | * 260 | J | 0 | 0 | ug/L |
| Y1443-01 | MW-101D | WATER | ACP3.00 | * 46 | A | 0 | 0 | ug/L |
| Total SVOC's: | | | | 1.70 | | | | |
| Total TIC's: | | | | 306.00 | | | | |
| Total SVOC's and TIC's: | | | | 307.70 | | | | |

Client ID: SB-219A(29-29)

| | | | | | | | | |
|-------------------------|----------------|------|--------------------------------|----------|----|-----|----|-------|
| Y1443-02 | SB-219A(29-29) | SOIL | Phenol | 240 | J | 390 | 59 | ug/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | bis(2-Ethylhexyl)phthalate | 130 | J | 390 | 75 | ug/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | ACP3.97 | * 5800 | AB | 0 | 0 | ug/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | Ethane, hexachloro- | * 600 | J | 0 | 0 | ug/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | 2- Chloropropionic acid, octad | * 150 | J | 0 | 0 | ug/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | Bromoacetic acid, hexadecyl e | * 250 | J | 0 | 0 | ug/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | Heptadecane, 2,6,10,15-tetra | * 150 | J | 0 | 0 | ug/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | Heptadecane | * 540 | J | 0 | 0 | ug/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | Tricosane | * 1500 | J | 0 | 0 | ug/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | Tetracosane | * 2600 | J | 0 | 0 | ug/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | Octadecane | * 2900 | J | 0 | 0 | ug/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | 2-Bromo dodecane | * 1400 | J | 0 | 0 | ug/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | Octadecane, 5,14-dibutyl- | * 410 | J | 0 | 0 | ug/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | Squalene | * 380 | J | 0 | 0 | ug/Kg |
| Total SVOC's: | | | | 370.00 | | | | |
| Total TIC's: | | | | 16680.00 | | | | |
| Total SVOC's and TIC's: | | | | 17050.00 | | | | |

Note: The asterisk "*" flag next to a parameter signifies a TIC parameter.

CHEMTECH

Lab Chronicle

Order ID: Y1443
Client: Shaw E & I
Contact: Randi Rothmel

Order Date: 2/6/2007 2:10:55 PM
Project: Textron
Location: B63

| Lab ID | Client ID | Matrix | Test | Method | Sample Date | PrepDate | AnalDate | Received |
|----------|-----------------------------|--------|------------------|--------|-------------|----------|----------|----------|
| Y1443-01 | MW-101D | WATER | SVOC-TCL BNA -20 | 8270 | 12/16/06 | 02/08/07 | 02/09/07 | 02/06/07 |
| Y1443-02 | SB-219A(2 ⁷ -29) | SOIL | SVOC-TCL BNA -20 | 8270 | 12/16/06 | 02/08/07 | 02/08/07 | 02/06/07 |

8/9
2/21

Report of Analysis

Client: Shaw E & I

Date Collected: 12/16/2006

Project: Textron

Date Received: 2/6/2007

Client Sample ID: MW-101D

SDG No.: Y1443

Lab Sample ID: Y1443-01

Matrix: WATER

% Solids: 0.00

| CAS No. | Analyte | Conc. | Qualifier | Units | DL | Dilution | Date Prep | Date Anal. | Method |
|-----------|-----------|--------|-----------|-------|-------|----------|-----------|------------|-----------------|
| 7440-36-0 | Antimony | 6.940 | J | ug/L | 3.300 | 1 | 2/8/2007 | 2/12/2007 | EPA SW-846 6010 |
| 7440-38-2 | Arsenic | 5.000 | U | ug/L | 5.000 | 1 | 2/8/2007 | 2/12/2007 | EPA SW-846 6010 |
| 7440-41-7 | Beryllium | 0.100 | U | ug/L | 0.100 | 1 | 2/8/2007 | 2/12/2007 | EPA SW-846 6010 |
| 7440-43-9 | Cadmium | 2.500 | J | ug/L | 0.500 | 1 | 2/8/2007 | 2/12/2007 | EPA SW-846 6010 |
| 7440-47-3 | Chromium | 3.000 | U | ug/L | 3.000 | 1 | 2/8/2007 | 2/12/2007 | EPA SW-846 6010 |
| 7440-50-8 | Copper | 3.640 | U | ug/L | 3.640 | 1 | 2/8/2007 | 2/12/2007 | EPA SW-846 6010 |
| 7439-92-1 | Lead | 4.380 | J | ug/L | 2.000 | 1 | 2/8/2007 | 2/12/2007 | EPA SW-846 6010 |
| 7439-97-6 | Mercury | 0.1100 | U | ug/L | 0.110 | 1 | 2/13/2007 | 2/13/2007 | EPA SW-846 7470 |
| 7440-02-0 | Nickel | 1.800 | U | ug/L | 1.800 | 1 | 2/8/2007 | 2/12/2007 | EPA SW-846 6010 |
| 7782-49-2 | Selenium | 2.200 | U | ug/L | 2.200 | 1 | 2/8/2007 | 2/12/2007 | EPA SW-846 6010 |
| 7440-22-4 | Silver | 1.100 | U | ug/L | 1.100 | 1 | 2/8/2007 | 2/12/2007 | EPA SW-846 6010 |
| 7440-28-0 | Thallium | 4.300 | U | ug/L | 4.300 | 1 | 2/8/2007 | 2/12/2007 | EPA SW-846 6010 |
| 7440-66-6 | Zinc | 94.9 | | ug/L | 1.800 | 1 | 2/8/2007 | 2/12/2007 | EPA SW-846 6010 |

Comments:

U = Not Detected
DL = Method Detection Limit or Instrument Detection Limit

J = Estimated Value
B = Analyte Found In Associated Method Blank
N = Spiked sample recovery not within control limits

Report of Analysis

Client: Shaw E & I

Date Collected: 12/16/2006

Project: Textron

Date Received: 2/6/2007

Client Sample ID: SB-219A²⁷(29-29)

SDG No.: Y1443

Lab Sample ID: Y1443-02

Matrix: SOIL

% Solids: 85.30

| CAS No. | Analyte | Conc. | Qualifier | Units | DL | Dilution | Date Prep | Date Anal. | Method |
|-----------|-----------|-------|-----------|-------|-------|----------|-----------|------------|-----------------|
| 7440-36-0 | Antimony | 0.40 | U | mg/Kg | 0.40 | 1 | 2/8/2007 | 2/8/2007 | EPA SW-846 6010 |
| 7440-38-2 | Arsenic | 2.3 | U | mg/Kg | 2.3 | 1 | 2/8/2007 | 2/8/2007 | EPA SW-846 6010 |
| 7440-41-7 | Beryllium | 0.15 | J | mg/Kg | 0.01 | 1 | 2/8/2007 | 2/8/2007 | EPA SW-846 6010 |
| 7440-43-9 | Cadmium | 0.29 | U | mg/Kg | 0.29 | 1 | 2/8/2007 | 2/8/2007 | EPA SW-846 6010 |
| 7440-47-3 | Chromium | 4.8 | | mg/Kg | 0.13 | 1 | 2/8/2007 | 2/8/2007 | EPA SW-846 6010 |
| 7440-50-8 | Copper | 9.1 | | mg/Kg | 0.15 | 1 | 2/8/2007 | 2/8/2007 | EPA SW-846 6010 |
| 7439-92-1 | Lead | 15.8 | | mg/Kg | 0.16 | 1 | 2/8/2007 | 2/8/2007 | EPA SW-846 6010 |
| 7439-97-6 | Mercury | 0.007 | U | mg/Kg | 0.007 | 1 | 2/8/2007 | 2/8/2007 | EPA SW-846 7471 |
| 7440-02-0 | Nickel | 6.8 | | mg/Kg | 0.14 | 1 | 2/8/2007 | 2/8/2007 | EPA SW-846 6010 |
| 7782-49-2 | Selenium | 0.21 | U | mg/Kg | 0.21 | 1 | 2/8/2007 | 2/8/2007 | EPA SW-846 6010 |
| 7440-22-4 | Silver | 0.21 | U | mg/Kg | 0.21 | 1 | 2/8/2007 | 2/8/2007 | EPA SW-846 6010 |
| 7440-28-0 | Thallium | 0.62 | U | mg/Kg | 0.62 | 1 | 2/8/2007 | 2/8/2007 | EPA SW-846 6010 |
| 7440-66-6 | Zinc | 19.9 | N | mg/Kg | 0.15 | 1 | 2/8/2007 | 2/8/2007 | EPA SW-846 6010 |

Comments:

U = Not Detected
DL = Method Detection Limit or Instrument Detection Limit

J = Estimated Value
B = Analyte Found In Associated Method Blank
N = Spiked sample recovery not within control limits

Chemtech Consulting Group

Hit Summary Sheet SW-846

SDG No.: Y1443

Order ID: Y1443

Client: Shaw E & I

Project ID: Textron

| Sample ID | Client ID | Matrix | Parameter | Concentration | C | RDL | MDL | Units |
|------------|-----------------------|--------|-----------|---------------|---|-------|-------|-------|
| Client ID: | MW-101D | | | | | | | |
| Y1443-01 | MW-101D | WATER | Antimony | 6.940 | J | 25.0 | 3.300 | ug/L |
| Y1443-01 | MW-101D | WATER | Cadmium | 2.500 | J | 3.000 | 0.500 | ug/L |
| Y1443-01 | MW-101D | WATER | Lead | 4.380 | J | 6.000 | 2.000 | ug/L |
| Y1443-01 | MW-101D | WATER | Zinc | 94.9 | | 20.0 | 1.800 | ug/L |
| Client ID: | SB-219A(29-29) | | | | | | | |
| Y1443-02 | SB-219A(29-29) | SOIL | Beryllium | 0.15 | J | 0.35 | 0.01 | mg/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | Chromium | 4.8 | | 0.59 | 0.13 | mg/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | Copper | 9.1 | | 1.2 | 0.15 | mg/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | Lead | 15.8 | | 0.70 | 0.16 | mg/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | Nickel | 6.8 | | 2.3 | 0.14 | mg/Kg |
| Y1443-02 | SB-219A(29-29) | SOIL | Zinc | 19.9 | | 2.3 | 0.15 | mg/Kg |

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CHEMTECH

Lab Chronicle

Order ID: Y1443 Order Date: 2/6/2007 2:10:55 PM
Client: Shaw E & I Project: Textron
Contact: Randi Rothmel Location: B63

| Lab ID | Client ID | Matrix | Test | Method | Sample Date | PrepDate | AnalDate | Received |
|----------|----------------|--------|--|--------------|-------------|----------------------|----------------------|----------|
| Y1443-01 | MW-101D | WATER | <u>Mercury</u> <u>Metals ICP-PP</u> | 7470 6010 | 12/16/06 | 02/13/07 02/08/07 | 02/13/07 02/12/07 | 02/06/07 |
| Y1443-02 | SB-219A(29-29) | SOIL | <u>Mercury</u> <u>Metals ICP-PP</u> | 7471 6010 | 12/16/06 | 02/08/07 02/08/07 | 02/08/07 02/08/07 | 02/06/07 |

04/2/07

CHEMTECH

284 Sheffield Street Mountainside, NJ 07092

Tel . (908) 789-8900 Fax (908) 789-8922

END OF ANALYTICAL RESULTS

APPENDIX C

Laboratory Analytical Report: Groundwater Sampling



March 01, 2007

ANALYTICAL TEST RESULTS

Ed VanDoren
SHAW E & I, Inc.
11 Northeastern Boulevard
Salem, NH 030791953
TEL: (603) 870-4500
FAX: (603) 870-4501

Subject: 101960 Textron Gorham

Workorder No.: 0702072

Dear Ed VanDoren:

AMRO Environmental Laboratories Corp. received 24 samples on 2/19/07 for the analyses presented in the following report.

AMRO is accredited in accordance with NELAC and certifies that these test results meet all the requirements of NELAC, where applicable, unless otherwise noted in the case narrative.

The enclosed Sample Receipt Checklist details the condition of your sample(s) upon receipt. Please be advised that any unused sample volume and sample extracts will be stored for a period of 60 days from sample receipt date (90 days for samples from New York). After this time, AMRO will properly dispose of the remaining sample(s). If you require further analysis, or need the samples held for a longer period, please contact us immediately.

This report consists of a total of 90 pages. This letter is an integral part of your data report. All results in this project relate only to the sample(s) as received by the laboratory and documented in the Chain-of-Custody. This report shall not be reproduced except in full, without the written approval of the laboratory. If you have any questions regarding this project in the future, please refer to the Workorder Number above.

Sincerely,

Nancy Stewart
Vice President

State Certifications: NH (NELAC): 1001, MA: M-NH012, CT: PH-0758, NY: 11278 (NELAC), ME: NH012 and 1001, NJ: NH125, RI: 00105, U.S. Army Corps of Engineers (USACE), Naval Facilities Engineering Service Center (NFESC).

Hard copy of the State Certification is available upon request.



CLIENT: SHAW E & I, Inc.
Project: 101960 Textron Gorham
Lab Order: 0702072
Date Received: 2/19/07

Work Order Sample Summary

| Lab Sample ID | Client Sample ID | Collection Date | Collection Time |
|----------------------|-------------------------|------------------------|------------------------|
| 0702072-01A | MW-207S | 2/17/07 | 7:00 AM |
| 0702072-02A | MW-207D | 2/17/07 | 7:30 AM |
| 0702072-03A | MW-202D | 2/17/07 | 8:00 AM |
| 0702072-04A | MW-202S | 2/17/07 | 8:30 AM |
| 0702072-05A | MW-101S | 2/17/07 | 9:00 AM |
| 0702072-06A | MW-101S DUP | 2/17/07 | 9:00 AM |
| 0702072-07A | MW-101D | 2/17/07 | 9:30 AM |
| 0702072-08A | MW-201D | 2/17/07 | 10:00 AM |
| 0702072-09A | MW-216S | 2/17/07 | 10:30 AM |
| 0702072-10A | MW-216D | 2/17/07 | 11:00 AM |
| 0702072-11A | MW-217D | 2/17/07 | 11:30 AM |
| 0702072-12A | MW-217S | 2/17/07 | 12:00 PM |
| 0702072-13A | MW-218S | 2/17/07 | 12:30 PM |
| 0702072-14A | MW-218D | 2/17/07 | 1:00 PM |
| 0702072-15A | MW-209D | 2/17/07 | 1:30 PM |
| 0702072-16A | MW-112 | 2/17/07 | 2:00 PM |
| 0702072-17A | MW-116S | 2/17/07 | 2:30 PM |
| 0702072-18A | MW-116D | 2/17/07 | 3:00 PM |
| 0702072-19A | CW-1 | 2/17/07 | 12:15 PM |
| 0702072-20A | CW-2 | 2/17/07 | 12:45 PM |
| 0702072-21A | GZA-5 | 2/17/07 | 2:30 PM |
| 0702072-21B | GZA-5 | 2/17/07 | 2:30 PM |
| 0702072-22A | GZA-5 DUP | 2/17/07 | 2:30 PM |
| 0702072-23A | CW-6 | 2/17/07 | 1:45 PM |
| 0702072-24A | CW-6 DUP | 2/17/07 | 1:45 PM |

AMRO Environmental Laboratories Corp.

01-Mar-07

DATES REPORT

Lab Order: 0702072

Client: SHAW E & I, Inc.

Project: 101960 Textron Gorham

| Sample ID | Client Sample ID | Collection Date | Matrix | Analytical Test Name Preparatory Test Name | Prep Date | Analysis Date Batch ID | TCLP Date |
|-------------|------------------|--------------------|-------------|---|-----------|---------------------------|-----------|
| 0702072-01A | MW-207S | 2/17/07 7:00:00 AM | Groundwater | EPA 8260B VOLATILES by GC/MS EPA 5030B | 2/17/07 | 2/21/07 R35962 | |
| 0702072-02A | MW-207D | 2/17/07 7:30:00 AM | | EPA 8260B VOLATILES by GC/MS | 2/17/07 | 2/21/07 R35962 | |
| | | | | EPA 8260B VOLATILES by GC/MS | 2/23/07 | 2/23/07 R35993 | |
| 0702072-03A | MW-202D | 2/17/07 8:00:00 AM | | EPA 8260B VOLATILES by GC/MS | 2/17/07 | 2/21/07 R35962 | |
| 0702072-04A | MW-202S | 2/17/07 8:30:00 AM | | EPA 8260B VOLATILES by GC/MS | 2/17/07 | 2/21/07 R35962 | |
| | | | | EPA 8260B VOLATILES by GC/MS | 2/22/07 | 2/22/07 R35964 | |
| 0702072-05A | MW-101S | 2/17/07 9:00:00 AM | | EPA 8260B VOLATILES by GC/MS | 2/17/07 | 2/21/07 R35962 | |
| | | | | EPA 8260B VOLATILES by GC/MS | 2/17/07 | 2/22/07 R35964 | |
| 0702072-06A | MW-101S DUP | | | EPA 8260B VOLATILES by GC/MS | 2/17/07 | 2/21/07 R35962 | |
| | | | | EPA 8260B VOLATILES by GC/MS | 2/17/07 | 2/22/07 R35964 | |
| 0702072-07A | MW-101D | 2/17/07 9:30:00 AM | | EPA 8260B VOLATILES by GC/MS | 2/17/07 | 2/22/07 R35964 | |
| | | | | EPA 8260B VOLATILES by GC/MS | 2/17/07 | 2/22/07 R35964 | |

AMRO Environmental Laboratories Corp.

01-Mar-07

DATES REPORT

Lab Order: 0702072

Client: SHAW E & I, Inc.

Project: 101960 Textron Gorham

| Sample ID | Client Sample ID | Collection Date | Matrix | Analytical Test Name | Preparatory Test Name | Prep Date | Batch ID | Analysis Date | TCLP Date |
|-------------|------------------|---------------------|-------------|------------------------------|-----------------------|-----------|----------|---------------|-----------|
| 0702072-08A | MW-201D | 2/17/07 10:00:00 AM | Groundwater | EPA 8260B VOLATILES by GC/MS | EPA 5030B | 2/17/07 | R35964 | 2/22/07 | |
| 0702072-09A | MW-216S | 2/17/07 10:30:00 AM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | R35964 | 2/22/07 | |
| 0702072-10A | MW-216D | 2/17/07 11:00:00 AM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | R35964 | 2/22/07 | |
| 0702072-11A | MW-217D | 2/17/07 11:30:00 AM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | R35964 | 2/22/07 | |
| 0702072-12A | MW-217S | 2/17/07 12:00:00 PM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | R35964 | 2/22/07 | |
| 0702072-13A | MW-218S | 2/17/07 12:30:00 PM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | R35964 | 2/22/07 | |
| 0702072-14A | MW-218D | 2/17/07 1:00:00 PM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | R35964 | 2/22/07 | |
| 0702072-15A | MW-209D | 2/17/07 1:30:00 PM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | R35964 | 2/22/07 | |
| 0702072-16A | MW-112 | 2/17/07 2:00:00 PM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | R35993 | 2/23/07 | |
| 0702072-17A | MW-116S | 2/17/07 2:30:00 PM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | R35993 | 2/23/07 | |
| 0702072-18A | MW-116D | 2/17/07 3:00:00 PM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | R35993 | 2/23/07 | |
| 0702072-19A | CW-1 | 2/17/07 12:15:00 PM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | R35993 | 2/23/07 | |

AMRO Environmental Laboratories Corp.

01-Mar-07

Lab Order: 0702072

Client: SHAW E & I, Inc.

Project: 101960 Textron Gorham

DATES REPORT

| Sample ID | Client Sample ID | Collection Date | Matrix | Analytical Test Name | Preparatory Test Name | Prep Date | Analysis Date | Batch ID | TCLP Date |
|-------------|------------------|---------------------|-------------|---|-----------------------|-----------|---------------|----------|-----------|
| 0702072-20A | CW-2 | 2/17/07 12:45:00 PM | Groundwater | EPA 8260B VOLATILES by GC/MS EPA 5030B | | 2/17/07 | 2/23/07 | R35993 | |
| 0702072-23A | CW-6 | 2/17/07 1:45:00 PM | | TPH by GC/FID (modified 8015B) AQPREP SEP FUNNEL: FING | | 2/23/07 | 2/23/07 | 16813 | |
| 0702072-24A | CW-6 DUP | | | TPH by GC/FID (modified 8015B) | | 2/23/07 | 2/23/07 | 16813 | |

| | | | | | | | | | |
|---|--------------|---------------------------------|---|---------------------------------|---|--|---|--|--|
| Project No: 101960 | | Project Name: Textron Gorham | | Project Manager: Ed Vandoren | | Samples/Signature: <i>[Signature]</i> | | AMRO Project No.: 0702072 | |
| P.O.#: 157431 | | Results Needed by: | | Project State: RI | | Requested Analyses | | Remarks | |
| Seal Intact? Yes No N/A | | Date/Time Sampled | | Matrix | | Total # of Cont. & Size | | Requested Analyses | |
| Sample ID: | | Date/Time Sampled | | Matrix | | Total # of Cont. & Size | | Requested Analyses | |
| MW-2075 | 2/17/07 0700 | GW | 2 | 2 | ✓ | 2 | 2 | 2 | EDD to: Catherine.Joe@shawgrp.com and PDF of report. |
| MW-207D | 2/17/07 0730 | | 2 | 2 | ✓ | 2 | 2 | 2 | |
| MW-208D | 2/17/07 0800 | | 2 | 2 | ✓ | 2 | 2 | 2 | |
| MW-202S | 2/17/07 0830 | | 2 | 2 | ✓ | 2 | 2 | 2 | |
| MW-101S | 2/17/07 0900 | | 2 | 2 | ✓ | 2 | 2 | 2 | |
| MW-101S DUP | 2/17/07 0900 | | 2 | 2 | ✓ | 2 | 2 | 2 | |
| MW-101D | 2/17/07 0930 | | 2 | 2 | ✓ | 2 | 2 | 2 | |
| MW-201D | 2/17/07 1000 | | 2 | 2 | ✓ | 2 | 2 | 2 | |
| MW-216S | 2/17/07 1030 | | 2 | 2 | ✓ | 2 | 2 | 2 | |
| MW-216D | 2/17/07 1100 | ✓ | 2 | 2 | ✓ | 2 | 2 | 2 | * MW18# 1100 |
| Preservative: Cl-HCl, MeOH, N-HNO3, S-H2SO4, Na-NaOH, O- Other | | | | | | | | | |
| Priority Turnaround Time Authorization | | | | | | | | | |
| Before submitting samples for expedited TAT, you must have a coded AUTHORIZATION NUMBER | | | | | | | | | |
| AUTHORIZATION No.: BY: | | | | | | | | | |
| PHONE #: 603-870-4530 FAX #: 603-870-4501 | | | | | | | | | |
| E-mail: Edward.Vandoren@shawgrp.com | | | | | | | | | |
| Relinquished By: | | Date/Time: | | Received By: | | MCP Presumptive Certainty Required? | | MCP Methods Needed: | |
| <i>[Signature]</i> | | 2/19/07 10:00 | | <i>[Signature]</i> | | YES <input type="checkbox"/> NO <input type="checkbox"/> | | YES <input type="checkbox"/> NO <input type="checkbox"/> | |
| <i>[Signature]</i> | | 2/19/07 13:00 | | <i>[Signature]</i> | | AMRO report package level needed: | | AMRO report package level needed: | |
| | | | | | | YES <input type="checkbox"/> NO <input type="checkbox"/> | | S-1 <input type="checkbox"/> GW-1 <input type="checkbox"/> | |
| | | | | | | EDD required: | | S-2 <input type="checkbox"/> GW-2 <input type="checkbox"/> | |
| | | | | | | GISKEY | | S-3 <input type="checkbox"/> GW-3 <input type="checkbox"/> | |
| | | | | | | AMRO policy requires notification in writing to the laboratory in cases where the samples were collected from highly contaminated sites. | | Other: <input type="checkbox"/> | |
| | | | | | | AMROCC2004, Rev.3 08/18/04 | | KNOWN SITE CONTAMINATION: | |

White: Lab Copy

Yellow: Accompanies Report

Pink: Client Copy

SHEET

OF

AMROCC2004, Rev.3 08/18/04

AMRO Environmental Laboratories Corporation
111 Herrick Street
Merrimack, NH 03054

CHAIN-OF-CUSTODY RECORD

No 50750

Office: (603) 424-2022
Fax: (603) 429-8496
web: www.amrolabs.com

| | | | | | | | | | |
|---|-----------------|------------------------------|--|-------------------------|--|--|--|---|--|
| Project No.: 101960 | | Project Name: Textron Gorham | | Project State: RI | | Project Manager: Ed Vandoren | | AMRO Project No.: 0702012 | |
| P.O.#: 157431 | | Results Needed by: | | Total # of Cont. & Size | | Requested Analyses | | Remarks | |
| QUOTE #: | | Seal Intact? Yes No N/A | | Matrix | | Grab | | Email GISKey formatted EDD to: Catherine Joe@shawgrp.com and PDF of report. | |
| Sample ID.: | | Date/Time Sampled | | GW | | Comp. | | | |
| MW-217D | 2/17/17 1130 | | | | | | | | |
| MW-217S | 2/17/17 1200 | | | | | | | | |
| MW-218S | 2/17/17 1230 | | | | | | | | |
| MW-218D | 2/17/17 1300 | | | | | | | | |
| MW-209D | 2/17/17 1330 | | | | | | | | |
| MW-112 | 2/17/17 1400 | | | | | | | | |
| MW-116S | 2/17/17 1430 | | | | | | | | |
| MW-116D | 2/17/17 1530 | | | | | | | | |
| TRIP BANK | provided by LAB | | | | | | | | |
| Preservative: Cl-HCl, MeOH, N-HNO3, S-H2SO4, Na-NaOH, O-Other | | | | | | | | | |
| Send Results To: Ed Vandoren | | | | | | | | | |
| Shaw Environmental, Inc. | | | | | | | | | |
| 11 Northeastern Blvd. | | | | | | | | | |
| Salem, MA 03079-1953 | | | | | | | | | |
| PHONE #: 603-870-4530 FAX #: 603-870-4501 | | | | | | | | | |
| E-mail: edward.vandoren@shawgrp.com | | | | | | | | | |
| Relinquished By | | Date/Time | | Received By | | MCP Presumptive Certainty Required? | | Required Reporting Limits: | |
| [Signature] | | 2/19/07 12:00 | | [Signature] | | YES <input type="checkbox"/> NO <input type="checkbox"/> | | S-1 <input type="checkbox"/> GW-1 <input type="checkbox"/> | |
| [Signature] | | 2/19/07 13:00 | | [Signature] | | YES <input type="checkbox"/> NO <input type="checkbox"/> | | S-2 <input type="checkbox"/> GW-2 <input type="checkbox"/> | |
| [Signature] | | 2/19/07 13:00 | | [Signature] | | YES <input type="checkbox"/> NO <input type="checkbox"/> | | S-3 <input type="checkbox"/> GW-3 <input type="checkbox"/> | |
| Please print clearly, legibly and completely. Samples can not be logged in and the turnaround time clock will not start until any ambiguities are resolved. | | | | | | | | | |
| White: Lab Copy | | Yellow: Accompanies Report | | Pink: Client Copy | | SHEET | | OF | |
| AMRO policy requires notification in writing to the laboratory in cases where the samples were collected from highly contaminated sites. | | | | | | | | | |
| AMROCC2004, Rev.3 08/18/04 | | | | | | | | | |

AMRO Environmental Laboratories Corporation
 111 Herrick Street
 Merrimack, NH 03054

CHAIN-OF-CUSTODY RECORD

No 50752

Office: (603) 424-2022
 Fax: (603) 429-8496
 web: www.amrolabs.com

| | | | | | | | | | | | |
|--|--------------|------------------------------|----|-------------------------------------|---|--|---|--|---|----------------------------|---------------|
| Project No.: 101960 | | Project Name: Textron Gorham | | Project State: RI | | Project Manager: Edward Vandoren | | Sampler's Signature: <i>Edward Vandoren</i> | | AMRO Project No.: 070207 | |
| P.O.#: 157431 | | Results Needed by: | | Total # of Cont. & Size | | Requested Analyses | | Requested Analyses | | Remarks | |
| QUOTE #: | | Seal Intact? Yes No N/A | | Comp. Grab | | Lead ICP Metals SW846 | | Lead ICP Metals SW846 | | Email GISKEY | |
| Sample ID.: | | Date/Time Sampled | | Matrix | | EPA 8260B (Vol) | | EPA 8260B (Vol) | | EDDMatted | |
| CW-1 | 2/17/1215 | 2 | GW | 2 | 2 | 2 | 2 | 2 | 2 | 2 | EDD and PDF |
| CW-2 | 2/17/1245 | 2 | | 2 | 2 | 2 | 2 | 2 | 2 | 2 | of report to: |
| GZA-5 | 2-17-07 1430 | 3 | | 3 | 3 | 3 | 3 | 3 | 3 | 3 | Catherine Joe |
| GZA-6 | | 3 | | 3 | 3 | 3 | 3 | 3 | 3 | 3 | @shawgrp.com |
| CW-6 | 2/17/1345 | 1 | | 1 | 1 | 1 | 1 | 1 | 1 | 1 | |
| GZA-5 DUP. | 2-17-07 1430 | 1 | | 1 | 1 | 1 | 1 | 1 | 1 | 1 | |
| CW-6 DUP. | 2/17/1345 | 1 | | 1 | 1 | 1 | 1 | 1 | 1 | 1 | |
| <p>Preservative: CI-HCl, MeOH, N-HNO3, S-H2SO4, Na-NaOH, O- Other: Cl N</p> <p>Send Results To: Edward Vandoren Shaw Environmental, Inc. 11 Northeastern Blvd. Salem, NH 03079-1953 PHONE #: 803-870-4530 FAX #: 603-870-4501 E-mail: Edward.Vandoren@shawgrp.com</p> <p>Priority Turnaround Time Authorization: Cl N Before submitting samples for expedited TAT, you must have a coded AUTHORIZATION NUMBER</p> <p>METALS: 8 RCRA <input type="checkbox"/> 13 PP <input type="checkbox"/> 23 TAL <input type="checkbox"/> 14 MCP <input type="checkbox"/> Method: 6010 <input type="checkbox"/> 200.7 <input checked="" type="checkbox"/> Other Metals: Lead <input type="checkbox"/></p> <p>Dissolved Metals Field Filtered? YES <input checked="" type="checkbox"/> NO <input type="checkbox"/></p> <p>MCP Presumptive Certainty Required? YES <input type="checkbox"/> NO <input type="checkbox"/></p> <p>MCP Methods Needed: YES <input type="checkbox"/> NO <input type="checkbox"/> AMRO report package level needed: YES <input type="checkbox"/> NO <input type="checkbox"/> EDD required: GISKEY</p> <p>Required Reporting Limits: S-1 <input type="checkbox"/> GW-1 <input type="checkbox"/> S-2 <input type="checkbox"/> GW-2 <input type="checkbox"/> S-3 <input type="checkbox"/> GW-3 <input type="checkbox"/> Other: GISKEY</p> | | | | | | | | | | | |
| Relinquished By: <i>Edward Vandoren</i> | | Date/Time: 2/19/07 1345 | | Received By: <i>William Swadlow</i> | | AMRO policy requires notification in writing to the laboratory in cases where the samples were collected from highly contaminated sites. | | AMRO policy requires notification in writing to the laboratory in cases where the samples were collected from highly contaminated sites. | | KNOWN SITE CONTAMINATION: | |
| White: Lab Copy | | Yellow: Accompanies Report | | Pink: Client Copy | | SHEET | | OF | | AMROCC2004, Rev.3 08/18/04 | |

Login Account for multiple users

From: VanDoren, Edward [Edward.VanDoren@shawgrp.com]
Sent: Thursday, February 22, 2007 2:13 PM
To: Login Account for multiple users
Subject: RE: Testron Gorham Sampling Date (AMRO 0702072)

Hi Connie-

Please do not analyze GZA-5 or GZA-5 Dup. The wrong well was sampled.

Thanks,
Ed

Edward Van Doren
Client Program Manager
Shaw Environmental, Inc.
11 Northeastern Boulevard
Salem, NH 03079-1953
603.870.4530 direct
603.870.4501 fax
edward.vandoren@shawgrp.com
www.shawgrp.com

From: Login Account for multiple users [mailto:login@amrolabs.com]
Sent: Tuesday, February 20, 2007 11:21 AM
To: VanDoren, Edward
Subject: FW: Testron Gorham Sampling Date (AMRO 0702072)

Hi Ed -

We had a call from Dan Leahy and he let us know that the samples were taken 2/17.

Connie in Receiving

From: Login Account for multiple users
Sent: Monday, February 19, 2007 3:11 PM
To: 'VanDoren, Edward'
Subject: Testron Gorham Sampling Date (AMRO 0702072)

Hi Ed -

We picked up these water samples today and I notice that all of the samples are dated **2/14/07** and the Chain reads **2/17/07**. Can you tell me which is the correct sampling date?

Also, the Trip Blank was found broken in the cooler, still inside the bubble bag, so we will be unable to run it for you.

Thanks for your help!

Connie in Receiving

| | |
|--|---------------------------|
| Client: <u>SHAW ENVIRONMENTAL, INC</u> | AMRO ID: <u>0702072</u> |
| Project Name: <u>TEXTRON GORHAM</u> | Date Rec.: <u>2-19-07</u> |
| Ship via: (circle one) Fed Ex., UPS, <u>(AMRO Courier)</u> | Date Due: <u>2-26-07</u> |
| Hand Del., Other Courier, Other: | |

Items to be Checked Upon Receipt

| Item | Yes | No | NA | Comments |
|---|-----|----|----|--|
| 1. Army Samples received in individual plastic bags? | | | ✓ | |
| 2. Custody Seals present? | | | ✓ | |
| 3. Custody Seals Intact? | | | ✓ | |
| 4. Air Bill included in folder if received? | | | ✓ | |
| 5. Is COC included with samples? | ✓ | | | |
| 6. Is COC signed and dated by client? | ✓ | | | |
| 7. Laboratory receipt temperature. TEMP = 3° Samples rec. with ice <input checked="" type="checkbox"/> ice packs <input type="checkbox"/> neither <input type="checkbox"/> | | | | |
| 8. Were samples received the same day they were sampled? Is client temperature 4°C ± 2°C? If no obtain authorization from the client for the analyses. Client authorization from: Date: Obtained by: | ✓ | ✓ | | |
| 9. Is the COC filled out correctly and completely? | ✓ | | | |
| 10. Does the info on the COC match the samples? | ✓ | ✓ | | ALL DATED 2/14/07 EMAIL TO CLIENT |
| 11. Were samples rec. within holding time? | ✓ | | | |
| 12. Were all samples properly labeled? | ✓ | | | |
| 13. Were all samples properly preserved? | ✓ | | | |
| 14. Were proper sample containers used? | ✓ | | | |
| 15. Were all samples received intact? (none broken or leaking) | ✓ | ✓ | | TRIP BLANK RECEIVED BROKEN - EMAIL TO CLIENT |
| 16. Were VOA vials rec. with no air bubbles? | ✓ | | | |
| 17. Were the sample volumes sufficient for requested analysis? | ✓ | | | |
| 18. Were all samples received? | ✓ | | | |

19. VPH and VOA Soils only:

Sampling Method VPH (circle one): M=Methanol, E=EnCore (air-tight container)

Sampling Method VOA (circle one): M=Methanol, SB=Sodium Bisulfate, E=EnCore, B=Bulk

If M or SB:

Does preservative cover the soil? If NO then client must be faxed.

Does preservation level come close to the fill line on the vial? If NO then client must be faxed.

Were vials provided by AMRO? If NO then weights MUST be obtained from client

Was dry weight aliquot provided? If NO then fax client and inform the VOA lab ASAP.

20. Subcontracted Samples:

What samples sent:

Where sent:

Date:

Analysis:

TAT:

21. Information entered into:

Internal Tracking Log?

Dry Weight Log?

Client Log?

Composite Log?

Filtration Log?

| | | | |
|------------------------|----------------------|----------------------------|----------------------|
| Received By: <u>CC</u> | Date: <u>2-19-07</u> | Logged in By: <u>CC MG</u> | Date: <u>2-19-07</u> |
| Labeled By: <u>CC</u> | Date: <u>2-19-07</u> | Checked By: <u>CC MG</u> | Date: <u>2-21-07</u> |

CLIENT: SHAW E & I, Inc.
Project: 101960 Textron Gorham
Lab Order: 0702072

CASE NARRATIVE

GC/MS VOLATILES:

1. A Matrix Spike (MS) and Matrix Spike Duplicate (MSD) were performed on sample MW-209D (0702072-15A) Batch ID: R35964.

1.1 The % Recovery for 4 analytes out of 67 analytes in the MS was outside the laboratory control limits.

2. A Matrix Spike (MS) and Matrix Spike Duplicate (MSD) were performed on sample CW-1 (0702072-19A) Batch ID: R35993.

2.1 The % Recovery for 2 analytes out of 67 analytes in the MS was outside the laboratory control limits.

2.2 The RPD for 1 analyte out of 67 analytes was outside the laboratory control limits.

DATA COMMENT PAGE

Organic Data Qualifiers

| | |
|----|--|
| ND | Indicates compound was analyzed for, but not detected at or above the reporting limit. |
| J | Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed, or when the data indicates the presence of a compound that meets the identification criteria but the result is less than the sample quantitation limit but greater than the method detection limit. |
| H | Method prescribed holding time exceeded. |
| E | This flag identifies compounds whose concentrations exceed the calibration range of the instrument for that specific analysis. |
| B | This flag is used when the analyte is found in the associated blank as well as in the sample. |
| R | RPD outside accepted recovery limits |
| RL | Reporting limit; defined as the lowest concentration the laboratory can accurately quantitate. |
| S | Spike Recovery outside accepted recovery limits. |
| # | See Case Narrative |

Micro Data Qualifiers

TNTC Too numerous to count

Inorganic Data Qualifiers

| | |
|---------|---|
| ND or U | Indicates element was analyzed for, but not detected at or above the reporting limit. |
| J | Indicates a value greater than or equal to the method detection limit, but less than the quantitation limit. |
| H | Indicates analytical holding time exceedance. |
| B | Indicates that the analyte is found in the associated blank, as well as in the sample. |
| MSA | Indicates value determined by the Method of Standard Addition |
| E | This flag identifies compounds whose concentrations exceed the calibration range of the instrument for that specific analysis. |
| R | RPD outside accepted recovery limits |
| RL | Reporting limit; defined as the lowest concentration the laboratory can accurately quantitate. |
| S | Spike Recovery outside accepted recovery limits. |
| W | Post-digestion spike for Furnace AA analysis is out of control limits (85-115), while sample absorbance is less than 50% of spike absorbance. |
| * | Duplicate analysis not within control limits. |
| + | Indicates the correlation coefficient for the Method of Standard Addition is less than 0.995 |
| # | See Case Narrative |

Report Comments:

1. Soil, sediment and sludge sample results are reported on a "dry weight" basis.
2. Reporting limits are adjusted for sample size used, dilutions and moisture content, if applicable.

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-01A

Client Sample ID: MW-207S
Collection Date: 2/17/07 7:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Vinyl chloride | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,1-Dichloroethene | ND | 100 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| cis-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Trichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Tetrachloroethene | 12,000 | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-01A

Client Sample ID: MW-207S
Collection Date: 2/17/07 7:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|--------------------|
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Surr: Dibromofluoromethane | 100 | 85-116 | | %REC | 100 | 2/21/07 8:31:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 108 | 77-127 | | %REC | 100 | 2/21/07 8:31:00 PM |
| Surr: Toluene-d8 | 101 | 86-114 | | %REC | 100 | 2/21/07 8:31:00 PM |
| Surr: 4-Bromofluorobenzene | 95.9 | 79-117 | | %REC | 100 | 2/21/07 8:31:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-02A

Client Sample ID: MW-207D
Collection Date: 2/17/07 7:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Vinyl chloride | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,1-Dichloroethene | ND | 100 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| cis-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Trichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Tetrachloroethene | 7,700 | 200 | | µg/L | 100 | 2/23/07 2:53:00 PM |
| Tetrachloroethene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-02A

Client Sample ID: MW-207D
Collection Date: 2/17/07 7:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|--------------------|
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Surr: Dibromofluoromethane | 97.3 | 85-116 | | %REC | 100 | 2/21/07 9:05:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 94.9 | 77-127 | | %REC | 100 | 2/21/07 9:05:00 PM |
| Surr: Toluene-d8 | 99.0 | 86-114 | | %REC | 100 | 2/21/07 9:05:00 PM |
| Surr: 4-Bromofluorobenzene | 91.2 | 79-117 | | %REC | 100 | 2/21/07 9:05:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-03A

Client Sample ID: MW-202D
Collection Date: 2/17/07 8:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Vinyl chloride | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,1-Dichloroethene | ND | 100 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| cis-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Trichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Tetrachloroethene | 29,000 | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-03A

Client Sample ID: MW-202D
Collection Date: 2/17/07 8:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|--------------------|
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Surr: Dibromofluoromethane | 102 | 85-116 | | %REC | 100 | 2/21/07 9:40:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 107 | 77-127 | | %REC | 100 | 2/21/07 9:40:00 PM |
| Surr: Toluene-d8 | 101 | 86-114 | | %REC | 100 | 2/21/07 9:40:00 PM |
| Surr: 4-Bromofluorobenzene | 95.9 | 79-117 | | %REC | 100 | 2/21/07 9:40:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-04A

Client Sample ID: MW-202S
Collection Date: 2/17/07 8:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|---------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Vinyl chloride | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,1-Dichloroethene | ND | 100 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| cis-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Trichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Tetrachloroethene | 51,000 | 1,000 | | µg/L | 500 | 2/22/07 3:12:00 PM |
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-04A

Client Sample ID: MW-202S
Collection Date: 2/17/07 8:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|---------------------|
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Surr: Dibromofluoromethane | 104 | 85-116 | | %REC | 100 | 2/21/07 10:14:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 107 | 77-127 | | %REC | 100 | 2/21/07 10:14:00 PM |
| Surr: Toluene-d8 | 100 | 86-114 | | %REC | 100 | 2/21/07 10:14:00 PM |
| Surr: 4-Bromofluorobenzene | 98.8 | 79-117 | | %REC | 100 | 2/21/07 10:14:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-05A

Client Sample ID: MW-101S
Collection Date: 2/17/07 9:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|---------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Vinyl chloride | 520 | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,1-Dichloroethene | ND | 100 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| cis-1,2-Dichloroethene | 510 | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Trichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Tetrachloroethene | 45,000 | 1,000 | | µg/L | 500 | 2/22/07 3:46:00 PM |
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-05A

Client Sample ID: MW-101S
Collection Date: 2/17/07 9:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|---------------------|
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Surr: Dibromofluoromethane | 104 | 85-116 | | %REC | 100 | 2/21/07 10:49:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 109 | 77-127 | | %REC | 100 | 2/21/07 10:49:00 PM |
| Surr: Toluene-d8 | 102 | 86-114 | | %REC | 100 | 2/21/07 10:49:00 PM |
| Surr: 4-Bromofluorobenzene | 96.6 | 79-117 | | %REC | 100 | 2/21/07 10:49:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-06A

Client Sample ID: MW-101S DUP
Collection Date: 2/17/07 9:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|---------------------|
| EPA 8260B VOLATILES BY GC/MS | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Vinyl chloride | 460 | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,1-Dichloroethene | ND | 100 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| cis-1,2-Dichloroethene | 500 | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Trichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Tetrachloroethene | 44,000 | 1,000 | | µg/L | 500 | 2/22/07 4:20:00 PM |
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-06A

Client Sample ID: MW-101S DUP
Collection Date: 2/17/07 9:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|---------------------|
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Surr: Dibromofluoromethane | 101 | 85-116 | | %REC | 100 | 2/21/07 11:23:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 108 | 77-127 | | %REC | 100 | 2/21/07 11:23:00 PM |
| Surr: Toluene-d8 | 102 | 86-114 | | %REC | 100 | 2/21/07 11:23:00 PM |
| Surr: 4-Bromofluorobenzene | 99.6 | 79-117 | | %REC | 100 | 2/21/07 11:23:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
 Lab Order: 0702072
 Project: 101960 Textron Gorham
 Lab ID: 0702072-07A

Client Sample ID: MW-101D
 Collection Date: 2/17/07 9:30:00 AM
 Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------------|-----|---------------------|
| EPA 8260B VOLATILES BY GC/MS | | SW8260B | | Analyst: SK | | |
| Dichlorodifluoromethane | ND | 50 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Chloromethane | ND | 50 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Vinyl chloride | 30 | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Chloroethane | ND | 50 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Bromomethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Trichlorofluoromethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Diethyl ether | ND | 50 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Acetone | ND | 100 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,1-Dichloroethene | 11 | 10 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Carbon disulfide | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Methylene chloride | ND | 50 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Methyl tert-butyl ether | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| trans-1,2-Dichloroethene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,1-Dichloroethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 2-Butanone | ND | 100 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 2,2-Dichloropropane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| cis-1,2-Dichloroethene | 280 | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Chloroform | 20 | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Tetrahydrofuran | ND | 100 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Bromochloromethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,1,1-Trichloroethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,1-Dichloropropene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Carbon tetrachloride | 32 | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2-Dichloroethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Benzene | ND | 10 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Trichloroethene | 160 | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2-Dichloropropane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Bromodichloromethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Dibromomethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 4-Methyl-2-pentanone | ND | 100 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| cis-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Toluene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| trans-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,1,2-Trichloroethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2-Dibromoethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 2-Hexanone | ND | 100 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,3-Dichloropropane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Tetrachloroethene | 19,000 | 200 | | µg/L | 100 | 2/22/07 4:54:00 PM |
| Dibromochloromethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-07A

Client Sample ID: MW-101D
Collection Date: 2/17/07 9:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|---------------------|
| Chlorobenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Ethylbenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| m,p-Xylene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| o-Xylene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Styrene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Bromoform | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Isopropylbenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2,3-Trichloropropane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Bromobenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| n-Propylbenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 2-Chlorotoluene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 4-Chlorotoluene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,3,5-Trimethylbenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| tert-Butylbenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2,4-Trimethylbenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| sec-Butylbenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 4-Isopropyltoluene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,3-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,4-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| n-Butylbenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 50 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2,4-Trichlorobenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Hexachlorobutadiene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Naphthalene | ND | 50 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2,3-Trichlorobenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Surr: Dibromofluoromethane | 102 | 85-116 | | %REC | 10 | 2/21/07 11:58:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 107 | 77-127 | | %REC | 10 | 2/21/07 11:58:00 PM |
| Surr: Toluene-d8 | 101 | 86-114 | | %REC | 10 | 2/21/07 11:58:00 PM |
| Surr: 4-Bromofluorobenzene | 95.3 | 79-117 | | %REC | 10 | 2/21/07 11:58:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-08A

Client Sample ID: MW-201D
Collection Date: 2/17/07 10:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Vinyl chloride | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,1-Dichloroethene | ND | 100 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| cis-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Trichloroethene | 970 | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Tetrachloroethene | 7,600 | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-08A

Client Sample ID: MW-201D
Collection Date: 2/17/07 10:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|--------------------|
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Surr: Dibromofluoromethane | 99.6 | 85-116 | | %REC | 100 | 2/22/07 7:47:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 101 | 77-127 | | %REC | 100 | 2/22/07 7:47:00 PM |
| Surr: Toluene-d8 | 98.0 | 86-114 | | %REC | 100 | 2/22/07 7:47:00 PM |
| Surr: 4-Bromofluorobenzene | 97.4 | 79-117 | | %REC | 100 | 2/22/07 7:47:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
 Lab Order: 0702072
 Project: 101960 Textron Gorham
 Lab ID: 0702072-09A

Client Sample ID: MW-216S
 Collection Date: 2/17/07 10:30:00 AM
 Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|------------------------------|--------|---------|------|-------------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | SW8260B | | Analyst: SK | | |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Trichlorofluoromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| cis-1,2-Dichloroethene | 140 | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Chloroform | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Trichloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Bromodichloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Toluene | 3.4 | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Tetrachloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-09A

Client Sample ID: MW-216S
Collection Date: 2/17/07 10:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Ethylbenzene | 2.6 | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| m,p-Xylene | 7.0 | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| o-Xylene | 8.9 | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,3,5-Trimethylbenzene | 10 | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2,4-Trimethylbenzene | 13 | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 4-Isopropyltoluene | 2.9 | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Naphthalene | 17 | 5.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Surr: Dibromofluoromethane | 102 | 85-116 | | %REC | 1 | 2/22/07 5:28:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 109 | 77-127 | | %REC | 1 | 2/22/07 5:28:00 PM |
| Surr: Toluene-d8 | 102 | 86-114 | | %REC | 1 | 2/22/07 5:28:00 PM |
| Surr: 4-Bromofluorobenzene | 99.4 | 79-117 | | %REC | 1 | 2/22/07 5:28:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-10A

Client Sample ID: MW-216D
Collection Date: 2/17/07 11:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Trichlorofluoromethane | 3.1 | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| cis-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Chloroform | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Trichloroethene | 6.4 | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Bromodichloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Toluene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Tetrachloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-10A

Client Sample ID: MW-216D
Collection Date: 2/17/07 11:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Ethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| m,p-Xylene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| o-Xylene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,3,5-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2,4-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Naphthalene | ND | 5.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Surr: Dibromofluoromethane | 98.6 | 85-116 | | %REC | 1 | 2/22/07 6:03:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 107 | 77-127 | | %REC | 1 | 2/22/07 6:03:00 PM |
| Surr: Toluene-d8 | 99.4 | 86-114 | | %REC | 1 | 2/22/07 6:03:00 PM |
| Surr: 4-Bromofluorobenzene | 100 | 79-117 | | %REC | 1 | 2/22/07 6:03:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-11A

Client Sample ID: MW-217D
Collection Date: 2/17/07 11:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Trichlorofluoromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| cis-1,2-Dichloroethene | 60 | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Chloroform | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Trichloroethene | 75 | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Bromodichloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Toluene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Tetrachloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-11A

Client Sample ID: MW-217D
Collection Date: 2/17/07 11:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Ethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| m,p-Xylene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| o-Xylene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,3,5-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2,4-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Naphthalene | ND | 5.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Surr: Dibromofluoromethane | 100 | 85-116 | | %REC | 1 | 2/22/07 6:38:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 107 | 77-127 | | %REC | 1 | 2/22/07 6:38:00 PM |
| Surr: Toluene-d8 | 100 | 86-114 | | %REC | 1 | 2/22/07 6:38:00 PM |
| Surr: 4-Bromofluorobenzene | 101 | 79-117 | | %REC | 1 | 2/22/07 6:38:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-12A

Client Sample ID: MW-217S
Collection Date: 2/17/07 12:00:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | SW8260B | | Analyst: SK | | |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Trichlorofluoromethane | 2.2 | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| cis-1,2-Dichloroethene | 9.8 | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Chloroform | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Trichloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Bromodichloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Toluene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Tetrachloroethene | 23 | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-12A

Client Sample ID: MW-217S
Collection Date: 2/17/07 12:00:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Ethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| m,p-Xylene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| o-Xylene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,3,5-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2,4-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Naphthalene | ND | 5.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Surr: Dibromofluoromethane | 98.9 | 85-116 | | %REC | 1 | 2/22/07 7:12:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 103 | 77-127 | | %REC | 1 | 2/22/07 7:12:00 PM |
| Surr: Toluene-d8 | 100 | 86-114 | | %REC | 1 | 2/22/07 7:12:00 PM |
| Surr: 4-Bromofluorobenzene | 95.0 | 79-117 | | %REC | 1 | 2/22/07 7:12:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-13A

Client Sample ID: MW-218S
Collection Date: 2/17/07 12:30:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 50 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Chloromethane | ND | 50 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Vinyl chloride | 50 | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Chloroethane | ND | 50 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Bromomethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Trichlorofluoromethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Diethyl ether | ND | 50 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Acetone | ND | 100 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,1-Dichloroethene | ND | 10 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Carbon disulfide | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Methylene chloride | ND | 50 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Methyl tert-butyl ether | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| trans-1,2-Dichloroethene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,1-Dichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 2-Butanone | ND | 100 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 2,2-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| cis-1,2-Dichloroethene | 650 | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Chloroform | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Tetrahydrofuran | ND | 100 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Bromochloromethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,1,1-Trichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,1-Dichloropropene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Carbon tetrachloride | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2-Dichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Benzene | ND | 10 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Trichloroethene | 49 | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Bromodichloromethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Dibromomethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 4-Methyl-2-pentanone | ND | 100 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| cis-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Toluene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| trans-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,1,2-Trichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2-Dibromoethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 2-Hexanone | ND | 100 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,3-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Tetrachloroethene | 370 | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Dibromochloromethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-13A

Client Sample ID: MW-218S
Collection Date: 2/17/07 12:30:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Ethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| m,p-Xylene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| o-Xylene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Styrene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Bromoform | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Isopropylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2,3-Trichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Bromobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| n-Propylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 2-Chlorotoluene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 4-Chlorotoluene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,3,5-Trimethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| tert-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2,4-Trimethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| sec-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 4-Isopropyltoluene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,3-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,4-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| n-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 50 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2,4-Trichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Hexachlorobutadiene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Naphthalene | ND | 50 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2,3-Trichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Surr: Dibromofluoromethane | 103 | 85-116 | | %REC | 10 | 2/22/07 8:21:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 108 | 77-127 | | %REC | 10 | 2/22/07 8:21:00 PM |
| Surr: Toluene-d8 | 103 | 86-114 | | %REC | 10 | 2/22/07 8:21:00 PM |
| Surr: 4-Bromofluorobenzene | 97.1 | 79-117 | | %REC | 10 | 2/22/07 8:21:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-14A

Client Sample ID: MW-218D
Collection Date: 2/17/07 1:00:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 50 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Chloromethane | ND | 50 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Vinyl chloride | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Chloroethane | ND | 50 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Bromomethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Trichlorofluoromethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Diethyl ether | ND | 50 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Acetone | ND | 100 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,1-Dichloroethene | 23 | 10 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Carbon disulfide | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Methylene chloride | ND | 50 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Methyl tert-butyl ether | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| trans-1,2-Dichloroethene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,1-Dichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 2-Butanone | ND | 100 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 2,2-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| cis-1,2-Dichloroethene | 28 | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Chloroform | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Tetrahydrofuran | ND | 100 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Bromochloromethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,1,1-Trichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,1-Dichloropropene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Carbon tetrachloride | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2-Dichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Benzene | ND | 10 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Trichloroethene | 840 | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Bromodichloromethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Dibromomethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 4-Methyl-2-pentanone | ND | 100 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| cis-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Toluene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| trans-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,1,2-Trichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2-Dibromoethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 2-Hexanone | ND | 100 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,3-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Tetrachloroethene | 600 | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Dibromochloromethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-14A

Client Sample ID: MW-218D
Collection Date: 2/17/07 1:00:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Ethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| m,p-Xylene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| o-Xylene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Styrene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Bromoform | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Isopropylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2,3-Trichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Bromobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| n-Propylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 2-Chlorotoluene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 4-Chlorotoluene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,3,5-Trimethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| tert-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2,4-Trimethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| sec-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 4-Isopropyltoluene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,3-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,4-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| n-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 50 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2,4-Trichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Hexachlorobutadiene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Naphthalene | ND | 50 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2,3-Trichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Surr: Dibromofluoromethane | 101 | 85-116 | | %REC | 10 | 2/22/07 8:56:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 106 | 77-127 | | %REC | 10 | 2/22/07 8:56:00 PM |
| Surr: Toluene-d8 | 102 | 86-114 | | %REC | 10 | 2/22/07 8:56:00 PM |
| Surr: 4-Bromofluorobenzene | 97.8 | 79-117 | | %REC | 10 | 2/22/07 8:56:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-15A

Client Sample ID: MW-209D
Collection Date: 2/17/07 1:30:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 50 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Chloromethane | ND | 50 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Vinyl chloride | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Chloroethane | ND | 50 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Bromomethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Trichlorofluoromethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Diethyl ether | ND | 50 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Acetone | ND | 100 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,1-Dichloroethene | ND | 10 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Carbon disulfide | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Methylene chloride | ND | 50 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Methyl tert-butyl ether | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| trans-1,2-Dichloroethene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,1-Dichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 2-Butanone | ND | 100 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 2,2-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| cis-1,2-Dichloroethene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Chloroform | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Tetrahydrofuran | ND | 100 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Bromochloromethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,1,1-Trichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,1-Dichloropropene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Carbon tetrachloride | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2-Dichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Benzene | ND | 10 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Trichloroethene | 110 | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Bromodichloromethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Dibromomethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 4-Methyl-2-pentanone | ND | 100 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| cis-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Toluene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| trans-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,1,2-Trichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2-Dibromoethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 2-Hexanone | ND | 100 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,3-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Tetrachloroethene | 430 | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Dibromochloromethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-15A

Client Sample ID: MW-209D
Collection Date: 2/17/07 1:30:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Ethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| m,p-Xylene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| o-Xylene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Styrene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Bromoform | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Isopropylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2,3-Trichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Bromobenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| n-Propylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 2-Chlorotoluene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 4-Chlorotoluene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,3,5-Trimethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| tert-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2,4-Trimethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| sec-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 4-Isopropyltoluene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,3-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,4-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| n-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 50 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2,4-Trichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Hexachlorobutadiene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Naphthalene | ND | 50 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2,3-Trichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Surr: Dibromofluoromethane | 98.9 | 85-116 | | %REC | 10 | 2/22/07 9:30:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 103 | 77-127 | | %REC | 10 | 2/22/07 9:30:00 PM |
| Surr: Toluene-d8 | 98.4 | 86-114 | | %REC | 10 | 2/22/07 9:30:00 PM |
| Surr: 4-Bromofluorobenzene | 96.8 | 79-117 | | %REC | 10 | 2/22/07 9:30:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-16A

Client Sample ID: MW-112
Collection Date: 2/17/07 2:00:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|--------------------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | SW8260B | | Analyst: SK | | |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Trichlorofluoromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Methyl tert-butyl ether | 17 | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| cis-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Chloroform | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Trichloroethene | 3.9 | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Bromodichloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Toluene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Tetrachloroethene | 42 | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-16A

Client Sample ID: MW-112
Collection Date: 2/17/07 2:00:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Ethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| m,p-Xylene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| o-Xylene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,3,5-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2,4-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Naphthalene | ND | 5.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Surr: Dibromofluoromethane | 97.7 | 85-116 | | %REC | 1 | 2/23/07 3:28:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 103 | 77-127 | | %REC | 1 | 2/23/07 3:28:00 PM |
| Surr: Toluene-d8 | 99.0 | 86-114 | | %REC | 1 | 2/23/07 3:28:00 PM |
| Surr: 4-Bromofluorobenzene | 97.7 | 79-117 | | %REC | 1 | 2/23/07 3:28:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
 Lab Order: 0702072
 Project: 101960 Textron Gorham
 Lab ID: 0702072-17A

Client Sample ID: MW-116S
 Collection Date: 2/17/07 2:30:00 PM
 Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | SW8260B | | Analyst: SK | | |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Trichlorofluoromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| cis-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Chloroform | 32 | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Trichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Bromodichloromethane | 4.2 | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Toluene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Tetrachloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-17A

Client Sample ID: MW-116S
Collection Date: 2/17/07 2:30:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Ethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| m,p-Xylene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| o-Xylene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,3,5-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2,4-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Naphthalene | ND | 5.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Surr: Dibromofluoromethane | 102 | 85-116 | | %REC | 1 | 2/23/07 4:03:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 104 | 77-127 | | %REC | 1 | 2/23/07 4:03:00 PM |
| Surr: Toluene-d8 | 99.4 | 86-114 | | %REC | 1 | 2/23/07 4:03:00 PM |
| Surr: 4-Bromofluorobenzene | 96.5 | 79-117 | | %REC | 1 | 2/23/07 4:03:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-18A

Client Sample ID: MW-116D
Collection Date: 2/17/07 3:00:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Trichlorofluoromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| cis-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Chloroform | 39 | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Trichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Bromodichloromethane | 4.3 | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Toluene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Tetrachloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-18A

Client Sample ID: MW-116D
Collection Date: 2/17/07 3:00:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Ethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| m,p-Xylene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| o-Xylene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,3,5-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2,4-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Naphthalene | ND | 5.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Surr: Dibromofluoromethane | 98.3 | 85-116 | | %REC | 1 | 2/23/07 4:37:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 104 | 77-127 | | %REC | 1 | 2/23/07 4:37:00 PM |
| Surr: Toluene-d8 | 99.8 | 86-114 | | %REC | 1 | 2/23/07 4:37:00 PM |
| Surr: 4-Bromofluorobenzene | 96.8 | 79-117 | | %REC | 1 | 2/23/07 4:37:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc. **Client Sample ID:** CW-1
Lab Order: 0702072 **Collection Date:** 2/17/07 12:15:00 PM
Project: 101960 Textron Gorham **Matrix:** GROUNDWATER
Lab ID: 0702072-19A

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Vinyl chloride | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,1-Dichloroethene | 180 | 100 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| cis-1,2-Dichloroethene | 350 | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Trichloroethene | 5,900 | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Tetrachloroethene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-19A

Client Sample ID: CW-1
Collection Date: 2/17/07 12:15:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|--------------------|
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Surr: Dibromofluoromethane | 100 | 85-116 | | %REC | 100 | 2/23/07 5:46:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 106 | 77-127 | | %REC | 100 | 2/23/07 5:46:00 PM |
| Surr: Toluene-d8 | 101 | 86-114 | | %REC | 100 | 2/23/07 5:46:00 PM |
| Surr: 4-Bromofluorobenzene | 97.2 | 79-117 | | %REC | 100 | 2/23/07 5:46:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-20A

Client Sample ID: CW-2
Collection Date: 2/17/07 12:45:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Trichlorofluoromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| cis-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Chloroform | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Trichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Bromodichloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Toluene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Tetrachloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-20A

Client Sample ID: CW-2
Collection Date: 2/17/07 12:45:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Ethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| m,p-Xylene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| o-Xylene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,3,5-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2,4-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Naphthalene | ND | 5.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Surr: Dibromofluoromethane | 101 | 85-116 | | %REC | 1 | 2/23/07 5:11:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 106 | 77-127 | | %REC | 1 | 2/23/07 5:11:00 PM |
| Surr: Toluene-d8 | 101 | 86-114 | | %REC | 1 | 2/23/07 5:11:00 PM |
| Surr: 4-Bromofluorobenzene | 97.0 | 79-117 | | %REC | 1 | 2/23/07 5:11:00 PM |

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Method Blank

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

Sample ID: mb-02/21/07 Batch ID: R35962 Test Code: SW8260B Units: µg/L Analysis Date 2/21/2007 3:55:00 PM Prep Date: 2/21/2007
Client ID: Run ID: V-1_070221A SeqNo: 596167

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | QC |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|----|
| Dichlorodifluoromethane | ND | 5.0 | µg/L | | | | | | | | | |
| Chloromethane | ND | 5.0 | µg/L | | | | | | | | | |
| Vinyl chloride | ND | 2.0 | µg/L | | | | | | | | | |
| Chloroethane | ND | 5.0 | µg/L | | | | | | | | | |
| Bromomethane | ND | 2.0 | µg/L | | | | | | | | | |
| Trichlorofluoromethane | ND | 2.0 | µg/L | | | | | | | | | |
| Diethyl ether | ND | 5.0 | µg/L | | | | | | | | | |
| Acetone | ND | 10 | µg/L | | | | | | | | | |
| 1,1-Dichloroethene | ND | 1.0 | µg/L | | | | | | | | | |
| Carbon disulfide | ND | 2.0 | µg/L | | | | | | | | | |
| Methylene chloride | ND | 5.0 | µg/L | | | | | | | | | |
| Methyl tert-butyl ether | ND | 2.0 | µg/L | | | | | | | | | |
| trans-1,2-Dichloroethene | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1-Dichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| 2-Butanone | ND | 10 | µg/L | | | | | | | | | |
| 2,2-Dichloropropane | ND | 2.0 | µg/L | | | | | | | | | |
| cis-1,2-Dichloroethene | ND | 2.0 | µg/L | | | | | | | | | |
| Chloroform | ND | 2.0 | µg/L | | | | | | | | | |
| Tetrahydrofuran | ND | 10 | µg/L | | | | | | | | | |
| Bromochloromethane | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1,1-Trichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1-Dichloropropene | ND | 2.0 | µg/L | | | | | | | | | |
| Carbon tetrachloride | ND | 2.0 | µg/L | | | | | | | | | |
| 1,2-Dichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| Benzene | ND | 1.0 | µg/L | | | | | | | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 06-Mar-07

QC SUMMARY REPORT
Method Blank

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

| | | | |
|---------------------------|----|-----|------|
| Trichloroethene | ND | 2.0 | µg/L |
| 1,2-Dichloropropane | ND | 2.0 | µg/L |
| Bromodichloromethane | ND | 2.0 | µg/L |
| Dibromomethane | ND | 2.0 | µg/L |
| 4-Methyl-2-pentanone | ND | 10 | µg/L |
| cis-1,3-Dichloropropene | ND | 1.0 | µg/L |
| Toluene | ND | 2.0 | µg/L |
| trans-1,3-Dichloropropene | ND | 1.0 | µg/L |
| 1,1,2-Trichloroethane | ND | 2.0 | µg/L |
| 1,2-Dibromoethane | ND | 2.0 | µg/L |
| 2-Hexanone | ND | 10 | µg/L |
| 1,3-Dichloropropane | ND | 2.0 | µg/L |
| Tetrachloroethene | ND | 2.0 | µg/L |
| Dibromochloromethane | ND | 2.0 | µg/L |
| Chlorobenzene | ND | 2.0 | µg/L |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | µg/L |
| Ethylbenzene | ND | 2.0 | µg/L |
| m,p-Xylene | ND | 2.0 | µg/L |
| o-Xylene | ND | 2.0 | µg/L |
| Styrene | ND | 2.0 | µg/L |
| Bromoform | ND | 2.0 | µg/L |
| Isopropylbenzene | ND | 2.0 | µg/L |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | µg/L |
| 1,2,3-Trichloropropane | ND | 2.0 | µg/L |
| Bromobenzene | ND | 2.0 | µg/L |
| n-Propylbenzene | ND | 2.0 | µg/L |
| 2-Chlorotoluene | ND | 2.0 | µg/L |
| 4-Chlorotoluene | ND | 2.0 | µg/L |
| 1,3,5-Trimethylbenzene | ND | 2.0 | µg/L |
| tert-Butylbenzene | ND | 2.0 | µg/L |
| 1,2,4-Trimethylbenzene | ND | 2.0 | µg/L |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc. **QC SUMMARY REPORT**
 Work Order: 0702072 Method Blank
 Project: 101960 Textron Gorham

| Compound | Reporting Limit | Concentration | Recovery | Acceptance | Concentration | Concentration |
|-----------------------------|-----------------|---------------|----------|------------|---------------|---------------|
| sec-Butylbenzene | ND | 2.0 | 0 | 100 | 85 | 116 |
| 4-Isopropyltoluene | ND | 2.0 | 0 | 105 | 77 | 127 |
| 1,3-Dichlorobenzene | ND | 2.0 | 0 | 99.2 | 86 | 114 |
| 1,4-Dichlorobenzene | ND | 2.0 | 0 | 96 | 79 | 117 |
| n-Butylbenzene | ND | 2.0 | 0 | | | |
| 1,2-Dichlorobenzene | ND | 2.0 | 0 | | | |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | 0 | | | |
| 1,2,4-Trichlorobenzene | ND | 2.0 | 0 | | | |
| Hexachlorobutadiene | ND | 2.0 | 0 | | | |
| Naphthalene | ND | 5.0 | 0 | | | |
| 1,2,3-Trichlorobenzene | ND | 2.0 | 0 | | | |
| Surr: Dibromofluoromethane | 25.05 | 2.0 | 25 | 100 | 85 | 116 |
| Surr: 1,2-Dichloroethane-d4 | 26.3 | 2.0 | 25 | 105 | 77 | 127 |
| Surr: Toluene-d8 | 24.81 | 2.0 | 25 | 99.2 | 86 | 114 |
| Surr: 4-Bromofluorobenzene | 24.01 | 2.0 | 25 | 96 | 79 | 117 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Method Blank

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

Sample ID: mb-02/22/07 Batch ID: R35964 Test Code: SW8260B Units: µg/L Analysis Date 2/22/2007 1:28:00 PM Prep Date: 2/22/2007
Client ID: Run ID: V-1_070222A SeqNo: 596193

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | ND | 5.0 | µg/L | | | | | | | | | |
| Chloromethane | ND | 5.0 | µg/L | | | | | | | | | |
| Vinyl chloride | ND | 2.0 | µg/L | | | | | | | | | |
| Chloroethane | ND | 5.0 | µg/L | | | | | | | | | |
| Bromomethane | ND | 2.0 | µg/L | | | | | | | | | |
| Trichlorofluoromethane | ND | 2.0 | µg/L | | | | | | | | | |
| Diethyl ether | ND | 5.0 | µg/L | | | | | | | | | |
| Acetone | ND | 10 | µg/L | | | | | | | | | |
| 1,1-Dichloroethene | ND | 1.0 | µg/L | | | | | | | | | |
| Carbon disulfide | ND | 2.0 | µg/L | | | | | | | | | |
| Methylene chloride | ND | 5.0 | µg/L | | | | | | | | | |
| Methyl tert-butyl ether | ND | 2.0 | µg/L | | | | | | | | | |
| trans-1,2-Dichloroethene | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1-Dichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| 2-Butanone | ND | 10 | µg/L | | | | | | | | | |
| 2,2-Dichloropropane | ND | 2.0 | µg/L | | | | | | | | | |
| cis-1,2-Dichloroethene | ND | 2.0 | µg/L | | | | | | | | | |
| Chloroform | ND | 2.0 | µg/L | | | | | | | | | |
| Tetrahydrofuran | ND | 10 | µg/L | | | | | | | | | |
| Bromochloromethane | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1,1-Trichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1-Dichloropropene | ND | 2.0 | µg/L | | | | | | | | | |
| Carbon tetrachloride | ND | 2.0 | µg/L | | | | | | | | | |
| 1,2-Dichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| Benzene | ND | 1.0 | µg/L | | | | | | | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Method Blank

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

| | | | |
|---------------------------|----|-----|------|
| Trichloroethene | ND | 2.0 | µg/L |
| 1,2-Dichloropropane | ND | 2.0 | µg/L |
| Bromodichloromethane | ND | 2.0 | µg/L |
| Dibromomethane | ND | 2.0 | µg/L |
| 4-Methyl-2-pentanone | ND | 10 | µg/L |
| cis-1,3-Dichloropropene | ND | 1.0 | µg/L |
| Toluene | ND | 2.0 | µg/L |
| trans-1,3-Dichloropropene | ND | 1.0 | µg/L |
| 1,1,2-Trichloroethane | ND | 2.0 | µg/L |
| 1,2-Dibromoethane | ND | 2.0 | µg/L |
| 2-Hexanone | ND | 10 | µg/L |
| 1,3-Dichloropropane | ND | 2.0 | µg/L |
| Tetrachloroethene | ND | 2.0 | µg/L |
| Dibromochloromethane | ND | 2.0 | µg/L |
| Chlorobenzene | ND | 2.0 | µg/L |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | µg/L |
| Ethylbenzene | ND | 2.0 | µg/L |
| m,p-Xylene | ND | 2.0 | µg/L |
| o-Xylene | ND | 2.0 | µg/L |
| Styrene | ND | 2.0 | µg/L |
| Bromoform | ND | 2.0 | µg/L |
| Isopropylbenzene | ND | 2.0 | µg/L |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | µg/L |
| 1,2,3-Trichloropropane | ND | 2.0 | µg/L |
| Bromobenzene | ND | 2.0 | µg/L |
| n-Propylbenzene | ND | 2.0 | µg/L |
| 2-Chlorotoluene | ND | 2.0 | µg/L |
| 4-Chlorotoluene | ND | 2.0 | µg/L |
| 1,3,5-Trimethylbenzene | ND | 2.0 | µg/L |
| tert-Butylbenzene | ND | 2.0 | µg/L |
| 1,2,4-Trimethylbenzene | ND | 2.0 | µg/L |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.
 Work Order: 0702072
 Project: 101960 Textron Gorham

QC SUMMARY REPORT
 Method Blank

| Compound | Limit (µg/L) | Result | Recovery (%) | Recovery Limit | Result | Recovery (%) | Recovery Limit |
|-----------------------------|--------------|--------|--------------|----------------|--------|--------------|----------------|
| sec-Butylbenzene | 2.0 | ND | | | | | |
| 4-Isopropyltoluene | 2.0 | ND | | | | | |
| 1,3-Dichlorobenzene | 2.0 | ND | | | | | |
| 1,4-Dichlorobenzene | 2.0 | ND | | | | | |
| n-Butylbenzene | 2.0 | ND | | | | | |
| 1,2-Dichlorobenzene | 2.0 | ND | | | | | |
| 1,2-Dibromo-3-chloropropane | 5.0 | ND | | | | | |
| 1,2,4-Trichlorobenzene | 2.0 | ND | | | | | |
| Hexachlorobutadiene | 2.0 | ND | | | | | |
| Naphthalene | 5.0 | ND | | | | | |
| 1,2,3-Trichlorobenzene | 2.0 | ND | | | | | |
| Surr: Dibromofluoromethane | 25.63 | 25.63 | 100 | 25 | 85 | 116 | 0 |
| Surr: 1,2-Dichloroethane-d4 | 26.63 | 26.63 | 100 | 25 | 77 | 127 | 0 |
| Surr: Toluene-d8 | 25.19 | 25.19 | 100 | 25 | 86 | 114 | 0 |
| Surr: 4-Bromofluorobenzene | 24.73 | 24.73 | 100 | 25 | 79 | 117 | 0 |

Qualifiers: ND - Not Detected at the Reporting Limit
 S - Spike Recovery outside accepted recovery limits
 B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits
 R - RPD outside accepted recovery limits
 NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Method Blank

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

Sample ID: mb-02/23/07 Batch ID: R35993 Test Code: SW8260B Units: µg/L Analysis Date 2/23/2007 2:19:00 PM Prep Date: 2/23/2007
Client ID: Run ID: V-1_070223A SeqNo: 596565

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | ND | 5.0 | µg/L | | | | | | | | | |
| Chloromethane | ND | 5.0 | µg/L | | | | | | | | | |
| Vinyl chloride | ND | 2.0 | µg/L | | | | | | | | | |
| Chloroethane | ND | 5.0 | µg/L | | | | | | | | | |
| Bromomethane | ND | 2.0 | µg/L | | | | | | | | | |
| Trichlorofluoromethane | ND | 2.0 | µg/L | | | | | | | | | |
| Diethyl ether | ND | 5.0 | µg/L | | | | | | | | | |
| Acetone | ND | 10 | µg/L | | | | | | | | | |
| 1,1-Dichloroethene | ND | 1.0 | µg/L | | | | | | | | | |
| Carbon disulfide | ND | 2.0 | µg/L | | | | | | | | | |
| Methylene chloride | ND | 5.0 | µg/L | | | | | | | | | |
| Methyl tert-butyl ether | ND | 2.0 | µg/L | | | | | | | | | |
| trans-1,2-Dichloroethene | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1-Dichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| 2-Butanone | ND | 10 | µg/L | | | | | | | | | |
| 2,2-Dichloropropane | ND | 2.0 | µg/L | | | | | | | | | |
| cis-1,2-Dichloroethene | ND | 2.0 | µg/L | | | | | | | | | |
| Chloroform | ND | 2.0 | µg/L | | | | | | | | | |
| Tetrahydrofuran | ND | 10 | µg/L | | | | | | | | | |
| Bromochloromethane | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1,1-Trichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1-Dichloropropene | ND | 2.0 | µg/L | | | | | | | | | |
| Carbon tetrachloride | ND | 2.0 | µg/L | | | | | | | | | |
| 1,2-Dichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| Benzene | ND | 1.0 | µg/L | | | | | | | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Method Blank

| | | | |
|---------------------------|----|-----|------|
| Trichloroethene | ND | 2.0 | µg/L |
| 1,2-Dichloropropane | ND | 2.0 | µg/L |
| Bromodichloromethane | ND | 2.0 | µg/L |
| Dibromomethane | ND | 2.0 | µg/L |
| 4-Methyl-2-pentanone | ND | 10 | µg/L |
| cis-1,3-Dichloropropene | ND | 1.0 | µg/L |
| Toluene | ND | 2.0 | µg/L |
| trans-1,3-Dichloropropene | ND | 1.0 | µg/L |
| 1,1,2-Trichloroethane | ND | 2.0 | µg/L |
| 1,2-Dibromoethane | ND | 2.0 | µg/L |
| 2-Hexanone | ND | 10 | µg/L |
| 1,3-Dichloropropane | ND | 2.0 | µg/L |
| Tetrachloroethene | ND | 2.0 | µg/L |
| Dibromochloromethane | ND | 2.0 | µg/L |
| Chlorobenzene | ND | 2.0 | µg/L |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | µg/L |
| Ethylbenzene | ND | 2.0 | µg/L |
| m,p-Xylene | ND | 2.0 | µg/L |
| o-Xylene | ND | 2.0 | µg/L |
| Styrene | ND | 2.0 | µg/L |
| Bromoform | ND | 2.0 | µg/L |
| Isopropylbenzene | ND | 2.0 | µg/L |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | µg/L |
| 1,2,3-Trichloropropane | ND | 2.0 | µg/L |
| Bromobenzene | ND | 2.0 | µg/L |
| n-Propylbenzene | ND | 2.0 | µg/L |
| 2-Chlorotoluene | ND | 2.0 | µg/L |
| 4-Chlorotoluene | ND | 2.0 | µg/L |
| 1,3,5-Trimethylbenzene | ND | 2.0 | µg/L |
| tert-Butylbenzene | ND | 2.0 | µg/L |
| 1,2,4-Trimethylbenzene | ND | 2.0 | µg/L |

Qualifiers: ND - Not Detected at the Reporting Limit

S - Spike Recovery outside accepted recovery limits

R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank

NA - Not applicable where J values or ND results occur

J - Analyte detected below quantitation limits

RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Method Blank

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

| Compound | Reporting Limit | Concentration (µg/L) | Recovery (%) | Acceptance |
|-----------------------------|-----------------|----------------------|--------------|------------|
| sec-Butylbenzene | ND | 2.0 | | |
| 4-Isopropyltoluene | ND | 2.0 | | |
| 1,3-Dichlorobenzene | ND | 2.0 | | |
| 1,4-Dichlorobenzene | ND | 2.0 | | |
| n-Butylbenzene | ND | 2.0 | | |
| 1,2-Dichlorobenzene | ND | 2.0 | | |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | |
| Hexachlorobutadiene | ND | 2.0 | | |
| Naphthalene | ND | 5.0 | | |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | |
| Surr: Dibromofluoromethane | 24.83 | 2.0 | 99.3 | 85 |
| Surr: 1,2-Dichloroethane-d4 | 26.39 | 2.0 | 106 | 77 |
| Surr: Toluene-d8 | 25.4 | 2.0 | 102 | 86 |
| Surr: 4-Bromofluorobenzene | 23.56 | 2.0 | 94.2 | 79 |
| | | | | 116 |
| | | | | 127 |
| | | | | 114 |
| | | | | 117 |
| | | | | 0 |
| | | | | 0 |
| | | | | 0 |
| | | | | 0 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Laboratory Control Spike - Full List

Sample ID: lcsf-02/21/07 Batch ID: R35962 Test Code: SW8260B Units: µg/L Analysis Date 2/21/2007 2:46:00 PM Prep Date: 2/21/2007
 Client ID: Run ID: V-1_070221A SeqNo: 596168

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | 21.08 | 5.0 | µg/L | 20 | 0 | 105 | 10 | 150 | 0 | 0 | | |
| Chloromethane | 19.17 | 5.0 | µg/L | 20 | 0 | 95.8 | 37 | 150 | 0 | 0 | | |
| Vinyl chloride | 21.2 | 2.0 | µg/L | 20 | 0 | 106 | 48 | 150 | 0 | 0 | | |
| Chloroethane | 20.01 | 5.0 | µg/L | 20 | 0 | 100 | 54 | 142 | 0 | 0 | | |
| Bromomethane | 20.46 | 2.0 | µg/L | 20 | 0 | 102 | 51 | 137 | 0 | 0 | | |
| Trichlorofluoromethane | 23.7 | 2.0 | µg/L | 20 | 0 | 118 | 62 | 141 | 0 | 0 | | |
| Diethyl ether | 20.27 | 5.0 | µg/L | 20 | 0 | 101 | 68 | 134 | 0 | 0 | | |
| Acetone | 22.62 | 10 | µg/L | 20 | 0 | 113 | 9 | 150 | 0 | 0 | | |
| 1,1-Dichloroethene | 20.45 | 1.0 | µg/L | 20 | 0 | 102 | 68 | 146 | 0 | 0 | | |
| Carbon disulfide | 21.64 | 2.0 | µg/L | 20 | 0 | 108 | 52 | 131 | 0 | 0 | | |
| Methylene chloride | 22.19 | 5.0 | µg/L | 20 | 0 | 111 | 67 | 138 | 0 | 0 | | |
| Methyl tert-butyl ether | 21.1 | 2.0 | µg/L | 20 | 0 | 106 | 63 | 139 | 0 | 0 | | |
| trans-1,2-Dichloroethene | 24.74 | 2.0 | µg/L | 20 | 0 | 124 | 81 | 126 | 0 | 0 | | |
| 1,1-Dichloroethane | 23.63 | 2.0 | µg/L | 20 | 0 | 118 | 78 | 124 | 0 | 0 | | |
| 2-Butanone | 17.79 | 10 | µg/L | 20 | 0 | 89 | 41 | 150 | 0 | 0 | | |
| 2,2-Dichloropropane | 27.5 | 2.0 | µg/L | 20 | 0 | 138 | 71 | 150 | 0 | 0 | | |
| cis-1,2-Dichloroethene | 20.11 | 2.0 | µg/L | 20 | 0 | 101 | 78 | 121 | 0 | 0 | | |
| Chloroform | 21.68 | 2.0 | µg/L | 20 | 0 | 108 | 82 | 123 | 0 | 0 | | |
| Tetrahydrofuran | 20.82 | 10 | µg/L | 20 | 0 | 104 | 51 | 146 | 0 | 0 | | |
| Bromochloromethane | 19.97 | 2.0 | µg/L | 20 | 0 | 99.8 | 77 | 131 | 0 | 0 | | |
| 1,1,1-Trichloroethane | 24.06 | 2.0 | µg/L | 20 | 0 | 120 | 81 | 127 | 0 | 0 | | |
| 1,1-Dichloropropene | 21.81 | 2.0 | µg/L | 20 | 0 | 109 | 76 | 119 | 0 | 0 | | |
| Carbon tetrachloride | 21.4 | 2.0 | µg/L | 20 | 0 | 107 | 76 | 129 | 0 | 0 | | |
| 1,2-Dichloroethane | 21.64 | 2.0 | µg/L | 20 | 0 | 108 | 76 | 127 | 0 | 0 | | |
| Benzene | 21.83 | 1.0 | µg/L | 20 | 0 | 109 | 81 | 118 | 0 | 0 | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT

Laboratory Control Spike - Full List

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

| Compound | Reporting Limit | Concentration (µg/L) | Recovery (%) | Recovery Status | Method Blank |
|---------------------------|-----------------|----------------------|--------------|-----------------|--------------|
| Trichloroethene | 23.22 | 2.0 | 81 | 0 | 0 |
| 1,2-Dichloropropane | 21.68 | 2.0 | 79 | 0 | 0 |
| Bromodichloromethane | 21.01 | 2.0 | 77 | 0 | 0 |
| Dibromomethane | 19.94 | 2.0 | 76 | 0 | 0 |
| 4-Methyl-2-pentanone | 18.5 | 1.0 | 51 | 0 | 0 |
| cis-1,3-Dichloropropene | 20.94 | 1.0 | 76 | 0 | 0 |
| Toluene | 21.95 | 2.0 | 83 | 0 | 0 |
| trans-1,3-Dichloropropene | 20.93 | 1.0 | 66 | 0 | 0 |
| 1,1,2-Trichloroethane | 20.34 | 2.0 | 74 | 0 | 0 |
| 1,2-Dibromoethane | 20.65 | 2.0 | 72 | 0 | 0 |
| 2-Hexanone | 18.18 | 1.0 | 31 | 0 | 0 |
| 1,3-Dichloropropane | 20.59 | 2.0 | 76 | 0 | 0 |
| Tetrachloroethene | 22.29 | 2.0 | 81 | 0 | 0 |
| Dibromochloromethane | 19.81 | 2.0 | 63 | 0 | 0 |
| Chlorobenzene | 19.91 | 2.0 | 84 | 0 | 0 |
| 1,1,1,2-Tetrachloroethane | 20.76 | 2.0 | 73 | 0 | 0 |
| Ethylbenzene | 21.45 | 2.0 | 83 | 0 | 0 |
| m,p-Xylene | 43.07 | 2.0 | 85 | 0 | 0 |
| o-Xylene | 19.94 | 2.0 | 84 | 0 | 0 |
| Styrene | 20.05 | 2.0 | 81 | 0 | 0 |
| Bromoforn | 17.31 | 2.0 | 55 | 0 | 0 |
| Isopropylbenzene | 23.27 | 2.0 | 77 | 0 | 0 |
| 1,1,2,2-Tetrachloroethane | 20.1 | 2.0 | 62 | 0 | 0 |
| 1,2,3-Trichloropropane | 20 | 2.0 | 62 | 0 | 0 |
| Bromobenzene | 19.36 | 2.0 | 78 | 0 | 0 |
| n-Propylbenzene | 22.06 | 2.0 | 77 | 0 | 0 |
| 2-Chlorotoluene | 21.4 | 2.0 | 78 | 0 | 0 |
| 4-Chlorotoluene | 21.95 | 2.0 | 77 | 0 | 0 |
| 1,3,5-Trimethylbenzene | 22.57 | 2.0 | 80 | 0 | 0 |
| tert-Butylbenzene | 23.1 | 2.0 | 81 | 0 | 0 |
| 1,2,4-Trimethylbenzene | 22.04 | 2.0 | 80 | 0 | 0 |

Qualifiers: ND - Not Detected at the Reporting Limit

S - Spike Recovery outside accepted recovery limits

B - Analyte detected in the associated Method Blank

J - Analyte detected below quantitation limits

R - RPD outside accepted recovery limits

NA - Not applicable where J values or ND results occur

RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Laboratory Control Spike - Full List

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

| Compound | Concentration (µg/L) | Recovery (%) | Acceptance | Reporting Limit (µg/L) | Concentration (µg/L) | Recovery (%) | Acceptance |
|-----------------------------|----------------------|--------------|------------|------------------------|----------------------|--------------|------------|
| sec-Butylbenzene | 19.73 | 2.0 | µg/L | 20 | 98.6 | 82 | 123 |
| 4-Isopropyltoluene | 22.98 | 2.0 | µg/L | 20 | 115 | 80 | 126 |
| 1,3-Dichlorobenzene | 20.22 | 2.0 | µg/L | 20 | 101 | 84 | 115 |
| 1,4-Dichlorobenzene | 20.42 | 2.0 | µg/L | 20 | 102 | 79 | 117 |
| n-Butylbenzene | 19.34 | 2.0 | µg/L | 20 | 96.7 | 76 | 128 |
| 1,2-Dichlorobenzene | 19.92 | 2.0 | µg/L | 20 | 99.6 | 81 | 117 |
| 1,2-Dibromo-3-chloropropane | 18.61 | 5.0 | µg/L | 20 | 93 | 47 | 136 |
| 1,2,4-Trichlorobenzene | 20.38 | 2.0 | µg/L | 20 | 102 | 73 | 126 |
| Hexachlorobutadiene | 20.28 | 2.0 | µg/L | 20 | 101 | 77 | 134 |
| Naphthalene | 20.73 | 5.0 | µg/L | 20 | 104 | 58 | 138 |
| 1,2,3-Trichlorobenzene | 20.62 | 2.0 | µg/L | 20 | 103 | 76 | 124 |
| Surr: Dibromofluoromethane | 25.69 | 2.0 | µg/L | 25 | 103 | 85 | 116 |
| Surr: 1,2-Dichloroethane-d4 | 27.39 | 2.0 | µg/L | 25 | 110 | 77 | 127 |
| Surr: Toluene-d8 | 25.13 | 2.0 | µg/L | 25 | 101 | 86 | 114 |
| Surr: 4-Bromofluorobenzene | 24.25 | 2.0 | µg/L | 25 | 97 | 79 | 117 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Laboratory Control Spike - Full List

Sample ID: lcsf-02/22/07 Batch ID: R35964 Test Code: SW8260B Units: µg/L Analysis Date 2/22/2007 12:18:00 PM Prep Date: 2/22/2007
 Client ID: Run ID: V-1_070222A SeqNo: 596194

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | 16.24 | 5.0 | µg/L | 20 | 0 | 81.2 | 10 | 150 | 0 | 0 | 0 | |
| Chloromethane | 16.21 | 5.0 | µg/L | 20 | 0 | 81 | 37 | 150 | 0 | 0 | 0 | |
| Vinyl chloride | 17.6 | 2.0 | µg/L | 20 | 0 | 88 | 48 | 150 | 0 | 0 | 0 | |
| Chloroethane | 18.68 | 5.0 | µg/L | 20 | 0 | 93.4 | 54 | 142 | 0 | 0 | 0 | |
| Bromomethane | 18.53 | 2.0 | µg/L | 20 | 0 | 92.6 | 51 | 137 | 0 | 0 | 0 | |
| Trichlorofluoromethane | 19.6 | 2.0 | µg/L | 20 | 0 | 98 | 62 | 141 | 0 | 0 | 0 | |
| Diethyl ether | 20.64 | 5.0 | µg/L | 20 | 0 | 103 | 68 | 134 | 0 | 0 | 0 | |
| Acetone | 20.99 | 10 | µg/L | 20 | 0 | 105 | 9 | 150 | 0 | 0 | 0 | |
| 1,1-Dichloroethene | 17.2 | 1.0 | µg/L | 20 | 0 | 86 | 68 | 146 | 0 | 0 | 0 | |
| Carbon disulfide | 18.52 | 2.0 | µg/L | 20 | 0 | 92.6 | 52 | 131 | 0 | 0 | 0 | |
| Methylene chloride | 20.46 | 5.0 | µg/L | 20 | 0 | 102 | 67 | 138 | 0 | 0 | 0 | |
| Methyl tert-butyl ether | 20.94 | 2.0 | µg/L | 20 | 0 | 105 | 63 | 139 | 0 | 0 | 0 | |
| trans-1,2-Dichloroethene | 22.09 | 2.0 | µg/L | 20 | 0 | 110 | 81 | 126 | 0 | 0 | 0 | |
| 1,1-Dichloroethane | 21.87 | 2.0 | µg/L | 20 | 0 | 109 | 78 | 124 | 0 | 0 | 0 | |
| 2-Butanone | 21.48 | 10 | µg/L | 20 | 0 | 107 | 41 | 150 | 0 | 0 | 0 | |
| 2,2-Dichloropropane | 25.31 | 2.0 | µg/L | 20 | 0 | 127 | 71 | 150 | 0 | 0 | 0 | |
| cis-1,2-Dichloroethene | 18.83 | 2.0 | µg/L | 20 | 0 | 94.2 | 78 | 121 | 0 | 0 | 0 | |
| Chloroform | 20.98 | 2.0 | µg/L | 20 | 0 | 105 | 82 | 123 | 0 | 0 | 0 | |
| Tetrahydrofuran | 21.18 | 10 | µg/L | 20 | 0 | 106 | 51 | 146 | 0 | 0 | 0 | |
| Bromochloromethane | 20.02 | 2.0 | µg/L | 20 | 0 | 100 | 77 | 131 | 0 | 0 | 0 | |
| 1,1,1-Trichloroethane | 21.77 | 2.0 | µg/L | 20 | 0 | 109 | 81 | 127 | 0 | 0 | 0 | |
| 1,1-Dichloropropene | 19.82 | 2.0 | µg/L | 20 | 0 | 99.1 | 76 | 119 | 0 | 0 | 0 | |
| Carbon tetrachloride | 19.22 | 2.0 | µg/L | 20 | 0 | 96.1 | 76 | 129 | 0 | 0 | 0 | |
| 1,2-Dichloroethane | 22.06 | 2.0 | µg/L | 20 | 0 | 110 | 76 | 127 | 0 | 0 | 0 | |
| Benzene | 20.23 | 1.0 | µg/L | 20 | 0 | 101 | 81 | 118 | 0 | 0 | 0 | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT

Laboratory Control Spike - Full List

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

| Compound | Reporting Limit | Concentration (µg/L) | Recovery (%) | Recovery Outside Limits | Accepted Recovery Limits | Number of Samples | Number of Samples with Analyte |
|---------------------------|-----------------|----------------------|--------------|-------------------------|--------------------------|-------------------|--------------------------------|
| Trichloroethene | 20.45 | 2.0 | 102 | 0 | 0 | 81 | 119 |
| 1,2-Dichloropropane | 19.96 | 2.0 | 99.8 | 0 | 0 | 79 | 120 |
| Bromodichloromethane | 20.35 | 2.0 | 102 | 0 | 0 | 77 | 131 |
| Dibromomethane | 19.6 | 2.0 | 98 | 0 | 0 | 76 | 128 |
| 4-Methyl-2-pentanone | 19.03 | 1.0 | 95.2 | 0 | 0 | 51 | 141 |
| cis-1,3-Dichloropropene | 19.03 | 1.0 | 95.2 | 0 | 0 | 76 | 120 |
| Toluene | 19.8 | 2.0 | 99 | 0 | 0 | 83 | 119 |
| trans-1,3-Dichloropropene | 19.8 | 1.0 | 99 | 0 | 0 | 66 | 128 |
| 1,1,2-Trichloroethane | 20.39 | 2.0 | 102 | 0 | 0 | 74 | 123 |
| 1,2-Dibromoethane | 20.94 | 2.0 | 105 | 0 | 0 | 72 | 128 |
| 2-Hexanone | 18.04 | 1.0 | 90.2 | 0 | 0 | 31 | 148 |
| 1,3-Dichloropropane | 19.76 | 2.0 | 98.8 | 0 | 0 | 76 | 122 |
| Tetrachloroethene | 20.03 | 2.0 | 100 | 0 | 0 | 81 | 124 |
| Dibromochloromethane | 18.27 | 2.0 | 91.4 | 0 | 0 | 63 | 126 |
| Chlorobenzene | 18.47 | 2.0 | 92.4 | 0 | 0 | 84 | 113 |
| 1,1,1,2-Tetrachloroethane | 18.74 | 2.0 | 93.7 | 0 | 0 | 73 | 124 |
| Ethylbenzene | 19.73 | 2.0 | 98.6 | 0 | 0 | 83 | 118 |
| m,p-Xylene | 40.32 | 2.0 | 101 | 0 | 0 | 85 | 116 |
| o-Xylene | 18.71 | 2.0 | 93.6 | 0 | 0 | 84 | 115 |
| Styrene | 19.38 | 2.0 | 96.9 | 0 | 0 | 81 | 118 |
| Bromoform | 17.07 | 2.0 | 85.4 | 0 | 0 | 55 | 126 |
| Isopropylbenzene | 20.36 | 2.0 | 102 | 0 | 0 | 77 | 125 |
| 1,1,2,2-Tetrachloroethane | 18.21 | 2.0 | 91 | 0 | 0 | 62 | 134 |
| 1,2,3-Trichloropropane | 19.34 | 2.0 | 96.7 | 0 | 0 | 62 | 132 |
| Bromobenzene | 18.84 | 2.0 | 94.2 | 0 | 0 | 78 | 119 |
| n-Propylbenzene | 19.34 | 2.0 | 96.7 | 0 | 0 | 77 | 127 |
| 2-Chlorotoluene | 19.29 | 2.0 | 96.5 | 0 | 0 | 78 | 118 |
| 4-Chlorotoluene | 19.62 | 2.0 | 98.1 | 0 | 0 | 77 | 119 |
| 1,3,5-Trimethylbenzene | 20.01 | 2.0 | 100 | 0 | 0 | 80 | 120 |
| tert-Butylbenzene | 20.46 | 2.0 | 102 | 0 | 0 | 81 | 120 |
| 1,2,4-Trimethylbenzene | 19.7 | 2.0 | 98.5 | 0 | 0 | 80 | 118 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Laboratory Control Spike - Full List

Sample ID: **lcst-02/23/07** Batch ID: **R35993** Test Code: **SW8260B** Units: **µg/L** Analysis Date **2/23/2007 12:35:00 PM** Prep Date: **2/23/2007**
 Client ID: **596566** Run ID: **V-1_070223A** SeqNo: **596566**

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | 15.43 | 5.0 | µg/L | 20 | 0 | 77.2 | 10 | 150 | 0 | 0 | | |
| Chloromethane | 17.02 | 5.0 | µg/L | 20 | 0 | 85.1 | 37 | 150 | 0 | 0 | | |
| Vinyl chloride | 19.06 | 2.0 | µg/L | 20 | 0 | 95.3 | 48 | 150 | 0 | 0 | | |
| Chloroethane | 20.04 | 5.0 | µg/L | 20 | 0 | 100 | 54 | 142 | 0 | 0 | | |
| Bromomethane | 19.28 | 2.0 | µg/L | 20 | 0 | 96.4 | 51 | 137 | 0 | 0 | | |
| Trichlorofluoromethane | 21.68 | 2.0 | µg/L | 20 | 0 | 108 | 62 | 141 | 0 | 0 | | |
| Diethyl ether | 18.88 | 5.0 | µg/L | 20 | 0 | 94.4 | 68 | 134 | 0 | 0 | | |
| Acetone | 14.65 | 10 | µg/L | 20 | 0 | 73.2 | 9 | 150 | 0 | 0 | | |
| 1,1-Dichloroethene | 18.7 | 1.0 | µg/L | 20 | 0 | 93.5 | 68 | 146 | 0 | 0 | | |
| Carbon disulfide | 20.28 | 2.0 | µg/L | 20 | 0 | 101 | 52 | 131 | 0 | 0 | | |
| Methylene chloride | 21.88 | 5.0 | µg/L | 20 | 0 | 109 | 67 | 138 | 0 | 0 | | |
| Methyl tert-butyl ether | 21.55 | 2.0 | µg/L | 20 | 0 | 108 | 63 | 139 | 0 | 0 | | |
| trans-1,2-Dichloroethene | 22.93 | 2.0 | µg/L | 20 | 0 | 115 | 81 | 126 | 0 | 0 | | |
| 1,1-Dichloroethane | 22.97 | 2.0 | µg/L | 20 | 0 | 115 | 78 | 124 | 0 | 0 | | |
| 2-Butanone | 15.28 | 10 | µg/L | 20 | 0 | 76.4 | 41 | 150 | 0 | 0 | | |
| 2,2-Dichloropropane | 27.34 | 2.0 | µg/L | 20 | 0 | 137 | 71 | 150 | 0 | 0 | | |
| cis-1,2-Dichloroethene | 19.43 | 2.0 | µg/L | 20 | 0 | 97.2 | 78 | 121 | 0 | 0 | | |
| Chloroform | 21.77 | 2.0 | µg/L | 20 | 0 | 109 | 82 | 123 | 0 | 0 | | |
| Tetrahydrofuran | 16.64 | 10 | µg/L | 20 | 0 | 83.2 | 51 | 146 | 0 | 0 | | |
| Bromochloromethane | 21.29 | 2.0 | µg/L | 20 | 0 | 106 | 77 | 131 | 0 | 0 | | |
| 1,1,1-Trichloroethane | 23.4 | 2.0 | µg/L | 20 | 0 | 117 | 81 | 127 | 0 | 0 | | |
| 1,1-Dichloropropene | 20.56 | 2.0 | µg/L | 20 | 0 | 103 | 76 | 119 | 0 | 0 | | |
| Carbon tetrachloride | 21.12 | 2.0 | µg/L | 20 | 0 | 106 | 76 | 129 | 0 | 0 | | |
| 1,2-Dichloroethane | 21.42 | 2.0 | µg/L | 20 | 0 | 107 | 76 | 127 | 0 | 0 | | |
| Benzene | 21.03 | 1.0 | µg/L | 20 | 0 | 105 | 81 | 118 | 0 | 0 | | |

Qualifiers: ND - Not Detected at the Reporting Limit

J - Analyte detected below quantitation limits

RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

S - Spike Recovery outside accepted recovery limits

R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank

NA - Not applicable where J values or ND results occur

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

QC SUMMARY REPORT

Laboratory Control Spike - Full List

| Compound | Reported Limit | Concentration | Recovery | Accepted Limits | Recovery | Accepted Limits | Recovery | Accepted Limits | Recovery |
|---------------------------|----------------|---------------|----------|-----------------|----------|-----------------|----------|-----------------|----------|
| Trichloroethene | 22.08 | 2.0 | µg/L | 20 | 0 | 110 | 81 | 119 | 0 |
| 1,2-Dichloropropane | 20.44 | 2.0 | µg/L | 20 | 0 | 102 | 79 | 120 | 0 |
| Bromodichloromethane | 20.44 | 2.0 | µg/L | 20 | 0 | 102 | 77 | 131 | 0 |
| Dibromomethane | 19.16 | 2.0 | µg/L | 20 | 0 | 95.8 | 76 | 128 | 0 |
| 4-Methyl-2-pentanone | 14.8 | 10 | µg/L | 20 | 0 | 74 | 51 | 141 | 0 |
| cis-1,3-Dichloropropene | 20.12 | 1.0 | µg/L | 20 | 0 | 101 | 76 | 120 | 0 |
| Toluene | 21.6 | 2.0 | µg/L | 20 | 0 | 108 | 83 | 119 | 0 |
| trans-1,3-Dichloropropene | 20 | 1.0 | µg/L | 20 | 0 | 100 | 66 | 128 | 0 |
| 1,1,2-Trichloroethane | 20 | 2.0 | µg/L | 20 | 0 | 100 | 74 | 123 | 0 |
| 1,2-Dibromoethane | 19.75 | 2.0 | µg/L | 20 | 0 | 98.8 | 72 | 128 | 0 |
| 2-Hexanone | 15.04 | 10 | µg/L | 20 | 0 | 75.2 | 31 | 148 | 0 |
| 1,3-Dichloropropane | 19.93 | 2.0 | µg/L | 20 | 0 | 99.7 | 76 | 122 | 0 |
| Tetrachloroethene | 21.64 | 2.0 | µg/L | 20 | 0 | 108 | 81 | 124 | 0 |
| Dibromochloromethane | 19.61 | 2.0 | µg/L | 20 | 0 | 98 | 63 | 126 | 0 |
| Chlorobenzene | 19.88 | 2.0 | µg/L | 20 | 0 | 99.4 | 84 | 113 | 0 |
| 1,1,1,2-Tetrachloroethane | 20.01 | 2.0 | µg/L | 20 | 0 | 100 | 73 | 124 | 0 |
| Ethylbenzene | 21.26 | 2.0 | µg/L | 20 | 0 | 106 | 83 | 118 | 0 |
| m,p-Xylene | 42.16 | 2.0 | µg/L | 40 | 0 | 105 | 85 | 116 | 0 |
| o-Xylene | 19.94 | 2.0 | µg/L | 20 | 0 | 99.7 | 84 | 115 | 0 |
| Styrene | 20.37 | 2.0 | µg/L | 20 | 0 | 102 | 81 | 118 | 0 |
| Bromoform | 16.71 | 2.0 | µg/L | 20 | 0 | 83.6 | 55 | 126 | 0 |
| Isopropylbenzene | 22.03 | 2.0 | µg/L | 20 | 0 | 110 | 77 | 125 | 0 |
| 1,1,2,2-Tetrachloroethane | 16.78 | 2.0 | µg/L | 20 | 0 | 83.9 | 62 | 134 | 0 |
| 1,2,3-Trichloropropane | 18.26 | 2.0 | µg/L | 20 | 0 | 91.3 | 62 | 132 | 0 |
| Bromobenzene | 18.81 | 2.0 | µg/L | 20 | 0 | 94.1 | 78 | 119 | 0 |
| n-Propylbenzene | 21 | 2.0 | µg/L | 20 | 0 | 105 | 77 | 127 | 0 |
| 2-Chlorotoluene | 20.66 | 2.0 | µg/L | 20 | 0 | 103 | 78 | 118 | 0 |
| 4-Chlorotoluene | 20.8 | 2.0 | µg/L | 20 | 0 | 104 | 77 | 119 | 0 |
| 1,3,5-Trimethylbenzene | 21.4 | 2.0 | µg/L | 20 | 0 | 107 | 80 | 120 | 0 |
| tert-Butylbenzene | 21.88 | 2.0 | µg/L | 20 | 0 | 109 | 81 | 120 | 0 |
| 1,2,4-Trimethylbenzene | 20.8 | 2.0 | µg/L | 20 | 0 | 104 | 80 | 118 | 0 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

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Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Laboratory Control Spike - Full List

| Compound | 18.23 | 21.45 | 19.37 | 19.03 | 17.29 | 18.93 | 15.41 | 17.9 | 17.35 | 17.05 | 16.93 | 25.04 | 25.82 | 25.56 | 24.62 |
|-----------------------------|-------|-------|-------|-------|-------|-------|-------|------|-------|-------|-------|-------|-------|-------|-------|
| sec-Butylbenzene | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 4-Isopropyltoluene | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L |
| 1,3-Dichlorobenzene | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 25 | 25 | 25 | 25 |
| 1,4-Dichlorobenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| n-Butylbenzene | 91.2 | 107 | 96.8 | 95.2 | 86.5 | 94.6 | 77 | 89.5 | 86.8 | 85.2 | 84.6 | 100 | 103 | 102 | 98.5 |
| 1,2-Dichlorobenzene | 82 | 80 | 84 | 79 | 76 | 81 | 47 | 73 | 77 | 58 | 76 | 85 | 77 | 86 | 79 |
| 1,2-Dibromo-3-chloropropane | 123 | 126 | 115 | 117 | 128 | 117 | 136 | 126 | 134 | 138 | 124 | 116 | 127 | 114 | 117 |
| 1,2,4-Trichlorobenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Hexachlorobutadiene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Naphthalene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 1,2,3-Trichlorobenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Surr: Dibromofluoromethane | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Surr: 1,2-Dichloroethane-d4 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Surr: Toluene-d8 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Surr: 4-Bromofluorobenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |

Qualifiers: ND - Not Detected at the Reporting Limit
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S - Spike Recovery outside accepted recovery limits
 R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank
 NA - Not applicable where J values or ND results occur

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.
 Work Order: 0702072
 Project: 101960 Textron Gorham

QC SUMMARY REPORT
 Laboratory Control Spike Duplicate - Full List

| Sample ID: | Batch ID: | Test Code: | Units: | Analysis Date | Prep Date: | | | | | | | |
|--------------------------|------------------|------------|--------|----------------------|------------------------|-----------|----------|------|------------------------------|----------|--------|-------|
| iesdf-02/23/07 | R35993 | SW8260B | µg/L | 2/23/2007 1:09:00 PM | 2/23/2007 | | | | | | | |
| Client ID: | Run ID: | SeqNo: | | | | | | | | | | |
| | V-1_070223A | 596567 | | | | | | | | | | |
| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | HighLimit | LowLimit | %REC | Original Sample or MS Result | RPDLimit | RPD | Queue |
| Dichlorodifluoromethane | 16.78 | 5.0 | µg/L | 20 | 0 | 10 | 150 | 83.9 | 15.43 | 25 | 8.38 | 25 |
| Chloromethane | 17.69 | 5.0 | µg/L | 20 | 0 | 37 | 150 | 88.4 | 17.02 | 25 | 3.86 | 25 |
| Vinyl chloride | 19.34 | 2.0 | µg/L | 20 | 0 | 48 | 150 | 96.7 | 19.06 | 25 | 1.46 | 25 |
| Chloroethane | 18.49 | 5.0 | µg/L | 20 | 0 | 54 | 142 | 92.5 | 20.04 | 25 | 8.05 | 25 |
| Bromomethane | 19.3 | 2.0 | µg/L | 20 | 0 | 51 | 137 | 96.5 | 19.28 | 25 | 0.104 | 25 |
| Trichlorofluoromethane | 21.62 | 2.0 | µg/L | 20 | 0 | 62 | 141 | 108 | 21.68 | 25 | 0.277 | 25 |
| Acetone | 16.26 | 10 | µg/L | 20 | 0 | 9 | 150 | 81.3 | 14.65 | 25 | 10.4 | 25 |
| 1,1-Dichloroethene | 18.55 | 1.0 | µg/L | 20 | 0 | 68 | 146 | 92.8 | 18.7 | 25 | 0.805 | 25 |
| Carbon disulfide | 20.33 | 2.0 | µg/L | 20 | 0 | 52 | 131 | 102 | 20.28 | 25 | 0.246 | 25 |
| Methylene chloride | 22.14 | 5.0 | µg/L | 20 | 0 | 67 | 138 | 111 | 21.88 | 25 | 1.18 | 25 |
| Methyl tert-butyl ether | 21.45 | 2.0 | µg/L | 20 | 0 | 63 | 139 | 107 | 21.55 | 25 | 0.465 | 25 |
| trans-1,2-Dichloroethene | 23.12 | 2.0 | µg/L | 20 | 0 | 81 | 126 | 116 | 22.93 | 25 | 0.825 | 25 |
| 1,1-Dichloroethane | 23.22 | 2.0 | µg/L | 20 | 0 | 78 | 124 | 116 | 22.97 | 25 | 1.08 | 25 |
| 2-Butanone | 15.15 | 10 | µg/L | 20 | 0 | 41 | 150 | 75.8 | 15.28 | 25 | 0.854 | 25 |
| 2,2-Dichloropropane | 26.83 | 2.0 | µg/L | 20 | 0 | 71 | 150 | 134 | 27.34 | 25 | 1.88 | 25 |
| cis-1,2-Dichloroethene | 19.61 | 2.0 | µg/L | 20 | 0 | 78 | 121 | 98 | 19.43 | 25 | 0.922 | 25 |
| Chloroform | 21 | 2.0 | µg/L | 20 | 0 | 82 | 123 | 105 | 21.77 | 25 | 3.6 | 25 |
| Bromochloromethane | 20.65 | 2.0 | µg/L | 20 | 0 | 77 | 131 | 103 | 21.29 | 25 | 3.05 | 25 |
| 1,1,1-Trichloroethane | 23.22 | 2.0 | µg/L | 20 | 0 | 81 | 127 | 116 | 23.4 | 25 | 0.772 | 25 |
| 1,1-Dichloropropene | 20.55 | 2.0 | µg/L | 20 | 0 | 76 | 119 | 103 | 20.56 | 25 | 0.0486 | 25 |
| Carbon tetrachloride | 20.73 | 2.0 | µg/L | 20 | 0 | 76 | 129 | 104 | 21.12 | 25 | 1.86 | 25 |
| 1,2-Dichloroethane | 21.76 | 2.0 | µg/L | 20 | 0 | 76 | 127 | 109 | 21.42 | 25 | 1.57 | 25 |
| Benzene | 20.64 | 1.0 | µg/L | 20 | 0 | 81 | 118 | 103 | 21.03 | 25 | 1.87 | 25 |
| Trichloroethene | 21.44 | 2.0 | µg/L | 20 | 0 | 81 | 119 | 107 | 22.08 | 25 | 2.94 | 25 |
| 1,2-Dichloropropane | 20.98 | 2.0 | µg/L | 20 | 0 | 79 | 120 | 105 | 20.44 | 25 | 2.61 | 25 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
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AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Laboratory Control Spike Duplicate - Full List

| Compound | Reporting Limit | ND - Not Detected at the Reporting Limit | S - Spike Recovery outside accepted recovery limits | R - RPD outside accepted recovery limits | RL - Reporting Limit, defined as the lowest concentration the laboratory can accurately quantitate. | Qualifiers: | | | | | |
|---------------------------|-----------------|--|---|--|---|-------------|----|-----|-------|-------|----|
| Bromodichloromethane | 20.49 | 2.0 | µg/L | 20 | 0 | 102 | 77 | 131 | 20.44 | 0.244 | 25 |
| Dibromomethane | 18.82 | 2.0 | µg/L | 20 | 0 | 94.1 | 76 | 128 | 19.16 | 1.79 | 25 |
| 4-Methyl-2-pentanone | 16.64 | 10 | µg/L | 20 | 0 | 83.2 | 51 | 141 | 14.8 | 11.7 | 25 |
| cis-1,3-Dichloropropene | 19.67 | 1.0 | µg/L | 20 | 0 | 98.4 | 76 | 120 | 20.12 | 2.26 | 25 |
| Toluene | 21.28 | 2.0 | µg/L | 20 | 0 | 106 | 83 | 119 | 21.6 | 1.49 | 25 |
| trans-1,3-Dichloropropene | 19.79 | 1.0 | µg/L | 20 | 0 | 99 | 66 | 128 | 20 | 1.06 | 25 |
| 1,1,2-Trichloroethane | 19.36 | 2.0 | µg/L | 20 | 0 | 96.8 | 74 | 123 | 20 | 3.25 | 25 |
| 1,2-Dibromoethane | 19.54 | 2.0 | µg/L | 20 | 0 | 97.7 | 72 | 128 | 19.75 | 1.07 | 25 |
| 2-Hexanone | 14.97 | 10 | µg/L | 20 | 0 | 74.8 | 31 | 148 | 15.04 | 0.467 | 25 |
| 1,3-Dichloropropane | 19.47 | 2.0 | µg/L | 20 | 0 | 97.4 | 76 | 122 | 19.93 | 2.34 | 25 |
| Tetrachloroethene | 20.82 | 2.0 | µg/L | 20 | 0 | 104 | 81 | 124 | 21.64 | 3.86 | 25 |
| Dibromochloromethane | 18.49 | 2.0 | µg/L | 20 | 0 | 92.5 | 63 | 126 | 19.61 | 5.88 | 25 |
| Chlorobenzene | 19.43 | 2.0 | µg/L | 20 | 0 | 97.2 | 84 | 113 | 19.88 | 2.29 | 25 |
| 1,1,1,2-Tetrachloroethane | 19.75 | 2.0 | µg/L | 20 | 0 | 98.8 | 73 | 124 | 20.01 | 1.31 | 25 |
| Ethylbenzene | 20.76 | 2.0 | µg/L | 20 | 0 | 104 | 83 | 118 | 21.26 | 2.38 | 25 |
| m,p-Xylene | 41.65 | 2.0 | µg/L | 40 | 0 | 104 | 85 | 116 | 42.16 | 1.22 | 25 |
| o-Xylene | 19.69 | 2.0 | µg/L | 20 | 0 | 98.4 | 84 | 115 | 19.94 | 1.26 | 25 |
| Styrene | 19.55 | 2.0 | µg/L | 20 | 0 | 97.8 | 81 | 118 | 20.37 | 4.11 | 25 |
| Bromoforn | 15.81 | 2.0 | µg/L | 20 | 0 | 79 | 55 | 126 | 16.71 | 5.54 | 25 |
| Isopropylbenzene | 21.93 | 2.0 | µg/L | 20 | 0 | 110 | 77 | 125 | 22.03 | 0.455 | 25 |
| 1,1,2,2-Tetrachloroethane | 17.16 | 2.0 | µg/L | 20 | 0 | 85.8 | 62 | 134 | 16.78 | 2.24 | 25 |
| 1,2,3-Trichloropropane | 17.42 | 2.0 | µg/L | 20 | 0 | 87.1 | 62 | 132 | 18.26 | 4.71 | 25 |
| Bromobenzene | 18.77 | 2.0 | µg/L | 20 | 0 | 93.8 | 78 | 119 | 18.81 | 0.213 | 25 |
| n-Propylbenzene | 21.46 | 2.0 | µg/L | 20 | 0 | 107 | 77 | 127 | 21 | 2.17 | 25 |
| 2-Chlorotoluene | 20.62 | 2.0 | µg/L | 20 | 0 | 103 | 78 | 118 | 20.66 | 0.194 | 25 |
| 4-Chlorotoluene | 20.95 | 2.0 | µg/L | 20 | 0 | 105 | 77 | 119 | 20.8 | 0.719 | 25 |
| 1,3,5-Trimethylbenzene | 20.87 | 2.0 | µg/L | 20 | 0 | 104 | 80 | 120 | 21.4 | 2.51 | 25 |
| tert-Butylbenzene | 20.88 | 2.0 | µg/L | 20 | 0 | 104 | 81 | 120 | 21.88 | 4.68 | 25 |
| 1,2,4-Trimethylbenzene | 20.48 | 2.0 | µg/L | 20 | 0 | 102 | 80 | 118 | 20.8 | 1.55 | 25 |
| sec-Butylbenzene | 17.89 | 2.0 | µg/L | 20 | 0 | 89.4 | 82 | 123 | 18.23 | 1.88 | 25 |
| 4-Isopropyltoluene | 20.59 | 2.0 | µg/L | 20 | 0 | 103 | 80 | 126 | 21.45 | 4.09 | 25 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
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AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Laboratory Control Spike Duplicate - Full List

| | 19.17 | 2.0 | µg/L | 20 | 0 | 95.8 | 84 | 115 | 19.37 | 1.04 | 25 |
|-----------------------------|-------|-----|------|----|---|------|----|-----|-------|------|----|
| 1,3-Dichlorobenzene | 18.82 | 2.0 | µg/L | 20 | 0 | 94.1 | 79 | 117 | 19.03 | 1.11 | 25 |
| 1,4-Dichlorobenzene | 17.01 | 2.0 | µg/L | 20 | 0 | 85 | 76 | 128 | 17.29 | 1.63 | 25 |
| n-Butylbenzene | 18.01 | 2.0 | µg/L | 20 | 0 | 90 | 81 | 117 | 18.93 | 4.98 | 25 |
| 1,2-Dichlorobenzene | 15.7 | 5.0 | µg/L | 20 | 0 | 78.5 | 47 | 136 | 15.41 | 1.86 | 25 |
| 1,2-Dibromo-3-chloropropane | 16.55 | 2.0 | µg/L | 20 | 0 | 82.8 | 73 | 126 | 17.9 | 7.84 | 25 |
| 1,2,4-Trichlorobenzene | 16.55 | 2.0 | µg/L | 20 | 0 | 82.8 | 77 | 134 | 17.35 | 4.72 | 25 |
| Hexachlorobutadiene | 16.75 | 5.0 | µg/L | 20 | 0 | 83.8 | 58 | 138 | 17.05 | 1.78 | 25 |
| Naphthalene | 15.69 | 2.0 | µg/L | 20 | 0 | 78.5 | 76 | 124 | 16.93 | 7.6 | 25 |
| 1,2,3-Trichlorobenzene | 25.74 | 2.0 | µg/L | 25 | 0 | 103 | 85 | 116 | 0 | 0 | 0 |
| Surr: Dibromofluoromethane | 26.86 | 2.0 | µg/L | 25 | 0 | 107 | 77 | 127 | 0 | 0 | 0 |
| Surr: 1,2-Dichloroethane-d4 | 25.15 | 2.0 | µg/L | 25 | 0 | 101 | 86 | 114 | 0 | 0 | 0 |
| Surr: Toluene-d8 | 24.35 | 2.0 | µg/L | 25 | 0 | 97.4 | 79 | 117 | 0 | 0 | 0 |
| Surr: 4-Bromofluorobenzene | | | | | | | | | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Matrix Spike - Full List

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

Sample ID: 0702072-15Amsf Batch ID: R35964 Test Code: SW6260B Units: µg/L Analysis Date 2/22/2007 10:05:00 PM Prep Date: 2/17/2007
Client ID: MW-209D Run ID: V-1_070222A SeqNo: 596251

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | QC |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|----|
| Dichlorodifluoromethane | 179.5 | 50 | µg/L | 200 | 0 | 89.8 | 16 | 150 | 0 | | | |
| Chloromethane | 183.6 | 50 | µg/L | 200 | 0 | 91.8 | 35 | 150 | 0 | | | |
| Vinyl chloride | 210.2 | 20 | µg/L | 200 | 0 | 105 | 49 | 150 | 0 | | | |
| Chloroethane | 196.3 | 50 | µg/L | 200 | 0 | 98.2 | 58 | 147 | 0 | | | |
| Bromomethane | 188.6 | 20 | µg/L | 200 | 0 | 94.3 | 49 | 142 | 0 | | | |
| Trichlorofluoromethane | 215.4 | 20 | µg/L | 200 | 0 | 108 | 57 | 149 | 0 | | | |
| Diethyl ether | 176.1 | 50 | µg/L | 200 | 0 | 88 | 66 | 136 | 0 | | | |
| Acetone | 140.3 | 100 | µg/L | 200 | 0 | 70.2 | 16 | 150 | 0 | | | |
| 1,1-Dichloroethene | 184.2 | 10 | µg/L | 200 | 0 | 92.1 | 70 | 150 | 0 | | | |
| Carbon disulfide | 202.5 | 20 | µg/L | 200 | 0 | 101 | 47 | 135 | 0 | | | |
| Methylene chloride | 213.5 | 50 | µg/L | 200 | 0 | 107 | 66 | 142 | 0 | | | |
| Methyl tert-butyl ether | 210.6 | 20 | µg/L | 200 | 0 | 105 | 63 | 138 | 0 | | | |
| trans-1,2-Dichloroethene | 227.8 | 20 | µg/L | 200 | 0 | 114 | 78 | 135 | 0 | | | |
| 1,1-Dichloroethane | 221.6 | 20 | µg/L | 200 | 0 | 111 | 76 | 131 | 0 | | | |
| 2-Butanone | 123.5 | 100 | µg/L | 200 | 0 | 61.8 | 51 | 142 | 0 | | | |
| 2,2-Dichloropropane | 217.9 | 20 | µg/L | 200 | 0 | 109 | 60 | 149 | 0 | | | |
| cis-1,2-Dichloroethene | 196.5 | 20 | µg/L | 200 | 0 | 98.2 | 74 | 128 | 0 | | | |
| Chloroform | 194.4 | 20 | µg/L | 200 | 0 | 97.2 | 80 | 129 | 0 | | | |
| Tetrahydrofuran | 144.7 | 100 | µg/L | 200 | 0 | 72.4 | 53 | 145 | 0 | | | |
| Bromochloromethane | 177.2 | 20 | µg/L | 200 | 0 | 88.6 | 78 | 130 | 0 | | | |
| 1,1,1-Trichloroethane | 222.1 | 20 | µg/L | 200 | 0 | 111 | 77 | 139 | 0 | | | |
| 1,1-Dichloropropene | 185.4 | 20 | µg/L | 200 | 0 | 92.7 | 74 | 127 | 0 | | | |
| Carbon tetrachloride | 184.6 | 20 | µg/L | 200 | 0 | 92.3 | 73 | 138 | 0 | | | |
| 1,2-Dichloroethane | 205.2 | 20 | µg/L | 200 | 0 | 103 | 75 | 130 | 0 | | | |
| Benzene | 202.1 | 10 | µg/L | 200 | 0 | 101 | 79 | 123 | 0 | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike - Full List

| Compound | Reporting Limit | Concentration (µg/L) | Recovery (%) | Recovery Outside Limits | Accepted Recovery Limits | Method Blank |
|---------------------------|-----------------|----------------------|--------------|-------------------------|--------------------------|--------------|
| Trichloroethene | 316.2 | 20 | 110.5 | 103 | 79 | 126 |
| 1,2-Dichloropropane | 193.4 | 20 | 0 | 96.7 | 76 | 125 |
| Bromodichloromethane | 189.5 | 20 | 0 | 94.8 | 69 | 119 |
| Dibromomethane | 190.1 | 20 | 0 | 95 | 76 | 127 |
| 4-Methyl-2-pentanone | 146.8 | 100 | 0 | 73.4 | 53 | 141 |
| cis-1,3-Dichloropropene | 176.5 | 10 | 0 | 88.2 | 70 | 119 |
| Toluene | 205.1 | 20 | 0 | 103 | 82 | 124 |
| trans-1,3-Dichloropropene | 172.4 | 10 | 0 | 86.2 | 64 | 124 |
| 1,1,2-Trichloroethane | 184.4 | 20 | 0 | 92.2 | 73 | 127 |
| 1,2-Dibromoethane | 188.7 | 20 | 0 | 94.4 | 73 | 127 |
| 2-Hexanone | 130 | 100 | 0 | 65 | 37 | 145 |
| 1,3-Dichloropropane | 181.8 | 20 | 0 | 90.9 | 76 | 123 |
| Tetrachloroethene | 584.2 | 20 | 431.8 | 76.2 | 82 | 129 |
| Dibromochloromethane | 175.7 | 20 | 0 | 87.8 | 59 | 125 |
| Chlorobenzene | 187.5 | 20 | 0 | 93.8 | 80 | 120 |
| 1,1,1,2-Tetrachloroethane | 183.7 | 20 | 0 | 91.8 | 72 | 124 |
| Ethylbenzene | 199.4 | 20 | 0 | 99.7 | 83 | 123 |
| m,p-Xylene | 405.6 | 20 | 0 | 101 | 84 | 121 |
| o-Xylene | 188.3 | 20 | 0 | 94.2 | 83 | 119 |
| Styrene | 192.9 | 20 | 0 | 96.5 | 80 | 122 |
| Bromoform | 139.7 | 20 | 0 | 69.8 | 54 | 119 |
| Isopropylbenzene | 212.8 | 20 | 0 | 106 | 75 | 131 |
| 1,1,2,2-Tetrachloroethane | 150.8 | 20 | 0 | 75.4 | 61 | 139 |
| 1,2,3-Trichloropropane | 157.1 | 20 | 0 | 78.6 | 66 | 130 |
| Bromobenzene | 175.6 | 20 | 0 | 87.8 | 77 | 124 |
| n-Propylbenzene | 199.4 | 20 | 0 | 99.7 | 76 | 131 |
| 2-Chlorotoluene | 185.3 | 20 | 0 | 92.6 | 78 | 125 |
| 4-Chlorotoluene | 190.8 | 20 | 0 | 95.4 | 75 | 124 |
| 1,3,5-Trimethylbenzene | 200.4 | 20 | 0 | 100 | 79 | 124 |
| tert-Butylbenzene | 203.4 | 20 | 0 | 102 | 79 | 126 |
| 1,2,4-Trimethylbenzene | 193.7 | 20 | 0 | 96.8 | 77 | 124 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike - Full List

| Compound | Concentration (µg/L) | Volume (µg/L) | Recovery (%) | Acceptance | Matrix Spike |
|-----------------------------|----------------------|---------------|--------------|------------|--------------|
| sec-Butylbenzene | 172.8 | 20 | 86.4 | 0 | 128 |
| 4-Isopropyltoluene | 195 | 20 | 97.5 | 0 | 128 |
| 1,3-Dichlorobenzene | 173.8 | 20 | 86.9 | 0 | 122 |
| 1,4-Dichlorobenzene | 168.8 | 20 | 84.4 | 0 | 123 |
| n-Butylbenzene | 159 | 20 | 79.5 | 0 | 130 |
| 1,2-Dichlorobenzene | 171.2 | 20 | 85.6 | 0 | 121 |
| 1,2-Dibromo-3-chloropropane | 128 | 50 | 64 | 0 | 127 |
| 1,2,4-Trichlorobenzene | 133.7 | 20 | 66.8 | 0 | 128 |
| Hexachlorobutadiene | 142.8 | 20 | 71.4 | 0 | 134 |
| Naphthalene | 130.6 | 50 | 65.3 | 0 | 131 |
| 1,2,3-Trichlorobenzene | 120.7 | 20 | 60.4 | 0 | 131 |
| Surr: Dibromofluoromethane | 252 | 20 | 101 | 0 | 116 |
| Surr: 1,2-Dichloroethane-d4 | 261.6 | 20 | 105 | 0 | 127 |
| Surr: Toluene-d8 | 255.4 | 20 | 102 | 0 | 114 |
| Surr: 4-Bromofluorobenzene | 242.9 | 20 | 97.2 | 0 | 117 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike Duplicate - Full List

Sample ID: 0702072-15Amsdf Batch ID: R35964 Test Code: SW6260B Units: µg/L Analysis Date 2/22/2007 10:39:00 PM Prep Date: 2/17/2007
 Client ID: MW-209D Run ID: V-1_070222A SeqNo: 596252

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | 198.3 | 50 | µg/L | 200 | 0 | 99.2 | 16 | 150 | 179.5 | 9.95 | 20 | |
| Chloromethane | 195.2 | 50 | µg/L | 200 | 0 | 97.6 | 35 | 150 | 183.6 | 6.12 | 20 | |
| Vinyl chloride | 222 | 20 | µg/L | 200 | 0 | 111 | 49 | 150 | 210.2 | 5.46 | 20 | |
| Chloroethane | 204.5 | 50 | µg/L | 200 | 0 | 102 | 58 | 147 | 196.3 | 4.09 | 20 | |
| Bromomethane | 203.4 | 20 | µg/L | 200 | 0 | 102 | 49 | 142 | 188.6 | 7.55 | 20 | |
| Trichlorofluoromethane | 234.4 | 20 | µg/L | 200 | 0 | 117 | 57 | 149 | 215.4 | 8.45 | 20 | |
| Diethyl ether | 192.4 | 50 | µg/L | 200 | 0 | 96.2 | 66 | 136 | 176.1 | 8.85 | 20 | |
| Acetone | 161.7 | 100 | µg/L | 200 | 0 | 80.8 | 16 | 150 | 140.3 | 14.2 | 20 | |
| 1,1-Dichloroethene | 209.7 | 10 | µg/L | 200 | 0 | 105 | 70 | 150 | 184.2 | 12.9 | 20 | |
| Carbon disulfide | 219.2 | 20 | µg/L | 200 | 0 | 110 | 47 | 135 | 202.5 | 7.92 | 20 | |
| Methylene chloride | 231.7 | 50 | µg/L | 200 | 0 | 116 | 66 | 142 | 213.5 | 8.18 | 20 | |
| Methyl tert-butyl ether | 228.8 | 20 | µg/L | 200 | 0 | 114 | 63 | 138 | 210.6 | 8.28 | 20 | |
| trans-1,2-Dichloroethene | 233.8 | 20 | µg/L | 200 | 0 | 117 | 78 | 135 | 227.8 | 2.6 | 20 | |
| 1,1-Dichloroethane | 239.2 | 20 | µg/L | 200 | 0 | 120 | 76 | 131 | 221.6 | 7.64 | 20 | |
| 2-Butanone | 145.3 | 100 | µg/L | 200 | 0 | 72.7 | 51 | 142 | 123.5 | 16.2 | 20 | |
| 2,2-Dichloropropane | 226.6 | 20 | µg/L | 200 | 0 | 113 | 60 | 149 | 217.9 | 3.91 | 20 | |
| cis-1,2-Dichloroethene | 203.8 | 20 | µg/L | 200 | 0 | 102 | 74 | 128 | 196.5 | 3.65 | 20 | |
| Chloroform | 208.9 | 20 | µg/L | 200 | 0 | 104 | 80 | 129 | 194.4 | 7.19 | 20 | |
| Tetrahydrofuran | 156.7 | 100 | µg/L | 200 | 0 | 78.4 | 53 | 145 | 144.7 | 7.96 | 20 | |
| Bromochloromethane | 190 | 20 | µg/L | 200 | 0 | 95 | 78 | 130 | 177.2 | 6.97 | 20 | |
| 1,1,1-Trichloroethane | 234.6 | 20 | µg/L | 200 | 0 | 117 | 77 | 139 | 222.1 | 5.47 | 20 | |
| 1,1-Dichloropropene | 192 | 20 | µg/L | 200 | 0 | 96 | 74 | 127 | 185.4 | 3.5 | 20 | |
| Carbon tetrachloride | 190 | 20 | µg/L | 200 | 0 | 95 | 73 | 138 | 184.6 | 2.88 | 20 | |
| 1,2-Dichloroethane | 218.7 | 20 | µg/L | 200 | 0 | 109 | 75 | 130 | 205.2 | 6.37 | 20 | |
| Benzene | 214.7 | 10 | µg/L | 200 | 0 | 107 | 79 | 123 | 202.1 | 6.05 | 20 | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike Duplicate - Full List

| Compound | Reporting Limit | Concentration | Recovery | Recovery Limits | Method | Matrix Spike Duplicate | Matrix Spike Duplicate |
|---------------------------|-----------------|---------------|----------|-----------------|--------|------------------------|------------------------|
| Trichloroethene | 331.8 | 20 | µg/L | 200 | 111 | 79 | 126 |
| 1,2-Dichloropropane | 216.5 | 20 | µg/L | 200 | 0 | 76 | 125 |
| Bromodichloromethane | 201.7 | 20 | µg/L | 200 | 0 | 69 | 119 |
| Dibromomethane | 191.7 | 20 | µg/L | 200 | 0 | 76 | 127 |
| 4-Methyl-2-pentanone | 154.6 | 100 | µg/L | 200 | 0 | 53 | 141 |
| cis-1,3-Dichloropropene | 189.6 | 10 | µg/L | 200 | 0 | 70 | 119 |
| Toluene | 217.9 | 20 | µg/L | 200 | 0 | 82 | 124 |
| trans-1,3-Dichloropropene | 185.3 | 10 | µg/L | 200 | 0 | 64 | 124 |
| 1,1,2-Trichloroethane | 197.5 | 20 | µg/L | 200 | 0 | 73 | 127 |
| 1,2-Dibromoethane | 198.4 | 20 | µg/L | 200 | 0 | 73 | 127 |
| 2-Hexanone | 140.8 | 100 | µg/L | 200 | 0 | 37 | 145 |
| 1,3-Dichloropropane | 194.1 | 20 | µg/L | 200 | 0 | 76 | 123 |
| Tetrachloroethene | 627.4 | 20 | µg/L | 200 | 431.8 | 82 | 129 |
| Dibromochloromethane | 181.7 | 20 | µg/L | 200 | 0 | 59 | 125 |
| Chlorobenzene | 195.1 | 20 | µg/L | 200 | 0 | 80 | 120 |
| 1,1,1,2-Tetrachloroethane | 192 | 20 | µg/L | 200 | 0 | 72 | 124 |
| Ethylbenzene | 215.1 | 20 | µg/L | 200 | 0 | 83 | 123 |
| m,p-Xylene | 431.4 | 20 | µg/L | 400 | 0 | 84 | 121 |
| o-Xylene | 200.3 | 20 | µg/L | 200 | 0 | 83 | 119 |
| Styrene | 207.6 | 20 | µg/L | 200 | 0 | 80 | 122 |
| Bromoform | 156.7 | 20 | µg/L | 200 | 0 | 54 | 119 |
| Isopropylbenzene | 235 | 20 | µg/L | 200 | 0 | 75 | 131 |
| 1,1,2,2-Tetrachloroethane | 168.7 | 20 | µg/L | 200 | 0 | 61 | 139 |
| 1,2,3-Trichloropropane | 175.7 | 20 | µg/L | 200 | 0 | 66 | 130 |
| Bromobenzene | 193.7 | 20 | µg/L | 200 | 0 | 77 | 124 |
| n-Propylbenzene | 220 | 20 | µg/L | 200 | 0 | 76 | 131 |
| 2-Chlorotoluene | 207.5 | 20 | µg/L | 200 | 0 | 78 | 125 |
| 4-Chlorotoluene | 206.4 | 20 | µg/L | 200 | 0 | 75 | 124 |
| 1,3,5-Trimethylbenzene | 216 | 20 | µg/L | 200 | 0 | 79 | 124 |
| tert-Butylbenzene | 222.7 | 20 | µg/L | 200 | 0 | 79 | 126 |
| 1,2,4-Trimethylbenzene | 207.7 | 20 | µg/L | 200 | 0 | 77 | 124 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike Duplicate - Full List

| Compound | Concentration (µg/L) | Volume (L) | Recovery (%) | Accepted Recovery Limits | Sample ID | Concentration (µg/L) | Volume (L) | Recovery (%) | Accepted Recovery Limits | Sample ID |
|-----------------------------|----------------------|------------|--------------|--------------------------|-----------|----------------------|------------|--------------|--------------------------|-----------|
| sec-Butylbenzene | 188.4 | 20 | 0 | 0 | 82 | 172.8 | 20 | 94.2 | 0 | 128 |
| 4-Isopropyltoluene | 214.6 | 20 | 0 | 0 | 77 | 195 | 20 | 107 | 0 | 128 |
| 1,3-Dichlorobenzene | 186.2 | 20 | 0 | 0 | 80 | 173.8 | 20 | 93.1 | 0 | 122 |
| 1,4-Dichlorobenzene | 181 | 20 | 0 | 0 | 78 | 168.8 | 20 | 90.5 | 0 | 123 |
| n-Butylbenzene | 172.2 | 20 | 0 | 0 | 74 | 159 | 20 | 86.1 | 0 | 130 |
| 1,2-Dichlorobenzene | 182.8 | 20 | 0 | 0 | 78 | 171.2 | 20 | 91.4 | 0 | 121 |
| 1,2-Dibromo-3-chloropropane | 139.1 | 50 | 0 | 0 | 50 | 128 | 20 | 69.6 | 0 | 127 |
| 1,2,4-Trichlorobenzene | 162.5 | 20 | 0 | 0 | 67 | 133.7 | 20 | 81.2 | 0 | 128 |
| Hexachlorobutadiene | 166 | 20 | 0 | 0 | 74 | 142.8 | 20 | 83 | 0 | 134 |
| Naphthalene | 158.6 | 50 | 0 | 0 | 57 | 130.6 | 20 | 79.3 | 0 | 131 |
| 1,2,3-Trichlorobenzene | 147 | 20 | 0 | 0 | 64 | 120.7 | 20 | 73.5 | 0 | 131 |
| Surr: Dibromofluoromethane | 249.3 | 20 | 0 | 0 | 85 | 0 | 20 | 99.7 | 0 | 116 |
| Surr: 1,2-Dichloroethane-d4 | 271 | 20 | 0 | 0 | 77 | 0 | 20 | 108 | 0 | 127 |
| Surr: Toluene-d8 | 253.8 | 20 | 0 | 0 | 86 | 0 | 20 | 102 | 0 | 114 |
| Surr: 4-Bromofluorobenzene | 249.1 | 20 | 0 | 0 | 79 | 0 | 20 | 99.6 | 0 | 117 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike - Full List

Sample ID: 0702072-19Amsf Batch ID: R35993

Test Code: SW8260B Units: µg/L

Analysis Date 2/23/2007 6:20:00 PM Prep Date: 2/17/2007

Client ID: CW-1

Run ID: V-1_070223A

SeqNo: 596559

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-------|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | 1581 | 500 | µg/L | 2000 | 0 | 79 | 16 | 150 | 0 | | | |
| Chloromethane | 1769 | 500 | µg/L | 2000 | 0 | 88.4 | 35 | 150 | 0 | | | |
| Vinyl chloride | 1983 | 200 | µg/L | 2000 | 0 | 99.2 | 49 | 150 | 0 | | | |
| Chloroethane | 1976 | 500 | µg/L | 2000 | 0 | 98.8 | 58 | 147 | 0 | | | |
| Bromomethane | 1922 | 200 | µg/L | 2000 | 0 | 96.1 | 49 | 142 | 0 | | | |
| Trichlorofluoromethane | 2221 | 200 | µg/L | 2000 | 0 | 111 | 57 | 149 | 0 | | | |
| Diethyl ether | 1840 | 500 | µg/L | 2000 | 0 | 92 | 66 | 136 | 0 | | | |
| Acetone | 1338 | 1,000 | µg/L | 2000 | 0 | 66.9 | 16 | 150 | 0 | | | |
| 1,1-Dichloroethene | 1983 | 100 | µg/L | 2000 | 177 | 90.3 | 70 | 150 | 0 | | | |
| Carbon disulfide | 1891 | 200 | µg/L | 2000 | 0 | 94.6 | 47 | 135 | 0 | | | |
| Methylene chloride | 2150 | 500 | µg/L | 2000 | 0 | 108 | 66 | 142 | 0 | | | |
| Methyl tert-butyl ether | 2045 | 200 | µg/L | 2000 | 0 | 102 | 63 | 138 | 0 | | | |
| trans-1,2-Dichloroethene | 2206 | 200 | µg/L | 2000 | 0 | 110 | 78 | 135 | 0 | | | |
| 1,1-Dichloroethane | 2235 | 200 | µg/L | 2000 | 0 | 112 | 76 | 131 | 0 | | | |
| 2-Butanone | 1331 | 1,000 | µg/L | 2000 | 0 | 66.6 | 51 | 142 | 0 | | | |
| 2,2-Dichloropropane | 2473 | 200 | µg/L | 2000 | 0 | 124 | 60 | 149 | 0 | | | |
| cis-1,2-Dichloroethene | 2230 | 200 | µg/L | 2000 | 352 | 93.9 | 74 | 128 | 0 | | | |
| Chloroform | 2051 | 200 | µg/L | 2000 | 0 | 103 | 80 | 129 | 0 | | | |
| Tetrahydrofuran | 1451 | 1,000 | µg/L | 2000 | 0 | 72.6 | 53 | 145 | 0 | | | |
| Bromochloromethane | 1803 | 200 | µg/L | 2000 | 0 | 90.2 | 78 | 130 | 0 | | | |
| 1,1,1-Trichloroethane | 2216 | 200 | µg/L | 2000 | 0 | 111 | 77 | 139 | 0 | | | |
| 1,1-Dichloropropene | 1863 | 200 | µg/L | 2000 | 0 | 93.2 | 74 | 127 | 0 | | | |
| Carbon tetrachloride | 1901 | 200 | µg/L | 2000 | 0 | 95 | 73 | 138 | 0 | | | |
| 1,2-Dichloroethane | 2109 | 200 | µg/L | 2000 | 0 | 105 | 75 | 130 | 0 | | | |
| Benzene | 1961 | 100 | µg/L | 2000 | 0 | 98 | 79 | 123 | 0 | | | |

Qualifiers: ND - Not Detected at the Reporting Limit

S - Spike Recovery outside accepted recovery limits

B - Analyte detected in the associated Method Blank

J - Analyte detected below quantitation limits

R - RPD outside accepted recovery limits

NA - Not applicable where J values or ND results occur

RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike - Full List

| Compound | Reporting Limit | Concentration (µg/L) | Recovery | Acceptance | Recovery Limits | Spikes | Spikes (µg/L) | Spikes (µg/L) | Spikes (µg/L) |
|---------------------------|-----------------|----------------------|----------|------------|-----------------|--------|---------------|---------------|---------------|
| Trichloroethene | 7887 | 200 | µg/L | 2000 | 5939 | 79 | 97.4 | 126 | 0 |
| 1,2-Dichloropropane | 1986 | 200 | µg/L | 2000 | 0 | 76 | 99.3 | 125 | 0 |
| Bromodichloromethane | 1878 | 200 | µg/L | 2000 | 0 | 69 | 93.9 | 119 | 0 |
| Dibromomethane | 1849 | 200 | µg/L | 2000 | 0 | 76 | 92.5 | 127 | 0 |
| 4-Methyl-2-pentanone | 1296 | 1,000 | µg/L | 2000 | 0 | 53 | 64.8 | 141 | 0 |
| cis-1,3-Dichloropropene | 1856 | 100 | µg/L | 2000 | 0 | 70 | 92.8 | 119 | 0 |
| Toluene | 2050 | 200 | µg/L | 2000 | 0 | 82 | 103 | 124 | 0 |
| trans-1,3-Dichloropropene | 1805 | 100 | µg/L | 2000 | 0 | 64 | 90.2 | 124 | 0 |
| 1,1,2-Trichloroethane | 1905 | 200 | µg/L | 2000 | 0 | 73 | 95.2 | 127 | 0 |
| 1,2-Dibromoethane | 1922 | 200 | µg/L | 2000 | 0 | 73 | 96.1 | 127 | 0 |
| 2-Hexanone | 1311 | 1,000 | µg/L | 2000 | 0 | 37 | 65.6 | 145 | 0 |
| 1,3-Dichloropropane | 1834 | 200 | µg/L | 2000 | 0 | 76 | 91.7 | 123 | 0 |
| Tetrachloroethane | 2144 | 200 | µg/L | 2000 | 0 | 82 | 107 | 129 | 0 |
| Dibromochloromethane | 1767 | 200 | µg/L | 2000 | 0 | 59 | 88.4 | 125 | 0 |
| Chlorobenzene | 1835 | 200 | µg/L | 2000 | 0 | 80 | 91.8 | 120 | 0 |
| 1,1,1,2-Tetrachloroethane | 1782 | 200 | µg/L | 2000 | 0 | 72 | 89.1 | 124 | 0 |
| Ethylbenzene | 1944 | 200 | µg/L | 2000 | 0 | 83 | 97.2 | 123 | 0 |
| m,p-Xylene | 3889 | 200 | µg/L | 4000 | 0 | 84 | 97.2 | 121 | 0 |
| o-Xylene | 1877 | 200 | µg/L | 2000 | 0 | 83 | 93.8 | 119 | 0 |
| Styrene | 1976 | 200 | µg/L | 2000 | 0 | 80 | 98.8 | 122 | 0 |
| Bromoform | 1496 | 200 | µg/L | 2000 | 0 | 54 | 74.8 | 119 | 0 |
| Isopropylbenzene | 2081 | 200 | µg/L | 2000 | 0 | 75 | 104 | 131 | 0 |
| 1,1,2,2-Tetrachloroethane | 1564 | 200 | µg/L | 2000 | 0 | 61 | 78.2 | 139 | 0 |
| 1,2,3-Trichloropropane | 1645 | 200 | µg/L | 2000 | 0 | 66 | 82.2 | 130 | 0 |
| Bromobenzene | 1799 | 200 | µg/L | 2000 | 0 | 77 | 90 | 124 | 0 |
| n-Propylbenzene | 1959 | 200 | µg/L | 2000 | 0 | 76 | 98 | 131 | 0 |
| 2-Chlorotoluene | 1911 | 200 | µg/L | 2000 | 0 | 78 | 95.6 | 125 | 0 |
| 4-Chlorotoluene | 1934 | 200 | µg/L | 2000 | 0 | 75 | 96.7 | 124 | 0 |
| 1,3,5-Trimethylbenzene | 1940 | 200 | µg/L | 2000 | 0 | 79 | 97 | 124 | 0 |
| tert-Butylbenzene | 2002 | 200 | µg/L | 2000 | 0 | 79 | 100 | 126 | 0 |
| 1,2,4-Trimethylbenzene | 1947 | 200 | µg/L | 2000 | 0 | 77 | 97.4 | 124 | 0 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank

J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur

RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike - Full List

| Compound | Reporting Limit | Concentration (µg/L) | Recovery (%) | Acceptance | Notes |
|-----------------------------|-----------------|----------------------|--------------|------------|-------|
| sec-Butylbenzene | 1694 | 200 | 84.7 | 0 | 128 |
| 4-Isopropyltoluene | 1920 | 200 | 96 | 0 | 128 |
| 1,3-Dichlorobenzene | 1802 | 200 | 90.1 | 0 | 122 |
| 1,4-Dichlorobenzene | 1781 | 200 | 89.1 | 0 | 123 |
| n-Butylbenzene | 1563 | 200 | 78.2 | 0 | 130 |
| 1,2-Dichlorobenzene | 1749 | 200 | 87.5 | 0 | 121 |
| 1,2-Dibromo-3-chloropropane | 1116 | 500 | 55.8 | 0 | 127 |
| 1,2,4-Trichlorobenzene | 1397 | 200 | 69.8 | 0 | 128 |
| Hexachlorobutadiene | 1388 | 200 | 69.4 | 0 | 134 |
| Naphthalene | 1305 | 500 | 65.2 | 0 | 131 |
| 1,2,3-Trichlorobenzene | 1270 | 200 | 63.5 | 0 | 131 |
| Surr: Dibromofluoromethane | 2548 | 200 | 102 | 0 | 116 |
| Surr: 1,2-Dichloroethane-d4 | 2635 | 200 | 105 | 0 | 127 |
| Surr: Toluene-d8 | 2563 | 200 | 103 | 0 | 114 |
| Surr: 4-Bromofluorobenzene | 2493 | 200 | 99.7 | 0 | 117 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT
Matrix Spike Duplicate - Full List

Sample ID: 0702072-19Amsdf Batch ID: R35993

Test Code: SW8260B Units: µg/L

Analysis Date 2/23/2007 6:54:00 PM

Prep Date: 2/17/2007

Client ID: CW-1

Run ID: V-1_070223A

SeqNo: 596561

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Qua |
|--------------------------|------------------|-------|-------|-----------------|------------------------|------|----------|-----------|------------------------------|-------|----------|-----|
| Dichlorodifluoromethane | 1697 | 500 | µg/L | 2000 | 0 | 84.8 | 16 | 150 | 1581 | 7.08 | 20 | |
| Chloromethane | 1805 | 500 | µg/L | 2000 | 0 | 90.2 | 35 | 150 | 1769 | 2.01 | 20 | |
| Vinyl chloride | 2084 | 200 | µg/L | 2000 | 0 | 104 | 49 | 150 | 1983 | 4.97 | 20 | |
| Chloroethane | 1962 | 500 | µg/L | 2000 | 0 | 98.1 | 58 | 147 | 1976 | 0.711 | 20 | |
| Bromomethane | 1909 | 200 | µg/L | 2000 | 0 | 95.4 | 49 | 142 | 1922 | 0.679 | 20 | |
| Trichlorofluoromethane | 2260 | 200 | µg/L | 2000 | 0 | 113 | 57 | 149 | 2221 | 1.74 | 20 | |
| Diethyl ether | 1883 | 500 | µg/L | 2000 | 0 | 94.2 | 66 | 136 | 1840 | 2.31 | 20 | |
| Acetone | 1457 | 1,000 | µg/L | 2000 | 0 | 72.8 | 16 | 150 | 1338 | 8.52 | 20 | |
| 1,1-Dichloroethene | 2145 | 100 | µg/L | 2000 | 177 | 98.4 | 70 | 150 | 1983 | 7.85 | 20 | |
| Carbon disulfide | 2058 | 200 | µg/L | 2000 | 0 | 103 | 47 | 135 | 1891 | 8.46 | 20 | |
| Methylene chloride | 2222 | 500 | µg/L | 2000 | 0 | 111 | 66 | 142 | 2150 | 3.29 | 20 | |
| Methyl tert-butyl ether | 2128 | 200 | µg/L | 2000 | 0 | 106 | 63 | 138 | 2045 | 3.98 | 20 | |
| trans-1,2-Dichloroethene | 2336 | 200 | µg/L | 2000 | 0 | 117 | 78 | 135 | 2206 | 5.72 | 20 | |
| 1,1-Dichloroethane | 2437 | 200 | µg/L | 2000 | 0 | 122 | 76 | 131 | 2235 | 8.65 | 20 | |
| 2-Butanone | 1456 | 1,000 | µg/L | 2000 | 0 | 72.8 | 51 | 142 | 1331 | 8.97 | 20 | |
| 2,2-Dichloropropane | 2564 | 200 | µg/L | 2000 | 0 | 128 | 60 | 149 | 2473 | 3.61 | 20 | |
| cis-1,2-Dichloroethene | 2418 | 200 | µg/L | 2000 | 352 | 103 | 74 | 128 | 2230 | 8.09 | 20 | |
| Chloroform | 2097 | 200 | µg/L | 2000 | 0 | 105 | 80 | 129 | 2051 | 2.22 | 20 | |
| Tetrahydrofuran | 1545 | 1,000 | µg/L | 2000 | 0 | 77.2 | 53 | 145 | 1451 | 6.28 | 20 | |
| Bromochloromethane | 1974 | 200 | µg/L | 2000 | 0 | 98.7 | 78 | 130 | 1803 | 9.05 | 20 | |
| 1,1,1-Trichloroethane | 2404 | 200 | µg/L | 2000 | 0 | 120 | 77 | 139 | 2216 | 8.14 | 20 | |
| 1,1-Dichloropropene | 2026 | 200 | µg/L | 2000 | 0 | 101 | 74 | 127 | 1863 | 8.38 | 20 | |
| Carbon tetrachloride | 2046 | 200 | µg/L | 2000 | 0 | 102 | 73 | 138 | 1901 | 7.35 | 20 | |
| 1,2-Dichloroethane | 2155 | 200 | µg/L | 2000 | 0 | 108 | 75 | 130 | 2109 | 2.16 | 20 | |
| Benzene | 2135 | 100 | µg/L | 2000 | 0 | 107 | 79 | 123 | 1961 | 8.5 | 20 | |

Qualifiers: ND - Not Detected at the Reporting Limit
 J - Analyte detected below quantitation limits
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

S - Spike Recovery outside accepted recovery limits
 R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank
 NA - Not applicable where J values or ND results occur

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike Duplicate - Full List

| Compound | Reporting Limit | Concentration | Recovery | Acceptance | Recovery | Acceptance | Concentration | Recovery | Acceptance | Concentration | Recovery | Acceptance |
|---------------------------|-----------------|---------------|----------|------------|----------|------------|---------------|----------|------------|---------------|----------|------------|
| Trichloroethene | 8359 | 200 | µg/L | 2000 | 5939 | 121 | 79 | 126 | 7887 | 5.81 | 20 | 20 |
| 1,2-Dichloropropane | 2115 | 200 | µg/L | 2000 | 0 | 106 | 76 | 125 | 1986 | 6.29 | 20 | 20 |
| Bromodichloromethane | 2039 | 200 | µg/L | 2000 | 0 | 102 | 69 | 119 | 1878 | 8.22 | 20 | 20 |
| Dibromomethane | 1926 | 200 | µg/L | 2000 | 0 | 96.3 | 76 | 127 | 1849 | 4.08 | 20 | 20 |
| 4-Methyl-2-pentanone | 1428 | 1,000 | µg/L | 2000 | 0 | 71.4 | 53 | 141 | 1296 | 9.69 | 20 | 20 |
| cis-1,3-Dichloropropene | 1954 | 100 | µg/L | 2000 | 0 | 97.7 | 70 | 119 | 1856 | 5.14 | 20 | 20 |
| Toluene | 2172 | 200 | µg/L | 2000 | 0 | 109 | 82 | 124 | 2050 | 5.78 | 20 | 20 |
| trans-1,3-Dichloropropene | 1947 | 100 | µg/L | 2000 | 0 | 97.4 | 64 | 124 | 1805 | 7.57 | 20 | 20 |
| 1,1,2-Trichloroethane | 1965 | 200 | µg/L | 2000 | 0 | 98.2 | 73 | 127 | 1905 | 3.1 | 20 | 20 |
| 1,2-Dibromoethane | 1874 | 200 | µg/L | 2000 | 0 | 93.7 | 73 | 127 | 1922 | 2.53 | 20 | 20 |
| 2-Hexanone | 1327 | 1,000 | µg/L | 2000 | 0 | 66.4 | 37 | 145 | 1311 | 1.21 | 20 | 20 |
| 1,3-Dichloropropane | 1892 | 200 | µg/L | 2000 | 0 | 94.6 | 76 | 123 | 1834 | 3.11 | 20 | 20 |
| Tetrachloroethene | 2137 | 200 | µg/L | 2000 | 0 | 107 | 82 | 129 | 2144 | 0.327 | 20 | 20 |
| Dibromochloromethane | 1761 | 200 | µg/L | 2000 | 0 | 88 | 59 | 125 | 1767 | 0.34 | 20 | 20 |
| Chlorobenzene | 1941 | 200 | µg/L | 2000 | 0 | 97 | 80 | 120 | 1835 | 5.61 | 20 | 20 |
| 1,1,1,2-Tetrachloroethane | 1885 | 200 | µg/L | 2000 | 0 | 94.2 | 72 | 124 | 1782 | 5.62 | 20 | 20 |
| Ethylbenzene | 2089 | 200 | µg/L | 2000 | 0 | 104 | 83 | 123 | 1944 | 7.19 | 20 | 20 |
| m,p-Xylene | 4285 | 200 | µg/L | 4000 | 0 | 107 | 84 | 121 | 3889 | 9.69 | 20 | 20 |
| o-Xylene | 1965 | 200 | µg/L | 2000 | 0 | 98.2 | 83 | 119 | 1877 | 4.58 | 20 | 20 |
| Styrene | 2049 | 200 | µg/L | 2000 | 0 | 102 | 80 | 122 | 1976 | 3.63 | 20 | 20 |
| Bromoform | 1495 | 200 | µg/L | 2000 | 0 | 74.8 | 54 | 119 | 1496 | 0.0669 | 20 | 20 |
| Isopropylbenzene | 2258 | 200 | µg/L | 2000 | 0 | 113 | 75 | 131 | 2081 | 8.16 | 20 | 20 |
| 1,1,2,2-Tetrachloroethane | 1580 | 200 | µg/L | 2000 | 0 | 79 | 61 | 139 | 1564 | 1.02 | 20 | 20 |
| 1,2,3-Trichloropropane | 1638 | 200 | µg/L | 2000 | 0 | 81.9 | 66 | 130 | 1645 | 0.426 | 20 | 20 |
| Bromobenzene | 1842 | 200 | µg/L | 2000 | 0 | 92.1 | 77 | 124 | 1799 | 2.36 | 20 | 20 |
| n-Propylbenzene | 2123 | 200 | µg/L | 2000 | 0 | 106 | 76 | 131 | 1959 | 8.04 | 20 | 20 |
| 2-Chlorotoluene | 1984 | 200 | µg/L | 2000 | 0 | 99.2 | 78 | 125 | 1911 | 3.75 | 20 | 20 |
| 4-Chlorotoluene | 2033 | 200 | µg/L | 2000 | 0 | 102 | 75 | 124 | 1934 | 4.99 | 20 | 20 |
| 1,3,5-Trimethylbenzene | 2116 | 200 | µg/L | 2000 | 0 | 106 | 79 | 124 | 1940 | 8.68 | 20 | 20 |
| tert-Butylbenzene | 2140 | 200 | µg/L | 2000 | 0 | 107 | 79 | 126 | 2002 | 6.66 | 20 | 20 |
| 1,2,4-Trimethylbenzene | 2050 | 200 | µg/L | 2000 | 0 | 103 | 77 | 124 | 1947 | 5.15 | 20 | 20 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike Duplicate - Full List

| Chemical Name | 1827 | 2041 | 1812 | 1791 | 1707 | 1818 | 1385 | 1521 | 1641 | 1416 | 1410 | 2560 | 2632 | 2568 | 2451 |
|-----------------------------|------|------|-------|------|------|------|------|------|------|------|------|------|------|------|------|
| sec-Butylbenzene | 200 | 200 | 200 | 200 | 200 | 200 | 500 | 200 | 200 | 500 | 200 | 200 | 200 | 200 | 200 |
| 4-Isopropyltoluene | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L |
| 1,3-Dichlorobenzene | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2500 | 2500 | 2500 | 2500 |
| 1,4-Dichlorobenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| n-Butylbenzene | 91.4 | 102 | 90.6 | 89.6 | 85.4 | 90.9 | 69.2 | 76 | 82 | 70.8 | 70.5 | 102 | 105 | 103 | 98 |
| 1,2-Dichlorobenzene | 82 | 77 | 80 | 78 | 74 | 78 | 50 | 67 | 74 | 57 | 64 | 85 | 77 | 86 | 79 |
| 1,2-Dibromo-3-chloropropane | 128 | 128 | 122 | 123 | 130 | 121 | 127 | 128 | 134 | 131 | 131 | 116 | 127 | 114 | 117 |
| 1,2,4-Trichlorobenzene | 1694 | 1920 | 1802 | 1781 | 1563 | 1749 | 1116 | 1397 | 1388 | 1305 | 1270 | 0 | 0 | 0 | 0 |
| Hexachlorobutadiene | 7.55 | 6.11 | 0.553 | 0.56 | 8.81 | 3.87 | 21.5 | 8.5 | 16.7 | 8.16 | 10.4 | 0 | 0 | 0 | 0 |
| Naphthalene | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 0 | 0 | 0 | 0 |
| 1,2,3-Trichlorobenzene | R | | | | | | | | | | | | | | |
| Surr: Dibromofluoromethane | | | | | | | | | | | | | | | |
| Surr: 1,2-Dichloroethane-d4 | | | | | | | | | | | | | | | |
| Surr: Toluene-d8 | | | | | | | | | | | | | | | |
| Surr: 4-Bromofluorobenzene | | | | | | | | | | | | | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-23A

Client Sample ID: CW-6
Tag Number:
Collection Date: 2/17/07 1:45:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|---------------------------------------|--------|----------------|------|-------|----|--------------------|
| TPH BY GC/FID (MODIFIED 8015B) | | SW8015B | | | | Analyst: FQ |
| Gasoline | ND | 0.051 | | mg/L | 1 | 2/23/07 6:08:00 PM |
| Mineral Spirits | ND | 0.051 | | mg/L | 1 | 2/23/07 6:08:00 PM |
| Kerosene | ND | 0.051 | | mg/L | 1 | 2/23/07 6:08:00 PM |
| Diesel Fuel/Fuel Oil #2 | ND | 0.051 | | mg/L | 1 | 2/23/07 6:08:00 PM |
| Motor Oil/Hydraulic Oil | ND | 0.10 | | mg/L | 1 | 2/23/07 6:08:00 PM |
| Unidentified Hydrocarbons | 8.5 | 0.20 | | mg/L | 1 | 2/23/07 6:08:00 PM |
| Surr: o-Terphenyl | 63.5 | 31-131 | | %REC | 1 | 2/23/07 6:08:00 PM |

Gasoline cannot be accurately determined by this method. Purge and trap sample introduction into a GC or GCMS is the recommended approach for gasoline. Due to the physical, chemical, and biological processes which affect the chemical composition of fuel mixtures exposed to the environment, the qualitative identity of a hydrocarbon mixture as a fuel product is not always conclusive by this method due to the method's reliance on chromatographic pattern recognition. A result provided for a specific fuel indicates that the mixture present in the sample has a chromatographic pattern similar to the laboratory's reference standard for that fuel mixture under specific GC operating conditions utilized at the time of analysis. A result identified as Unidentified Hydrocarbons is based upon the detector response obtained for the laboratory's Fuel Oil#2 reference standard and includes the entire chromatographic response for the sample between n-Alkanes of carbon numbers C9 to C36.

Qualifiers:

| | |
|---|---|
| ND - Not Detected at the Reporting Limit | S - Spike Recovery outside accepted recovery limits |
| J - Analyte detected below quantitation limits | R - RPD outside accepted recovery limits |
| B - Analyte detected in the associated Method Blank | E - Value above quantitation range |
| H - Method prescribed holding time exceeded. | # - See Case Narrative |
| RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate. | |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-24A

Client Sample ID: CW-6 DUP
Tag Number:
Collection Date: 2/17/07 1:45:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|----------|--------|----|------|-------|----|---------------|
|----------|--------|----|------|-------|----|---------------|

| TPH BY GC/FID (MODIFIED 8015B) | SW8015B | | Analyst: FQ | | | |
|--------------------------------|---------|--------|-------------|------|---|--------------------|
| Gasoline | ND | 0.052 | | mg/L | 1 | 2/23/07 6:44:00 PM |
| Mineral Spirits | ND | 0.052 | | mg/L | 1 | 2/23/07 6:44:00 PM |
| Kerosene | ND | 0.052 | | mg/L | 1 | 2/23/07 6:44:00 PM |
| Diesel Fuel/Fuel Oil #2 | ND | 0.052 | | mg/L | 1 | 2/23/07 6:44:00 PM |
| Motor Oil/Hydraulic Oil | ND | 0.10 | | mg/L | 1 | 2/23/07 6:44:00 PM |
| Unidentified Hydrocarbons | 8.6 | 0.21 | | mg/L | 1 | 2/23/07 6:44:00 PM |
| Surr: o-Terphenyl | 66.3 | 31-131 | | %REC | 1 | 2/23/07 6:44:00 PM |

Gasoline cannot be accurately determined by this method. Purge and trap sample introduction into a GC or GCMS is the recommended approach for gasoline. Due to the physical, chemical, and biological processes which affect the chemical composition of fuel mixtures exposed to the environment, the qualitative identity of a hydrocarbon mixture as a fuel product is not always conclusive by this method due to the method's reliance on chromatographic pattern recognition. A result provided for a specific fuel indicates that the mixture present in the sample has a chromatographic pattern similar to the laboratory's reference standard for that fuel mixture under specific GC operating conditions utilized at the time of analysis. A result identified as Unidentified Hydrocarbons is based upon the detector response obtained for the laboratory's Fuel Oil#2 reference standard and includes the entire chromatographic response for the sample between n-Alkanes of carbon numbers C9 to C36.

Qualifiers:

| | |
|---|---|
| ND - Not Detected at the Reporting Limit | S - Spike Recovery outside accepted recovery limits |
| J - Analyte detected below quantitation limits | R - RPD outside accepted recovery limits |
| B - Analyte detected in the associated Method Blank | E - Value above quantitation range |
| H - Method prescribed holding time exceeded. | # - See Case Narrative |
| RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate. | |

Date: 28-Feb-07

AMRO Environmental Laboratories Corp.

QC SUMMARY REPORT Method Blank

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

Sample ID: MB-16813 Batch ID: 16813 Test Code: SW8015B Units: mg/L Analysis Date 2/23/2007 4:18:00 PM Prep Date: 2/23/2007
Client ID: Run ID: GC-FING1_070223A SeqNo: 596574

| Analyte | QC Sample Result | RL | Units | QC Spike Original Sample Amount | Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|---------------------------|------------------|-------|-------|---------------------------------|--------|------|----------|-----------|------------------------------|------|----------|-----|
| Gasoline | ND | 0.050 | mg/L | | | | | | | | | |
| Mineral Spirits | ND | 0.050 | mg/L | | | | | | | | | |
| Kerosene | ND | 0.050 | mg/L | | | | | | | | | |
| Diesel Fuel/Fuel Oil #2 | ND | 0.050 | mg/L | | | | | | | | | |
| Motor Oil/Hydraulic Oil | ND | 0.10 | mg/L | | | | | | | | | |
| Unidentified Hydrocarbons | ND | 0.20 | mg/L | | | | | | | | | |
| Surr: o-Terphenyl | 0.06958 | 0 | mg/L | 0.1 | 0 | 69.6 | 31 | 131 | 0 | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Laboratory Control Spike

Sample ID: LCS-16813 Batch ID: 16813 Test Code: SW8015B Units: mg/L Analysis Date 2/23/2007 4:55:00 PM Prep Date: 2/23/2007
 Client ID: Run ID: GC-FING1_070223A SeqNo: 596575

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|-------------------------|------------------|-------|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Diesel Fuel/Fuel Oil #2 | 1.236 | 0.050 | mg/L | 2 | 0 | 61.8 | 42 | 119 | 0 | | | |
| Surr: o-Terphenyl | 0.07476 | 0 | mg/L | 0.1 | 0 | 74.8 | 31 | 131 | 0 | | | |

Sample ID: LCSD-16813 Batch ID: 16813 Test Code: SW8015B Units: mg/L Analysis Date 2/23/2007 5:31:00 PM Prep Date: 2/23/2007
 Client ID: Run ID: GC-FING1_070223A SeqNo: 596576

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|-------------------------|------------------|-------|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Diesel Fuel/Fuel Oil #2 | 1.184 | 0.050 | mg/L | 2 | 0 | 59.2 | 42 | 119 | 1.236 | 4.29 | 40 | |
| Surr: o-Terphenyl | 0.06378 | 0 | mg/L | 0.1 | 0 | 63.8 | 31 | 131 | 0 | 0 | 0 | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.



March 19, 2007

ANALYTICAL TEST RESULTS

Ed VanDoren
SHAW E & I, Inc.
11 Northeastern Boulevard
Salem, NH 030791953
TEL: (603) 870-4500
FAX: (603) 870-4501

Subject: 101960 Textron Gorham

Workorder No.: 0702072

Dear Ed VanDoren:

AMRO Environmental Laboratories Corp. received 24 samples on 2/19/2007 for the analyses presented in the following report.

The enclosed sample results are revised based upon further review of the the analytical data or legitimate changes made at your request.

AMRO is accredited in accordance with NELAC and certifies that these test results meet all the requirements of NELAC, where applicable, unless otherwise noted in the case narrative.

Please be advised that any unused sample volume and sample extracts will be stored for a period of thirty (30) days from this report date. After this time, AMRO will properly dispose of the remaining sample(s). If you require further analysis, or need the samples held for a longer period, please contact us immediately.

This report consists of a total of 3 pages. This letter is an integral part of your data report. If you have any questions regarding this project in the future, please refer to the Order Number above.

Sincerely,

Nancy Stewart
Vice President

State Certifications: NH (NELAC): 1001, MA: M-NH012, CT: PH-0758, NY: 11278 (NELAC), ME: NH012 and 1001, NJ: NH125, RI: 00105, U.S. Army Corps of Engineers (USACE), Naval Facilities Engineering Service Center (NFESC).

Hard copy of the State Certification is available upon request.



AMRO Environmental Laboratories Corp.

Date: 19-Mar-07

CLIENT: SHAW E & I, Inc.
 Lab Order: 0702072
 Project: 101960 Textron Gorham
 Lab ID: 0702072-02A

Client Sample ID: MW-207D
 Collection Date: 2/17/2007 7:30:00 AM
 Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|----------------------|
| EPA 8260B VOLATILES BY GC/MS | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Vinyl chloride | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,1-Dichloroethene | ND | 100 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| cis-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Trichloroethene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Tetrachloroethene | 7,700 | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |

AMRO Environmental Laboratories Corp.

Date: 19-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-02A

Client Sample ID: MW-207D
Collection Date: 2/17/2007 7:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|----------------------|
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/23/2007 2:53:00 PM |
| Surr: Dibromofluoromethane | 98.4 | 85-116 | | %REC | 100 | 2/23/2007 2:53:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 104 | 77-127 | | %REC | 100 | 2/23/2007 2:53:00 PM |
| Surr: Toluene-d8 | 100 | 86-114 | | %REC | 100 | 2/23/2007 2:53:00 PM |
| Surr: 4-Bromofluorobenzene | 98.7 | 79-117 | | %REC | 100 | 2/23/2007 2:53:00 PM |



111 Herrick Street, Merrimack, NH 03054
TEL: (603) 424-2022 • FAX: (603) 429-8496
www.amrolabs.com

March 01, 2007

ANALYTICAL TEST RESULTS

Ed VanDoren
SHAW E & I, Inc.
11 Northeastern Boulevard
Salem, NH 030791953
TEL: (603) 870-4500
FAX: (603) 870-4501

Subject: 101960 Textron Gorham

Workorder No.: 0702072

Dear Ed VanDoren:

AMRO Environmental Laboratories Corp. received 24 samples on 2/19/07 for the analyses presented in the following report.

AMRO is accredited in accordance with NELAC and certifies that these test results meet all the requirements of NELAC, where applicable, unless otherwise noted in the case narrative.

The enclosed Sample Receipt Checklist details the condition of your sample(s) upon receipt. Please be advised that any unused sample volume and sample extracts will be stored for a period of 60 days from sample receipt date (90 days for samples from New York). After this time, AMRO will properly dispose of the remaining sample(s). If you require further analysis, or need the samples held for a longer period, please contact us immediately.

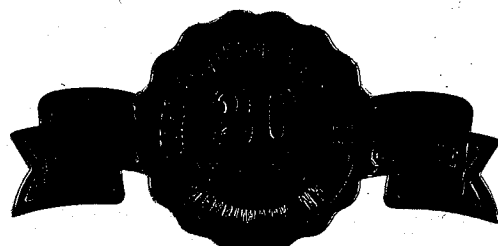
This report consists of a total of 90 pages. This letter is an integral part of your data report. All results in this project relate only to the sample(s) as received by the laboratory and documented in the Chain-of-Custody. This report shall not be reproduced except in full, without the written approval of the laboratory. If you have any questions regarding this project in the future, please refer to the Workorder Number above.

Sincerely,

Nancy Stewart
Vice President

State Certifications: NH (NELAC): 1001, MA: M-NH012, CT: PH-0758, NY: 11278 (NELAC), ME: NH012 and 1001, NJ: NH125, RI: 00105, U.S. Army Corps of Engineers (USACE), Naval Facilities Engineering Service Center (NFESC).

Hard copy of the State Certification is available upon request.



CLIENT: SHAW E & I, Inc.
Project: 101960 Textron Gorham
Lab Order: 0702072
Date Received: 2/19/07

Work Order Sample Summary

| Lab Sample ID | Client Sample ID | Collection Date | Collection Time |
|---------------|------------------|-----------------|-----------------|
| 0702072-01A | MW-207S | 2/17/07 | 7:00 AM |
| 0702072-02A | MW-207D | 2/17/07 | 7:30 AM |
| 0702072-03A | MW-202D | 2/17/07 | 8:00 AM |
| 0702072-04A | MW-202S | 2/17/07 | 8:30 AM |
| 0702072-05A | MW-101S | 2/17/07 | 9:00 AM |
| 0702072-06A | MW-101S DUP | 2/17/07 | 9:00 AM |
| 0702072-07A | MW-101D | 2/17/07 | 9:30 AM |
| 0702072-08A | MW-201D | 2/17/07 | 10:00 AM |
| 0702072-09A | MW-216S | 2/17/07 | 10:30 AM |
| 0702072-10A | MW-216D | 2/17/07 | 11:00 AM |
| 0702072-11A | MW-217D | 2/17/07 | 11:30 AM |
| 0702072-12A | MW-217S | 2/17/07 | 12:00 PM |
| 0702072-13A | MW-218S | 2/17/07 | 12:30 PM |
| 0702072-14A | MW-218D | 2/17/07 | 1:00 PM |
| 0702072-15A | MW-209D | 2/17/07 | 1:30 PM |
| 0702072-16A | MW-112 | 2/17/07 | 2:00 PM |
| 0702072-17A | MW-116S | 2/17/07 | 2:30 PM |
| 0702072-18A | MW-116D | 2/17/07 | 3:00 PM |
| 0702072-19A | CW-1 | 2/17/07 | 12:15 PM |
| 0702072-20A | CW-2 | 2/17/07 | 12:45 PM |
| 0702072-21A | GZA-5 | 2/17/07 | 2:30 PM |
| 0702072-21B | GZA-5 | 2/17/07 | 2:30 PM |
| 0702072-22A | GZA-5 DUP | 2/17/07 | 2:30 PM |
| 0702072-23A | CW-6 | 2/17/07 | 1:45 PM |
| 0702072-24A | CW-6 DUP | 2/17/07 | 1:45 PM |

AMRO Environmental Laboratories Corp.

01-Mar-07

DATES REPORT

Lab Order: 0702072

Client: SHAW E & I, Inc.

Project: 101960 Textron Gorham

| Sample ID | Client Sample ID | Collection Date | Matrix | Analytical Test Name | Preparatory Test Name | Prep Date | Analysis Date | Batch ID | TCLP Date |
|-------------|------------------|--------------------|-------------|------------------------------|-----------------------|-----------|---------------|----------|-----------|
| 0702072-01A | MW-207S | 2/17/07 7:00:00 AM | Groundwater | EPA 8260B VOLATILES by GC/MS | EPA 5030B | 2/17/07 | 2/21/07 | R35962 | |
| 0702072-02A | MW-207D | 2/17/07 7:30:00 AM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/21/07 | R35962 | |
| 0702072-03A | MW-202D | 2/17/07 8:00:00 AM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/23/07 | R35993 | |
| 0702072-04A | MW-202S | 2/17/07 8:30:00 AM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/21/07 | R35962 | |
| 0702072-05A | MW-101S | 2/17/07 9:00:00 AM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/22/07 | R35964 | |
| 0702072-06A | MW-101S DUP | | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/21/07 | R35962 | |
| 0702072-07A | MW-101D | 2/17/07 9:30:00 AM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/22/07 | R35964 | |
| | | | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/21/07 | R35962 | |

AMRO Environmental Laboratories Corp.

01-Mar-07

DATES REPORT

Lab Order: 0702072

Client: SHAW E & I, Inc.

Project: 101960 Textron Gorham

| Sample ID | Client Sample ID | Collection Date | Matrix | Analytical Test Name | Preparatory Test Name | Prep Date | Analysis Date | Batch ID | TCLP Date |
|-------------|------------------|---------------------|-------------|------------------------------|-----------------------|-----------|---------------|----------|-----------|
| 0702072-08A | MW-201D | 2/17/07 10:00:00 AM | Groundwater | EPA 8260B VOLATILES by GC/MS | EPA 5030B | 2/17/07 | 2/22/07 | R35964 | |
| 0702072-09A | MW-216S | 2/17/07 10:30:00 AM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/22/07 | R35964 | |
| 0702072-10A | MW-216D | 2/17/07 11:00:00 AM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/22/07 | R35964 | |
| 0702072-11A | MW-217D | 2/17/07 11:30:00 AM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/22/07 | R35964 | |
| 0702072-12A | MW-217S | 2/17/07 12:00:00 PM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/22/07 | R35964 | |
| 0702072-13A | MW-218S | 2/17/07 12:30:00 PM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/22/07 | R35964 | |
| 0702072-14A | MW-218D | 2/17/07 1:00:00 PM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/22/07 | R35964 | |
| 0702072-15A | MW-209D | 2/17/07 1:30:00 PM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/22/07 | R35964 | |
| 0702072-16A | MW-112 | 2/17/07 2:00:00 PM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/23/07 | R35993 | |
| 0702072-17A | MW-116S | 2/17/07 2:30:00 PM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/23/07 | R35993 | |
| 0702072-18A | MW-116D | 2/17/07 3:00:00 PM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/23/07 | R35993 | |
| 0702072-19A | CW-1 | 2/17/07 12:15:00 PM | | EPA 8260B VOLATILES by GC/MS | | 2/17/07 | 2/23/07 | R35993 | |

AMRO Environmental Laboratories Corp.

01-Mar-07

Lab Order: 0702072
Client: SHAW E & I, Inc.
Project: 101960 Textron Gorham

DATES REPORT

| Sample ID | Client Sample ID | Collection Date | Matrix | Analytical Test Name | Preparatory Test Name | Prep Date | Analysis Date | Batch ID | TCLP Date |
|-------------|------------------|---------------------|-------------|---|-----------------------|-----------|---------------|----------|-----------|
| 0702072-20A | CW-2 | 2/17/07 12:45:00 PM | Groundwater | EPA 8260B VOLATILES by GC/MS EPA 5030B | | 2/17/07 | 2/23/07 | R35993 | |
| 0702072-23A | CW-6 | 2/17/07 1:45:00 PM | | TPH by GC/FID (modified 8015B) AQPREP SEP FUNNEL: FING | | 2/23/07 | 2/23/07 | 16813 | |
| 0702072-24A | CW-6 DUP | | | TPH by GC/FID (modified 8015B) | | 2/23/07 | 2/23/07 | 16813 | |

AMRO Environmental Laboratories Corporation
 111 Herrick Street
 Merrimack, NH 03054

CHAIN-OF-CUSTODY RECORD

No 50751

Office: (603) 424-2022
 Fax: (603) 429-8496
 web: www.amrolabs.com

| | | | | |
|---|---------------------------------|-------------------------|---------------------------------|------------------------------|
| Project No.: 101960 | Project Name: Textron Gorham | Project State: RI | Project Manager: Ed Vandoren | AMRO Project No.: 0702072 |
| P.O.#: 157431 | Results Needed by: | Total # of Cont. & Size | Requested Analyses | Remarks |
| QUOTE #: | Seal Intact? Yes No N/A | Matrix | Requested Analyses | Remarks |
| Sample ID: | Date/Time Sampled | Matrix | Requested Analyses | Remarks |
| MW-2075 | 2/17/0700 | GW | | |
| MW-207D | 2/17/0730 | | | |
| MW-208D | 2/17/0800 | | | |
| MW-202S | 2/17/0830 | | | |
| MW-101S | 2/17/0900 | | | |
| MW-101S DUP | 2/17/0900 | | | |
| MW-101D | 2/17/0930 | | | |
| MW-201D | 2/17/1000 | | | |
| MW-216S | 2/17/1030 | | | |
| MW-216D | 2/17/1100 | | | |
| Preservative: Cl-HCl, MeOH, N-HNO3, S-H2SO4, Na-NaOH, O-Other | | | | |
| Priority Turnaround Time Authorization Before submitting samples for expedited TAT, you must have a coded AUTHORIZATION NUMBER | | | | |
| Send Results To: Edward Vandoren Shaw Environmental, Inc. 11 Northeastern Blvd. Salem, MA 03079-1953 | | | | |
| PHONE #: 603-870-4530 FAX #: 603-870-4501 E-mail: Edward.Vandoren@Shawgrp.com | | | | |
| Relinquished By: [Signature] Date/Time: 2/19/07 10:00 Received By: [Signature] Date/Time: 2/19/07 13:00 | | | | |
| Please print clearly, legibly and completely. Samples can not be logged in and the turnaround time clock will not start until any ambiguities are resolved. | | | | |
| White: Lab Copy Yellow: Accompanies Report Pink: Client Copy SHEET OF SHEET AMROCOC2004, Rev.3 08/18/04 | | | | |

Requested Analyses

Requested Analyses

Requested Analyses

Requested Analyses

Requested Analyses

Requested Analyses

Requested Analyses

Requested Analyses

Requested Analyses

Requested Analyses

AMRO Environmental Laboratories Corporation
 111 Herrick Street
 Merrimack, NH 03054

CHAIN-OF-CUSTODY RECORD

No. 50750

Office: (603) 424-2022
 Fax: (603) 429-8496
 web: www.amrolabs.com

| | | | | |
|--|------------------------------|-------------------------|------------------------------|---|
| Project No.: 101960 | Project Name: Textron Gorham | Project State: RI | Project Manager: Ed Vandoren | AMRO Project No.: 0702012 |
| P.O.#: 157431 | Results Needed by: | Project State: RI | Requested Analyses: | Remarks: |
| QUOTE #: | Seal Intact? Yes No N/A | Comp. Grab | Requested Analyses: | Email GISKey formatted EDD to: Catherine.Joe@shawgrp.com and PDF of report. |
| Sample ID: | Date/Time Sampled | Total # of Cont. & Size | Requested Analyses: | |
| MW-217D | 2/17/17 1130 | 2 | | |
| MW-217S | 2/17/17 1200 | 2 | | |
| MW-218S | 2/17/17 1230 | 2 | | |
| MW-218D | 2/17/17 1300 | 2 | | |
| MW-209D | 2/17/17 1330 | 2 | | |
| MW-112 | 2/17/17 1400 | 2 | | |
| MW-116S | 2/17/17 1430 | 2 | | |
| MW-116D | 2/17/17 1500 | 2 | | |
| TRIP BANK | Received by LAB | 1 | | |
| Preservative: Cl-HCl, MeOH, N-HNO3, S-H2SO4, Na-NaOH, O-Other | | | | |
| Priority Turnaround Time Authorization | | | | |
| Send Results To: Ed Vandoren | | | | |
| Shaw Environmental, Inc. | | | | |
| 11 Northeastern Blvd. | | | | |
| Salem, MA 03079-1953 | | | | |
| PHONE #: 603-870-4530 FAX #: 603-870-4501 | | | | |
| E-mail: edward.vandoren@shawgrp.com | | | | |
| Received By: [Signature] | | | | |
| Date/Time: 2/17/07 12:00 | | | | |
| 2/19/07 10:00 | | | | |
| 2/19/07 1300 | | | | |
| Samples arriving after 12:00 noon will be tracked and billed as received on the following day. | | | | |
| White: Lab Copy Yellow: Accompanies Report Pink: Client Copy SHEET OF | | | | |
| AMROCOC2004, Rev.3 08/18/04 | | | | |

Login Account for multiple users

From: VanDoren, Edward [Edward.VanDoren@shawgrp.com]
Sent: Thursday, February 22, 2007 2:13 PM
To: Login Account for multiple users
Subject: RE: Testron Gorham Sampling Date (AMRO 0702072)

Hi Connie-

Please do not analyze GZA-5 or GZA-5 Dup. The wrong well was sampled.

Thanks,
Ed

Edward Van Doren
Client Program Manager
Shaw Environmental, Inc.
11 Northeastern Boulevard
Salem, NH 03079-1953
603.870.4530 direct
603.870.4501 fax
edward.vandoren@shawgrp.com
www.shawgrp.com

From: Login Account for multiple users [mailto:login@amrolabs.com]
Sent: Tuesday, February 20, 2007 11:21 AM
To: VanDoren, Edward
Subject: FW: Testron Gorham Sampling Date (AMRO 0702072)

Hi Ed -

We had a call from Dan Leahy and he let us know that the samples were taken 2/17.

Connie in Receiving

From: Login Account for multiple users
Sent: Monday, February 19, 2007 3:11 PM
To: 'VanDoren, Edward'
Subject: Testron Gorham Sampling Date (AMRO 0702072)

Hi Ed -

We picked up these water samples today and I notice that all of the samples are dated **2/14/07** and the Chain reads **2/17/07**. Can you tell me which is the correct sampling date?

Also, the Trip Blank was found broken in the cooler, still inside the bubble bag, so we will be unable to run it for you.

Thanks for your help!

Connie in Receiving

| | |
|--|---------------------------|
| Client: <u>SHAW ENVIRONMENTAL, INC</u> | AMRO ID: <u>0702072</u> |
| Project Name: <u>TEXTRON GORHAM</u> | Date Rec.: <u>2-19-07</u> |
| Ship via: (circle one) Fed Ex., UPS, <u>(AMRO Courier)</u> | Date Due: <u>2-26-07</u> |
| Hand Del., Other Courier, Other: | |

Items to be Checked Upon Receipt

1. Army Samples received in individual plastic bags?
2. Custody Seals present?
3. Custody Seals Intact?
4. Air Bill included in folder if received?
5. Is COC included with samples?
6. Is COC signed and dated by client?
7. Laboratory receipt temperature. TEMP = 3°
 Samples rec. with ice ice packs neither
8. Were samples received the same day they were sampled?
 Is client temperature 4°C ± 2°C?
If no obtain authorization from the client for the analyses.
 Client authorization from: _____ Date: _____ Obtained by: _____
9. Is the COC filled out correctly and completely?
10. Does the info on the COC match the samples?
11. Were samples rec. within holding time?
12. Were all samples properly labeled?
13. Were all samples properly preserved?
14. Were proper sample containers used?
15. Were all samples received intact? (none broken or leaking)
16. Were VOA vials rec. with no air bubbles?
17. Were the sample volumes sufficient for requested analysis?
18. Were all samples received?

| Yes | No | NA | Comments |
|-----|----|----|--|
| | | ✓ | |
| | | ✓ | |
| | | ✓ | |
| | | ✓ | |
| ✓ | | | |
| ✓ | | | |
| | ✓ | | |
| ✓ | | | |
| | | | |
| ✓ | ✓ | | ALL DATED 2/14/07 EMAIL TO CLIENT |
| ✓ | | | |
| ✓ | | | |
| ✓ | | | |
| ✓ | ✓ | | TRIP BLANK RECEIVED BROKEN - EMAIL TO CLIENT |
| ✓ | | | |
| ✓ | | | |

19. VPH and VOA Soils only:
 Sampling Method VPH (circle one): M=Methanol, E=EnCore (air-tight container)
 Sampling Method VOA (circle one): M=Methanol, SB=Sodium Bisulfate, E=EnCore, B=Bulk
 If M or SB:
 Does preservative cover the soil?
If NO then client must be faxed.
 Does preservation level come close to the fill line on the vial?
If NO then client must be faxed.
 Were vials provided by AMRO?
If NO then weights MUST be obtained from client
 Was dry weight aliquot provided?
If NO then fax client and inform the VOA lab ASAP.

| Yes | No | NA | Comments |
|-----|----|----|----------|
| | | ✓ | |
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| | | | |

20. Subcontracted Samples:
 What samples sent:
 Where sent:
 Date:
 Analysis:
 TAT:

| Yes | No | NA | Comments |
|-----|----|----|----------|
| | | ✓ | |
| | | | |
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| | | | |
| | | | |

21. Information entered into:
 Internal Tracking Log?
 Dry Weight Log?
 Client Log?
 Composite Log?
 Filtration Log?

| Yes | No | NA | Comments |
|-----|----|----|----------|
| ✓ | | | |
| | | ✓ | |
| | | ✓ | |
| | | ✓ | |
| | | ✓ | |

| | | | |
|------------------------|----------------------|----------------------------|----------------------|
| Received By: <u>CC</u> | Date: <u>2-19-07</u> | Logged in By: <u>CC MG</u> | Date: <u>2-19-07</u> |
| Labeled By: <u>CC</u> | Date: <u>2-19-07</u> | Checked By: _____ | Date: <u>2-21-07</u> |

CLIENT: SHAW E & I, Inc.
Project: 101960 Textron Gorham
Lab Order: 0702072

CASE NARRATIVE

GC/MS VOLATILES:

1. A Matrix Spike (MS) and Matrix Spike Duplicate (MSD) were performed on sample MW-209D (0702072-15A) Batch ID: R35964.

1.1 The % Recovery for 4 analytes out of 67 analytes in the MS was outside the laboratory control limits.

2. A Matrix Spike (MS) and Matrix Spike Duplicate (MSD) were performed on sample CW-1 (0702072-19A) Batch ID: R35993.

2.1 The % Recovery for 2 analytes out of 67 analytes in the MS was outside the laboratory control limits.

2.2 The RPD for 1 analyte out of 67 analytes was outside the laboratory control limits.

DATA COMMENT PAGE

Organic Data Qualifiers

| | |
|----|--|
| ND | Indicates compound was analyzed for, but not detected at or above the reporting limit. |
| J | Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed, or when the data indicates the presence of a compound that meets the identification criteria but the result is less than the sample quantitation limit but greater than the method detection limit. |
| H | Method prescribed holding time exceeded. |
| E | This flag identifies compounds whose concentrations exceed the calibration range of the instrument for that specific analysis. |
| B | This flag is used when the analyte is found in the associated blank as well as in the sample. |
| R | RPD outside accepted recovery limits |
| RL | Reporting limit; defined as the lowest concentration the laboratory can accurately quantitate. |
| S | Spike Recovery outside accepted recovery limits. |
| # | See Case Narrative |

Micro Data Qualifiers

TNTC Too numerous to count

Inorganic Data Qualifiers

| | |
|---------|---|
| ND or U | Indicates element was analyzed for, but not detected at or above the reporting limit. |
| J | Indicates a value greater than or equal to the method detection limit, but less than the quantitation limit. |
| H | Indicates analytical holding time exceedance. |
| B | Indicates that the analyte is found in the associated blank, as well as in the sample. |
| MSA | Indicates value determined by the Method of Standard Addition |
| E | This flag identifies compounds whose concentrations exceed the calibration range of the instrument for that specific analysis. |
| R | RPD outside accepted recovery limits |
| RL | Reporting limit; defined as the lowest concentration the laboratory can accurately quantitate. |
| S | Spike Recovery outside accepted recovery limits. |
| W | Post-digestion spike for Furnace AA analysis is out of control limits (85-115), while sample absorbance is less than 50% of spike absorbance. |
| * | Duplicate analysis not within control limits. |
| + | Indicates the correlation coefficient for the Method of Standard Addition is less than 0.995 |
| # | See Case Narrative |

Report Comments:

1. Soil, sediment and sludge sample results are reported on a "dry weight" basis.
2. Reporting limits are adjusted for sample size used, dilutions and moisture content, if applicable.

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
 Lab Order: 0702072
 Project: 101960 Textron Gorham
 Lab ID: 0702072-01A

Client Sample ID: MW-207S
 Collection Date: 2/17/07 7:00:00 AM
 Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Vinyl chloride | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,1-Dichloroethene | ND | 100 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| cis-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Trichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Tetrachloroethene | 12,000 | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-01A

Client Sample ID: MW-207S
Collection Date: 2/17/07 7:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|--------------------|
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 8:31:00 PM |
| Surr: Dibromofluoromethane | 100 | 85-116 | | %REC | 100 | 2/21/07 8:31:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 108 | 77-127 | | %REC | 100 | 2/21/07 8:31:00 PM |
| Surr: Toluene-d8 | 101 | 86-114 | | %REC | 100 | 2/21/07 8:31:00 PM |
| Surr: 4-Bromofluorobenzene | 95.9 | 79-117 | | %REC | 100 | 2/21/07 8:31:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-02A

Client Sample ID: MW-207D
Collection Date: 2/17/07 7:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Vinyl chloride | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,1-Dichloroethene | ND | 100 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| cis-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Trichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Tetrachloroethene | 7,700 | 200 | | µg/L | 100 | 2/23/07 2:53:00 PM |
| Tetrachloroethene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
 Lab Order: 0702072
 Project: 101960 Textron Gorham
 Lab ID: 0702072-02A

Client Sample ID: MW-207D
 Collection Date: 2/17/07 7:30:00 AM
 Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|--------------------|
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:05:00 PM |
| Surr: Dibromofluoromethane | 97.3 | 85-116 | | %REC | 100 | 2/21/07 9:05:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 94.9 | 77-127 | | %REC | 100 | 2/21/07 9:05:00 PM |
| Surr: Toluene-d8 | 99.0 | 86-114 | | %REC | 100 | 2/21/07 9:05:00 PM |
| Surr: 4-Bromofluorobenzene | 91.2 | 79-117 | | %REC | 100 | 2/21/07 9:05:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-03A

Client Sample ID: MW-202D
Collection Date: 2/17/07 8:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Vinyl chloride | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,1-Dichloroethene | ND | 100 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| cis-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Trichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Tetrachloroethene | 29,000 | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-03A

Client Sample ID: MW-202D
Collection Date: 2/17/07 8:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|--------------------|
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 9:40:00 PM |
| Surr: Dibromofluoromethane | 102 | 85-116 | | %REC | 100 | 2/21/07 9:40:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 107 | 77-127 | | %REC | 100 | 2/21/07 9:40:00 PM |
| Surr: Toluene-d8 | 101 | 86-114 | | %REC | 100 | 2/21/07 9:40:00 PM |
| Surr: 4-Bromofluorobenzene | 95.9 | 79-117 | | %REC | 100 | 2/21/07 9:40:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-04A

Client Sample ID: MW-202S
Collection Date: 2/17/07 8:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|---------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Vinyl chloride | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,1-Dichloroethene | ND | 100 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| cis-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Trichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Tetrachloroethene | 51,000 | 1,000 | | µg/L | 500 | 2/22/07 3:12:00 PM |
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-04A

Client Sample ID: MW-202S
Collection Date: 2/17/07 8:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|---------------------|
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:14:00 PM |
| Surr: Dibromofluoromethane | 104 | 85-116 | | %REC | 100 | 2/21/07 10:14:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 107 | 77-127 | | %REC | 100 | 2/21/07 10:14:00 PM |
| Surr: Toluene-d8 | 100 | 86-114 | | %REC | 100 | 2/21/07 10:14:00 PM |
| Surr: 4-Bromofluorobenzene | 98.8 | 79-117 | | %REC | 100 | 2/21/07 10:14:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-05A

Client Sample ID: MW-101S
Collection Date: 2/17/07 9:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|---------------------|
| EPA 8260B VOLATILES BY GC/MS | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Vinyl chloride | 520 | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,1-Dichloroethene | ND | 100 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| cis-1,2-Dichloroethene | 510 | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Trichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Tetrachloroethene | 45,000 | 1,000 | | µg/L | 500 | 2/22/07 3:46:00 PM |
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.

Client Sample ID: MW-101S

Lab Order: 0702072

Collection Date: 2/17/07 9:00:00 AM

Project: 101960 Textron Gorham

Matrix: GROUNDWATER

Lab ID: 0702072-05A

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|---------------------|
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 10:49:00 PM |
| Surr: Dibromofluoromethane | 104 | 85-116 | | %REC | 100 | 2/21/07 10:49:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 109 | 77-127 | | %REC | 100 | 2/21/07 10:49:00 PM |
| Surr: Toluene-d8 | 102 | 86-114 | | %REC | 100 | 2/21/07 10:49:00 PM |
| Surr: 4-Bromofluorobenzene | 96.6 | 79-117 | | %REC | 100 | 2/21/07 10:49:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-06A

Client Sample ID: MW-101S DUP
Collection Date: 2/17/07 9:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|--------------------|---------------------|
| EPA 8260B VOLATILES BY GC/MS | | SW8260B | | | Analyst: SK | |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Vinyl chloride | 460 | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,1-Dichloroethene | ND | 100 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| cis-1,2-Dichloroethene | 500 | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Trichloroethene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Tetrachloroethene | 44,000 | 1,000 | | µg/L | 500 | 2/22/07 4:20:00 PM |
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-06A

Client Sample ID: MW-101S DUP
Collection Date: 2/17/07 9:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|---------------------|
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/21/07 11:23:00 PM |
| Surr: Dibromofluoromethane | 101 | 85-116 | | %REC | 100 | 2/21/07 11:23:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 108 | 77-127 | | %REC | 100 | 2/21/07 11:23:00 PM |
| Surr: Toluene-d8 | 102 | 86-114 | | %REC | 100 | 2/21/07 11:23:00 PM |
| Surr: 4-Bromofluorobenzene | 99.6 | 79-117 | | %REC | 100 | 2/21/07 11:23:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-07A

Client Sample ID: MW-101D
Collection Date: 2/17/07 9:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|---------------------|
| EPA 8260B VOLATILES BY GC/MS | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 50 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Chloromethane | ND | 50 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Vinyl chloride | 30 | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Chloroethane | ND | 50 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Bromomethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Trichlorofluoromethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Diethyl ether | ND | 50 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Acetone | ND | 100 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,1-Dichloroethene | 11 | 10 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Carbon disulfide | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Methylene chloride | ND | 50 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Methyl tert-butyl ether | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| trans-1,2-Dichloroethene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,1-Dichloroethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 2-Butanone | ND | 100 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 2,2-Dichloropropane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| cis-1,2-Dichloroethene | 280 | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Chloroform | 20 | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Tetrahydrofuran | ND | 100 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Bromochloromethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,1,1-Trichloroethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,1-Dichloropropene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Carbon tetrachloride | 32 | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2-Dichloroethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Benzene | ND | 10 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Trichloroethene | 160 | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2-Dichloropropane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Bromodichloromethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Dibromomethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 4-Methyl-2-pentanone | ND | 100 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| cis-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Toluene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| trans-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,1,2-Trichloroethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2-Dibromoethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 2-Hexanone | ND | 100 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,3-Dichloropropane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Tetrachloroethene | 19,000 | 200 | | µg/L | 100 | 2/22/07 4:54:00 PM |
| Dibromochloromethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-07A

Client Sample ID: MW-101D
Collection Date: 2/17/07 9:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|---------------------|
| Chlorobenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Ethylbenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| m,p-Xylene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| o-Xylene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Styrene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Bromoform | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Isopropylbenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2,3-Trichloropropane | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Bromobenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| n-Propylbenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 2-Chlorotoluene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 4-Chlorotoluene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,3,5-Trimethylbenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| tert-Butylbenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2,4-Trimethylbenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| sec-Butylbenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 4-Isopropyltoluene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,3-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,4-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| n-Butylbenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 50 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2,4-Trichlorobenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Hexachlorobutadiene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Naphthalene | ND | 50 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| 1,2,3-Trichlorobenzene | ND | 20 | | µg/L | 10 | 2/21/07 11:58:00 PM |
| Surr: Dibromofluoromethane | 102 | 85-116 | | %REC | 10 | 2/21/07 11:58:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 107 | 77-127 | | %REC | 10 | 2/21/07 11:58:00 PM |
| Surr: Toluene-d8 | 101 | 86-114 | | %REC | 10 | 2/21/07 11:58:00 PM |
| Surr: 4-Bromofluorobenzene | 95.3 | 79-117 | | %REC | 10 | 2/21/07 11:58:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-08A

Client Sample ID: MW-201D
Collection Date: 2/17/07 10:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Vinyl chloride | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,1-Dichloroethene | ND | 100 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| cis-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Trichloroethene | 970 | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Tetrachloroethene | 7,600 | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-08A

Client Sample ID: MW-201D
Collection Date: 2/17/07 10:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|--------------------|
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/22/07 7:47:00 PM |
| Surr: Dibromofluoromethane | 99.6 | 85-116 | | %REC | 100 | 2/22/07 7:47:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 101 | 77-127 | | %REC | 100 | 2/22/07 7:47:00 PM |
| Surr: Toluene-d8 | 98.0 | 86-114 | | %REC | 100 | 2/22/07 7:47:00 PM |
| Surr: 4-Bromofluorobenzene | 97.4 | 79-117 | | %REC | 100 | 2/22/07 7:47:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-09A

Client Sample ID: MW-216S
Collection Date: 2/17/07 10:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Trichlorofluoromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| cis-1,2-Dichloroethene | 140 | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Chloroform | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Trichloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Bromodichloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Toluene | 3.4 | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Tetrachloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-09A

Client Sample ID: MW-216S
Collection Date: 2/17/07 10:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Ethylbenzene | 2.6 | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| m,p-Xylene | 7.0 | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| o-Xylene | 8.9 | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,3,5-Trimethylbenzene | 10 | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2,4-Trimethylbenzene | 13 | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 4-Isopropyltoluene | 2.9 | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Naphthalene | 17 | 5.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 5:28:00 PM |
| Surr: Dibromofluoromethane | 102 | 85-116 | | %REC | 1 | 2/22/07 5:28:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 109 | 77-127 | | %REC | 1 | 2/22/07 5:28:00 PM |
| Surr: Toluene-d8 | 102 | 86-114 | | %REC | 1 | 2/22/07 5:28:00 PM |
| Surr: 4-Bromofluorobenzene | 99.4 | 79-117 | | %REC | 1 | 2/22/07 5:28:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-10A

Client Sample ID: MW-216D
Collection Date: 2/17/07 11:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Trichlorofluoromethane | 3.1 | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| cis-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Chloroform | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Trichloroethene | 6.4 | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Bromodichloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Toluene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Tetrachloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-10A

Client Sample ID: MW-216D
Collection Date: 2/17/07 11:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Ethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| m,p-Xylene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| o-Xylene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,3,5-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2,4-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Naphthalene | ND | 5.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:03:00 PM |
| Surr: Dibromofluoromethane | 98.6 | 85-116 | | %REC | 1 | 2/22/07 6:03:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 107 | 77-127 | | %REC | 1 | 2/22/07 6:03:00 PM |
| Surr: Toluene-d8 | 99.4 | 86-114 | | %REC | 1 | 2/22/07 6:03:00 PM |
| Surr: 4-Bromofluorobenzene | 100 | 79-117 | | %REC | 1 | 2/22/07 6:03:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-11A

Client Sample ID: MW-217D
Collection Date: 2/17/07 11:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | SW8260B | | Analyst: SK | | |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Trichlorofluoromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| cis-1,2-Dichloroethene | 60 | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Chloroform | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Trichloroethene | 75 | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Bromodichloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Toluene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Tetrachloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-11A

Client Sample ID: MW-217D
Collection Date: 2/17/07 11:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Ethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| m,p-Xylene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| o-Xylene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,3,5-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2,4-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Naphthalene | ND | 5.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 6:38:00 PM |
| Surr: Dibromofluoromethane | 100 | 85-116 | | %REC | 1 | 2/22/07 6:38:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 107 | 77-127 | | %REC | 1 | 2/22/07 6:38:00 PM |
| Surr: Toluene-d8 | 100 | 86-114 | | %REC | 1 | 2/22/07 6:38:00 PM |
| Surr: 4-Bromofluorobenzene | 101 | 79-117 | | %REC | 1 | 2/22/07 6:38:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-12A

Client Sample ID: MW-217S
Collection Date: 2/17/07 12:00:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Trichlorofluoromethane | 2.2 | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| cis-1,2-Dichloroethene | 9.8 | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Chloroform | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Trichloroethene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Bromodichloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Toluene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Tetrachloroethene | 23 | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-12A

Client Sample ID: MW-217S
Collection Date: 2/17/07 12:00:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Ethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| m,p-Xylene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| o-Xylene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,3,5-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2,4-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Naphthalene | ND | 5.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/22/07 7:12:00 PM |
| Surr: Dibromofluoromethane | 98.9 | 85-116 | | %REC | 1 | 2/22/07 7:12:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 103 | 77-127 | | %REC | 1 | 2/22/07 7:12:00 PM |
| Surr: Toluene-d8 | 100 | 86-114 | | %REC | 1 | 2/22/07 7:12:00 PM |
| Surr: 4-Bromofluorobenzene | 95.0 | 79-117 | | %REC | 1 | 2/22/07 7:12:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-13A

Client Sample ID: MW-218S
Collection Date: 2/17/07 12:30:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 50 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Chloromethane | ND | 50 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Vinyl chloride | 50 | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Chloroethane | ND | 50 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Bromomethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Trichlorofluoromethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Diethyl ether | ND | 50 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Acetone | ND | 100 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,1-Dichloroethene | ND | 10 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Carbon disulfide | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Methylene chloride | ND | 50 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Methyl tert-butyl ether | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| trans-1,2-Dichloroethene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,1-Dichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 2-Butanone | ND | 100 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 2,2-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| cis-1,2-Dichloroethene | 650 | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Chloroform | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Tetrahydrofuran | ND | 100 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Bromochloromethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,1,1-Trichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,1-Dichloropropene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Carbon tetrachloride | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2-Dichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Benzene | ND | 10 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Trichloroethene | 49 | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Bromodichloromethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Dibromomethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 4-Methyl-2-pentanone | ND | 100 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| cis-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Toluene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| trans-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,1,2-Trichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2-Dibromoethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 2-Hexanone | ND | 100 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,3-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Tetrachloroethene | 370 | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Dibromochloromethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-13A

Client Sample ID: MW-218S
Collection Date: 2/17/07 12:30:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Ethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| m,p-Xylene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| o-Xylene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Styrene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Bromoform | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Isopropylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2,3-Trichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Bromobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| n-Propylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 2-Chlorotoluene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 4-Chlorotoluene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,3,5-Trimethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| tert-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2,4-Trimethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| sec-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 4-Isopropyltoluene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,3-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,4-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| n-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 50 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2,4-Trichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Hexachlorobutadiene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Naphthalene | ND | 50 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| 1,2,3-Trichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:21:00 PM |
| Surr: Dibromofluoromethane | 103 | 85-116 | | %REC | 10 | 2/22/07 8:21:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 108 | 77-127 | | %REC | 10 | 2/22/07 8:21:00 PM |
| Surr: Toluene-d8 | 103 | 86-114 | | %REC | 10 | 2/22/07 8:21:00 PM |
| Surr: 4-Bromofluorobenzene | 97.1 | 79-117 | | %REC | 10 | 2/22/07 8:21:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc. **Client Sample ID:** MW-218D
Lab Order: 0702072 **Collection Date:** 2/17/07 1:00:00 PM
Project: 101960 Textron Gorham **Matrix:** GROUNDWATER
Lab ID: 0702072-14A

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 50 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Chloromethane | ND | 50 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Vinyl chloride | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Chloroethane | ND | 50 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Bromomethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Trichlorofluoromethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Diethyl ether | ND | 50 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Acetone | ND | 100 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,1-Dichloroethene | 23 | 10 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Carbon disulfide | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Methylene chloride | ND | 50 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Methyl tert-butyl ether | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| trans-1,2-Dichloroethene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,1-Dichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 2-Butanone | ND | 100 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 2,2-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| cis-1,2-Dichloroethene | 28 | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Chloroform | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Tetrahydrofuran | ND | 100 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Bromochloromethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,1,1-Trichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,1-Dichloropropene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Carbon tetrachloride | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2-Dichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Benzene | ND | 10 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Trichloroethene | 840 | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Bromodichloromethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Dibromomethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 4-Methyl-2-pentanone | ND | 100 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| cis-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Toluene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| trans-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,1,2-Trichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2-Dibromoethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 2-Hexanone | ND | 100 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,3-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Tetrachloroethene | 600 | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Dibromochloromethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-14A

Client Sample ID: MW-218D
Collection Date: 2/17/07 1:00:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Ethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| m,p-Xylene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| o-Xylene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Styrene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Bromoform | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Isopropylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2,3-Trichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Bromobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| n-Propylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 2-Chlorotoluene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 4-Chlorotoluene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,3,5-Trimethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| tert-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2,4-Trimethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| sec-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 4-Isopropyltoluene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,3-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,4-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| n-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 50 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2,4-Trichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Hexachlorobutadiene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Naphthalene | ND | 50 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| 1,2,3-Trichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 8:56:00 PM |
| Surr: Dibromofluoromethane | 101 | 85-116 | | %REC | 10 | 2/22/07 8:56:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 106 | 77-127 | | %REC | 10 | 2/22/07 8:56:00 PM |
| Surr: Toluene-d8 | 102 | 86-114 | | %REC | 10 | 2/22/07 8:56:00 PM |
| Surr: 4-Bromofluorobenzene | 97.8 | 79-117 | | %REC | 10 | 2/22/07 8:56:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-15A

Client Sample ID: MW-209D
Collection Date: 2/17/07 1:30:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 50 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Chloromethane | ND | 50 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Vinyl chloride | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Chloroethane | ND | 50 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Bromomethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Trichlorofluoromethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Diethyl ether | ND | 50 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Acetone | ND | 100 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,1-Dichloroethene | ND | 10 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Carbon disulfide | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Methylene chloride | ND | 50 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Methyl tert-butyl ether | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| trans-1,2-Dichloroethene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,1-Dichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 2-Butanone | ND | 100 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 2,2-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| cis-1,2-Dichloroethene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Chloroform | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Tetrahydrofuran | ND | 100 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Bromochloromethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,1,1-Trichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,1-Dichloropropene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Carbon tetrachloride | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2-Dichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Benzene | ND | 10 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Trichloroethene | 110 | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Bromodichloromethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Dibromomethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 4-Methyl-2-pentanone | ND | 100 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| cis-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Toluene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| trans-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,1,2-Trichloroethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2-Dibromoethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 2-Hexanone | ND | 100 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,3-Dichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Tetrachloroethene | 430 | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Dibromochloromethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-15A

Client Sample ID: MW-209D
Collection Date: 2/17/07 1:30:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Ethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| m,p-Xylene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| o-Xylene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Styrene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Bromoform | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Isopropylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2,3-Trichloropropane | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Bromobenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| n-Propylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 2-Chlorotoluene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 4-Chlorotoluene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,3,5-Trimethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| tert-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2,4-Trimethylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| sec-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 4-Isopropyltoluene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,3-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,4-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| n-Butylbenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2-Dichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 50 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2,4-Trichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Hexachlorobutadiene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Naphthalene | ND | 50 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| 1,2,3-Trichlorobenzene | ND | 20 | | µg/L | 10 | 2/22/07 9:30:00 PM |
| Surr: Dibromofluoromethane | 98.9 | 85-116 | | %REC | 10 | 2/22/07 9:30:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 103 | 77-127 | | %REC | 10 | 2/22/07 9:30:00 PM |
| Surr: Toluene-d8 | 98.4 | 86-114 | | %REC | 10 | 2/22/07 9:30:00 PM |
| Surr: 4-Bromofluorobenzene | 96.8 | 79-117 | | %REC | 10 | 2/22/07 9:30:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-16A

Client Sample ID: MW-112
Collection Date: 2/17/07 2:00:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Trichlorofluoromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Methyl tert-butyl ether | 17 | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| cis-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Chloroform | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Trichloroethene | 3.9 | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Bromodichloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Toluene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Tetrachloroethene | 42 | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-16A

Client Sample ID: MW-112
Collection Date: 2/17/07 2:00:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Ethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| m,p-Xylene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| o-Xylene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,3,5-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2,4-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Naphthalene | ND | 5.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 3:28:00 PM |
| Surr: Dibromofluoromethane | 97.7 | 85-116 | | %REC | 1 | 2/23/07 3:28:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 103 | 77-127 | | %REC | 1 | 2/23/07 3:28:00 PM |
| Surr: Toluene-d8 | 99.0 | 86-114 | | %REC | 1 | 2/23/07 3:28:00 PM |
| Surr: 4-Bromofluorobenzene | 97.7 | 79-117 | | %REC | 1 | 2/23/07 3:28:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-17A

Client Sample ID: MW-116S
Collection Date: 2/17/07 2:30:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Trichlorofluoromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| cis-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Chloroform | 32 | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Trichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Bromodichloromethane | 4.2 | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Toluene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Tetrachloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-17A

Client Sample ID: MW-116S
Collection Date: 2/17/07 2:30:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Ethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| m,p-Xylene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| o-Xylene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,3,5-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2,4-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Naphthalene | ND | 5.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:03:00 PM |
| Surr: Dibromofluoromethane | 102 | 85-116 | | %REC | 1 | 2/23/07 4:03:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 104 | 77-127 | | %REC | 1 | 2/23/07 4:03:00 PM |
| Surr: Toluene-d8 | 99.4 | 86-114 | | %REC | 1 | 2/23/07 4:03:00 PM |
| Surr: 4-Bromofluorobenzene | 96.5 | 79-117 | | %REC | 1 | 2/23/07 4:03:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-18A

Client Sample ID: MW-116D
Collection Date: 2/17/07 3:00:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Trichlorofluoromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| cis-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Chloroform | 39 | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Trichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Bromodichloromethane | 4.3 | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Toluene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Tetrachloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-18A

Client Sample ID: MW-116D
Collection Date: 2/17/07 3:00:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Ethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| m,p-Xylene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| o-Xylene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,3,5-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2,4-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Naphthalene | ND | 5.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 4:37:00 PM |
| Surr: Dibromofluoromethane | 98.3 | 85-116 | | %REC | 1 | 2/23/07 4:37:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 104 | 77-127 | | %REC | 1 | 2/23/07 4:37:00 PM |
| Surr: Toluene-d8 | 99.8 | 86-114 | | %REC | 1 | 2/23/07 4:37:00 PM |
| Surr: 4-Bromofluorobenzene | 96.8 | 79-117 | | %REC | 1 | 2/23/07 4:37:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-19A

Client Sample ID: CW-1
Collection Date: 2/17/07 12:15:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|-----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 500 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Chloromethane | ND | 500 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Vinyl chloride | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Chloroethane | ND | 500 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Bromomethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Trichlorofluoromethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Diethyl ether | ND | 500 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Acetone | ND | 1,000 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,1-Dichloroethene | 180 | 100 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Carbon disulfide | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Methylene chloride | ND | 500 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Methyl tert-butyl ether | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| trans-1,2-Dichloroethene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,1-Dichloroethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 2-Butanone | ND | 1,000 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 2,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| cis-1,2-Dichloroethene | 350 | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Chloroform | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Tetrahydrofuran | ND | 1,000 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Bromochloromethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,1,1-Trichloroethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,1-Dichloropropene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Carbon tetrachloride | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2-Dichloroethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Benzene | ND | 100 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Trichloroethene | 5,900 | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2-Dichloropropane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Bromodichloromethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Dibromomethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 4-Methyl-2-pentanone | ND | 1,000 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| cis-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Toluene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| trans-1,3-Dichloropropene | ND | 100 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,1,2-Trichloroethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2-Dibromoethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 2-Hexanone | ND | 1,000 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,3-Dichloropropane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Tetrachloroethene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Dibromochloromethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-19A

Client Sample ID: CW-1
Collection Date: 2/17/07 12:15:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|-----|--------------------|
| Chlorobenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Ethylbenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| m,p-Xylene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| o-Xylene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Styrene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Bromoform | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Isopropylbenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2,3-Trichloropropane | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Bromobenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| n-Propylbenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 2-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 4-Chlorotoluene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,3,5-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| tert-Butylbenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2,4-Trimethylbenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| sec-Butylbenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 4-Isopropyltoluene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,3-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,4-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| n-Butylbenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2-Dichlorobenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 500 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2,4-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Hexachlorobutadiene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Naphthalene | ND | 500 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| 1,2,3-Trichlorobenzene | ND | 200 | | µg/L | 100 | 2/23/07 5:46:00 PM |
| Surr: Dibromofluoromethane | 100 | 85-116 | | %REC | 100 | 2/23/07 5:46:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 106 | 77-127 | | %REC | 100 | 2/23/07 5:46:00 PM |
| Surr: Toluene-d8 | 101 | 86-114 | | %REC | 100 | 2/23/07 5:46:00 PM |
| Surr: 4-Bromofluorobenzene | 97.2 | 79-117 | | %REC | 100 | 2/23/07 5:46:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-20A

Client Sample ID: CW-2
Collection Date: 2/17/07 12:45:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Trichlorofluoromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| cis-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Chloroform | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Trichloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Bromodichloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Toluene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Tetrachloroethene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-20A

Client Sample ID: CW-2
Collection Date: 2/17/07 12:45:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|--------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Ethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| m,p-Xylene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| o-Xylene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,3,5-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2,4-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Naphthalene | ND | 5.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 2/23/07 5:11:00 PM |
| Surr: Dibromofluoromethane | 101 | 85-116 | | %REC | 1 | 2/23/07 5:11:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 106 | 77-127 | | %REC | 1 | 2/23/07 5:11:00 PM |
| Surr: Toluene-d8 | 101 | 86-114 | | %REC | 1 | 2/23/07 5:11:00 PM |
| Surr: 4-Bromofluorobenzene | 97.0 | 79-117 | | %REC | 1 | 2/23/07 5:11:00 PM |

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT

Method Blank

CLIENT: SHAW E & I, Inc.
 Work Order: 0702072
 Project: 101960 Textron Gorham

Sample ID: mb-02/21/07 Batch ID: R35962 Test Code: SW8260B Units: µg/L Analysis Date 2/21/2007 3:55:00 PM Prep Date: 2/21/2007
 Client ID: Run ID: V-1_070221A SeqNo: 596167

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | ND | 5.0 | µg/L | | | | | | | | | |
| Chloromethane | ND | 5.0 | µg/L | | | | | | | | | |
| Vinyl chloride | ND | 2.0 | µg/L | | | | | | | | | |
| Chloroethane | ND | 5.0 | µg/L | | | | | | | | | |
| Bromomethane | ND | 2.0 | µg/L | | | | | | | | | |
| Trichlorofluoromethane | ND | 2.0 | µg/L | | | | | | | | | |
| Diethyl ether | ND | 5.0 | µg/L | | | | | | | | | |
| Acetone | ND | 10 | µg/L | | | | | | | | | |
| 1,1-Dichloroethene | ND | 1.0 | µg/L | | | | | | | | | |
| Carbon disulfide | ND | 2.0 | µg/L | | | | | | | | | |
| Methylene chloride | ND | 5.0 | µg/L | | | | | | | | | |
| Methyl tert-butyl ether | ND | 2.0 | µg/L | | | | | | | | | |
| trans-1,2-Dichloroethene | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1-Dichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| 2-Butanone | ND | 10 | µg/L | | | | | | | | | |
| 2,2-Dichloropropane | ND | 2.0 | µg/L | | | | | | | | | |
| cis-1,2-Dichloroethene | ND | 2.0 | µg/L | | | | | | | | | |
| Chloroform | ND | 2.0 | µg/L | | | | | | | | | |
| Tetrahydrofuran | ND | 10 | µg/L | | | | | | | | | |
| Bromochloromethane | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1,1-Trichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1-Dichloropropene | ND | 2.0 | µg/L | | | | | | | | | |
| Carbon tetrachloride | ND | 2.0 | µg/L | | | | | | | | | |
| 1,2-Dichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| Benzene | ND | 1.0 | µg/L | | | | | | | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 06-Mar-07

QC SUMMARY REPORT
Method Blank

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

| | | | |
|---------------------------|----|-----|------|
| Trichloroethene | ND | 2.0 | µg/L |
| 1,2-Dichloropropane | ND | 2.0 | µg/L |
| Bromodichloromethane | ND | 2.0 | µg/L |
| Dibromomethane | ND | 2.0 | µg/L |
| 4-Methyl-2-pentanone | ND | 10 | µg/L |
| cis-1,3-Dichloropropene | ND | 1.0 | µg/L |
| Toluene | ND | 2.0 | µg/L |
| trans-1,3-Dichloropropene | ND | 1.0 | µg/L |
| 1,1,2-Trichloroethane | ND | 2.0 | µg/L |
| 1,2-Dibromoethane | ND | 2.0 | µg/L |
| 2-Hexanone | ND | 10 | µg/L |
| 1,3-Dichloropropane | ND | 2.0 | µg/L |
| Tetrachloroethene | ND | 2.0 | µg/L |
| Dibromochloromethane | ND | 2.0 | µg/L |
| Chlorobenzene | ND | 2.0 | µg/L |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | µg/L |
| Ethylbenzene | ND | 2.0 | µg/L |
| m,p-Xylene | ND | 2.0 | µg/L |
| o-Xylene | ND | 2.0 | µg/L |
| Styrene | ND | 2.0 | µg/L |
| Bromoform | ND | 2.0 | µg/L |
| Isopropylbenzene | ND | 2.0 | µg/L |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | µg/L |
| 1,2,3-Trichloropropane | ND | 2.0 | µg/L |
| Bromobenzene | ND | 2.0 | µg/L |
| n-Propylbenzene | ND | 2.0 | µg/L |
| 2-Chlorotoluene | ND | 2.0 | µg/L |
| 4-Chlorotoluene | ND | 2.0 | µg/L |
| 1,3,5-Trimethylbenzene | ND | 2.0 | µg/L |
| tert-Butylbenzene | ND | 2.0 | µg/L |
| 1,2,4-Trimethylbenzene | ND | 2.0 | µg/L |

Qualifiers: ND - Not Detected at the Reporting Limit

S - Spike Recovery outside accepted recovery limits

B - Analyte detected in the associated Method Blank

J - Analyte detected below quantitation limits

R - RPD outside accepted recovery limits

NA - Not applicable where J values or ND results occur

RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Method Blank

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

| Analyte | Reporting Limit | Concentration (µg/L) | Recovery (%) | Acceptance |
|-----------------------------|-----------------|----------------------|--------------|------------|
| sec-Butylbenzene | ND | 2.0 | µg/L | |
| 4-Isopropyltoluene | ND | 2.0 | µg/L | |
| 1,3-Dichlorobenzene | ND | 2.0 | µg/L | |
| 1,4-Dichlorobenzene | ND | 2.0 | µg/L | |
| n-Butylbenzene | ND | 2.0 | µg/L | |
| 1,2-Dichlorobenzene | ND | 2.0 | µg/L | |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | µg/L | |
| 1,2,4-Trichlorobenzene | ND | 2.0 | µg/L | |
| Hexachlorobutadiene | ND | 2.0 | µg/L | |
| Naphthalene | ND | 5.0 | µg/L | |
| 1,2,3-Trichlorobenzene | ND | 2.0 | µg/L | |
| Surr: Dibromofluoromethane | 25.05 | 2.0 | 85 | 116 |
| Surr: 1,2-Dichloroethane-d4 | 26.3 | 2.0 | 77 | 127 |
| Surr: Toluene-d8 | 24.81 | 2.0 | 86 | 114 |
| Surr: 4-Bromofluorobenzene | 24.01 | 2.0 | 79 | 117 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Method Blank

Sample ID: mb-02/22/07 Batch ID: R35964 Test Code: SW8260B Units: µg/L Analysis Date 2/22/2007 1:28:00 PM Prep Date: 2/22/2007
 Client ID: Run ID: V-1_070222A SeqNo: 596193

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | ND | 5.0 | µg/L | | | | | | | | | |
| Chloromethane | ND | 5.0 | µg/L | | | | | | | | | |
| Vinyl chloride | ND | 2.0 | µg/L | | | | | | | | | |
| Chloroethane | ND | 5.0 | µg/L | | | | | | | | | |
| Bromomethane | ND | 2.0 | µg/L | | | | | | | | | |
| Trichlorofluoromethane | ND | 2.0 | µg/L | | | | | | | | | |
| Diethyl ether | ND | 5.0 | µg/L | | | | | | | | | |
| Acetone | ND | 10 | µg/L | | | | | | | | | |
| 1,1-Dichloroethene | ND | 1.0 | µg/L | | | | | | | | | |
| Carbon disulfide | ND | 2.0 | µg/L | | | | | | | | | |
| Methylene chloride | ND | 5.0 | µg/L | | | | | | | | | |
| Methyl tert-butyl ether | ND | 2.0 | µg/L | | | | | | | | | |
| trans-1,2-Dichloroethene | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1-Dichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| 2-Butanone | ND | 10 | µg/L | | | | | | | | | |
| 2,2-Dichloropropane | ND | 2.0 | µg/L | | | | | | | | | |
| cis-1,2-Dichloroethene | ND | 2.0 | µg/L | | | | | | | | | |
| Chloroform | ND | 2.0 | µg/L | | | | | | | | | |
| Tetrahydrofuran | ND | 10 | µg/L | | | | | | | | | |
| Bromochloromethane | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1,1-Trichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1-Dichloropropene | ND | 2.0 | µg/L | | | | | | | | | |
| Carbon tetrachloride | ND | 2.0 | µg/L | | | | | | | | | |
| 1,2-Dichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| Benzene | ND | 1.0 | µg/L | | | | | | | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.
 Work Order: 0702072
 Project: 101960 Textron Gorham
QC SUMMARY REPORT
 Method Blank

| | | | |
|-----------------------------|----|-----|------|
| Trichloroethene | ND | 2.0 | µg/L |
| 1,2-Dichloropropane | ND | 2.0 | µg/L |
| Bromodichloromethane | ND | 2.0 | µg/L |
| Dibromomethane | ND | 2.0 | µg/L |
| 4-Methyl-2-pentanone | ND | 10 | µg/L |
| cis-1,3-Dichloropropene | ND | 1.0 | µg/L |
| Toluene | ND | 2.0 | µg/L |
| trans-1,3-Dichloropropene | ND | 1.0 | µg/L |
| 1,1,2-Trichloroethane | ND | 2.0 | µg/L |
| 1,2-Dibromoethane | ND | 2.0 | µg/L |
| 2-Hexanone | ND | 10 | µg/L |
| 1,3-Dichloropropane | ND | 2.0 | µg/L |
| Tetrachloroethene | ND | 2.0 | µg/L |
| Dibromochloromethane | ND | 2.0 | µg/L |
| Chlorobenzene | ND | 2.0 | µg/L |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | µg/L |
| Ethylbenzene | ND | 2.0 | µg/L |
| m,p-Xylene | ND | 2.0 | µg/L |
| o-Xylene | ND | 2.0 | µg/L |
| Styrene | ND | 2.0 | µg/L |
| Bromoform | ND | 2.0 | µg/L |
| Isopropylbenzene | ND | 2.0 | µg/L |
| 1,1,1,2,2-Tetrachloroethane | ND | 2.0 | µg/L |
| 1,2,3-Trichloropropane | ND | 2.0 | µg/L |
| Bromobenzene | ND | 2.0 | µg/L |
| n-Propylbenzene | ND | 2.0 | µg/L |
| 2-Chlorotoluene | ND | 2.0 | µg/L |
| 4-Chlorotoluene | ND | 2.0 | µg/L |
| 1,3,5-Trimethylbenzene | ND | 2.0 | µg/L |
| tert-Butylbenzene | ND | 2.0 | µg/L |
| 1,2,4-Trimethylbenzene | ND | 2.0 | µg/L |

Qualifiers: ND - Not Detected at the Reporting Limit
 S - Spike Recovery outside accepted recovery limits
 B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits
 R - RPD outside accepted recovery limits
 NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Method Blank

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

| | ND | 2.0 | µg/L | | 25 | 0 | 103 | 85 | 116 | 0 |
|-----------------------------|-------|-----|------|----|----|------|-----|-----|-----|---|
| sec-Butylbenzene | ND | 2.0 | µg/L | | | | | | | |
| 4-Isopropyltoluene | ND | 2.0 | µg/L | | | | | | | |
| 1,3-Dichlorobenzene | ND | 2.0 | µg/L | | | | | | | |
| 1,4-Dichlorobenzene | ND | 2.0 | µg/L | | | | | | | |
| n-Butylbenzene | ND | 2.0 | µg/L | | | | | | | |
| 1,2-Dichlorobenzene | ND | 2.0 | µg/L | | | | | | | |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | µg/L | | | | | | | |
| 1,2,4-Trichlorobenzene | ND | 2.0 | µg/L | | | | | | | |
| Hexachlorobutadiene | ND | 2.0 | µg/L | | | | | | | |
| Naphthalene | ND | 5.0 | µg/L | | | | | | | |
| 1,2,3-Trichlorobenzene | ND | 2.0 | µg/L | | | | | | | |
| Surr: Dibromofluoromethane | 25.63 | 2.0 | µg/L | 25 | 0 | 103 | 85 | 116 | | 0 |
| Surr: 1,2-Dichloroethane-d4 | 26.63 | 2.0 | µg/L | 25 | 0 | 107 | 77 | 127 | | 0 |
| Surr: Toluene-d8 | 25.19 | 2.0 | µg/L | 25 | 0 | 101 | 86 | 114 | | 0 |
| Surr: 4-Bromofluorobenzene | 24.73 | 2.0 | µg/L | 25 | 0 | 98.9 | 79 | 117 | | 0 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Method Blank

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

Sample ID: mb-02/23/07 Batch ID: R35993 Test Code: SW8260B Units: µg/L Analysis Date 2/23/2007 2:19:00 PM Prep Date: 2/23/2007
Client ID: Run ID: V-1_070223A SeqNo: 5965665

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | ND | 5.0 | µg/L | | | | | | | | | |
| Chloromethane | ND | 5.0 | µg/L | | | | | | | | | |
| Vinyl chloride | ND | 2.0 | µg/L | | | | | | | | | |
| Chloroethane | ND | 5.0 | µg/L | | | | | | | | | |
| Bromomethane | ND | 2.0 | µg/L | | | | | | | | | |
| Trichlorofluoromethane | ND | 2.0 | µg/L | | | | | | | | | |
| Diethyl ether | ND | 5.0 | µg/L | | | | | | | | | |
| Acetone | ND | 10 | µg/L | | | | | | | | | |
| 1,1-Dichloroethene | ND | 1.0 | µg/L | | | | | | | | | |
| Carbon disulfide | ND | 2.0 | µg/L | | | | | | | | | |
| Methylene chloride | ND | 5.0 | µg/L | | | | | | | | | |
| Methyl tert-butyl ether | ND | 2.0 | µg/L | | | | | | | | | |
| trans-1,2-Dichloroethene | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1-Dichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| 2-Butanone | ND | 10 | µg/L | | | | | | | | | |
| 2,2-Dichloropropane | ND | 2.0 | µg/L | | | | | | | | | |
| cis-1,2-Dichloroethene | ND | 2.0 | µg/L | | | | | | | | | |
| Chloroform | ND | 2.0 | µg/L | | | | | | | | | |
| Tetrahydrofuran | ND | 10 | µg/L | | | | | | | | | |
| Bromochloromethane | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1,1-Trichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| 1,1-Dichloropropene | ND | 2.0 | µg/L | | | | | | | | | |
| Carbon tetrachloride | ND | 2.0 | µg/L | | | | | | | | | |
| 1,2-Dichloroethane | ND | 2.0 | µg/L | | | | | | | | | |
| Benzene | ND | 1.0 | µg/L | | | | | | | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Method Blank

| | | | |
|---------------------------|----|-----|------|
| Trichloroethene | ND | 2.0 | µg/L |
| 1,2-Dichloropropane | ND | 2.0 | µg/L |
| Bromodichloromethane | ND | 2.0 | µg/L |
| Dibromomethane | ND | 2.0 | µg/L |
| 4-Methyl-2-pentanone | ND | 10 | µg/L |
| cis-1,3-Dichloropropene | ND | 1.0 | µg/L |
| Toluene | ND | 2.0 | µg/L |
| trans-1,3-Dichloropropene | ND | 1.0 | µg/L |
| 1,1,2-Trichloroethane | ND | 2.0 | µg/L |
| 1,2-Dibromoethane | ND | 2.0 | µg/L |
| 2-Hexanone | ND | 10 | µg/L |
| 1,3-Dichloropropane | ND | 2.0 | µg/L |
| Tetrachloroethene | ND | 2.0 | µg/L |
| Dibromochloromethane | ND | 2.0 | µg/L |
| Chlorobenzene | ND | 2.0 | µg/L |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | µg/L |
| Ethylbenzene | ND | 2.0 | µg/L |
| m,p-Xylene | ND | 2.0 | µg/L |
| o-Xylene | ND | 2.0 | µg/L |
| Styrene | ND | 2.0 | µg/L |
| Bromoform | ND | 2.0 | µg/L |
| Isopropylbenzene | ND | 2.0 | µg/L |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | µg/L |
| 1,2,3-Trichloropropane | ND | 2.0 | µg/L |
| Bromobenzene | ND | 2.0 | µg/L |
| n-Propylbenzene | ND | 2.0 | µg/L |
| 2-Chlorotoluene | ND | 2.0 | µg/L |
| 4-Chlorotoluene | ND | 2.0 | µg/L |
| 1,3,5-Trimethylbenzene | ND | 2.0 | µg/L |
| tert-Butylbenzene | ND | 2.0 | µg/L |
| 1,2,4-Trimethylbenzene | ND | 2.0 | µg/L |

Qualifiers: ND - Not Detected at the Reporting Limit

J - Analyte detected below quantitation limits

RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

S - Spike Recovery outside accepted recovery limits

R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank

NA - Not applicable where J values or ND results occur

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Method Blank

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

| Compound | Reporting Limit | Concentration (µg/L) | Recovery (%) | Acceptance |
|-----------------------------|-----------------|----------------------|--------------|------------|
| sec-Butylbenzene | ND | 2.0 | | |
| 4-Isopropyltoluene | ND | 2.0 | | |
| 1,3-Dichlorobenzene | ND | 2.0 | | |
| 1,4-Dichlorobenzene | ND | 2.0 | | |
| n-Butylbenzene | ND | 2.0 | | |
| 1,2-Dichlorobenzene | ND | 2.0 | | |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | |
| Hexachlorobutadiene | ND | 2.0 | | |
| Naphthalene | ND | 5.0 | | |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | |
| Surr: Dibromofluoromethane | 24.83 | 2.0 | 99.3 | 85 |
| Surr: 1,2-Dichloroethane-d4 | 26.39 | 2.0 | 106 | 77 |
| Surr: Toluene-d8 | 25.4 | 2.0 | 102 | 86 |
| Surr: 4-Bromofluorobenzene | 23.56 | 2.0 | 94.2 | 79 |
| | | | | 116 |
| | | | | 127 |
| | | | | 114 |
| | | | | 117 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Laboratory Control Spike - Full List

Sample ID: lcsf-02/21/07 Batch ID: R35962 Test Code: SW8260B Units: µg/L Analysis Date 2/21/2007 2:46:00 PM Prep Date: 2/21/2007

Client ID: Run ID: V-1_070221A SeqNo: 596168

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | 21.08 | 5.0 | µg/L | 20 | 0 | 105 | 10 | 150 | 0 | 0 | 150 | 0 |
| Chloromethane | 19.17 | 5.0 | µg/L | 20 | 0 | 95.8 | 37 | 150 | 0 | 0 | 150 | 0 |
| Vinyl chloride | 21.2 | 2.0 | µg/L | 20 | 0 | 106 | 48 | 150 | 0 | 0 | 150 | 0 |
| Chloroethane | 20.01 | 5.0 | µg/L | 20 | 0 | 100 | 54 | 142 | 0 | 0 | 142 | 0 |
| Bromomethane | 20.46 | 2.0 | µg/L | 20 | 0 | 102 | 51 | 137 | 0 | 0 | 137 | 0 |
| Trichlorofluoromethane | 23.7 | 2.0 | µg/L | 20 | 0 | 118 | 62 | 141 | 0 | 0 | 141 | 0 |
| Diethyl ether | 20.27 | 5.0 | µg/L | 20 | 0 | 101 | 68 | 134 | 0 | 0 | 134 | 0 |
| Acetone | 22.62 | 10 | µg/L | 20 | 0 | 113 | 9 | 150 | 0 | 0 | 150 | 0 |
| 1,1-Dichloroethene | 20.45 | 1.0 | µg/L | 20 | 0 | 102 | 68 | 146 | 0 | 0 | 146 | 0 |
| Carbon disulfide | 21.64 | 2.0 | µg/L | 20 | 0 | 108 | 52 | 131 | 0 | 0 | 131 | 0 |
| Methylene chloride | 22.19 | 5.0 | µg/L | 20 | 0 | 111 | 67 | 138 | 0 | 0 | 138 | 0 |
| Methyl tert-butyl ether | 21.1 | 2.0 | µg/L | 20 | 0 | 106 | 63 | 139 | 0 | 0 | 139 | 0 |
| trans-1,2-Dichloroethene | 24.74 | 2.0 | µg/L | 20 | 0 | 124 | 81 | 126 | 0 | 0 | 126 | 0 |
| 1,1-Dichloroethane | 23.63 | 2.0 | µg/L | 20 | 0 | 118 | 78 | 124 | 0 | 0 | 124 | 0 |
| 2-Butanone | 17.79 | 10 | µg/L | 20 | 0 | 89 | 41 | 150 | 0 | 0 | 150 | 0 |
| 2,2-Dichloropropane | 27.5 | 2.0 | µg/L | 20 | 0 | 138 | 71 | 150 | 0 | 0 | 150 | 0 |
| cis-1,2-Dichloroethene | 20.11 | 2.0 | µg/L | 20 | 0 | 101 | 78 | 121 | 0 | 0 | 121 | 0 |
| Chloroform | 21.68 | 2.0 | µg/L | 20 | 0 | 108 | 82 | 123 | 0 | 0 | 123 | 0 |
| Tetrahydrofuran | 20.82 | 10 | µg/L | 20 | 0 | 104 | 51 | 146 | 0 | 0 | 146 | 0 |
| Bromochloromethane | 19.97 | 2.0 | µg/L | 20 | 0 | 99.8 | 77 | 131 | 0 | 0 | 131 | 0 |
| 1,1,1-Trichloroethane | 24.06 | 2.0 | µg/L | 20 | 0 | 120 | 81 | 127 | 0 | 0 | 127 | 0 |
| 1,1-Dichloropropene | 21.81 | 2.0 | µg/L | 20 | 0 | 109 | 76 | 119 | 0 | 0 | 119 | 0 |
| Carbon tetrachloride | 21.4 | 2.0 | µg/L | 20 | 0 | 107 | 76 | 129 | 0 | 0 | 129 | 0 |
| 1,2-Dichloroethane | 21.64 | 2.0 | µg/L | 20 | 0 | 108 | 76 | 127 | 0 | 0 | 127 | 0 |
| Benzene | 21.83 | 1.0 | µg/L | 20 | 0 | 109 | 81 | 118 | 0 | 0 | 118 | 0 |

Qualifiers: ND - Not Detected at the Reporting Limit
 J - Analyte detected below quantitation limits
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

S - Spike Recovery outside accepted recovery limits
 R - RPD outside accepted recovery limits
 NA - Not applicable where J values or ND results occur

B - Analyte detected in the associated Method Blank

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT

Laboratory Control Spike - Full List

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

| Chemical Name | Concentration (µg/L) | Recovery (%) | Acceptance | Reporting Limit (µg/L) | Qualifiers | | | | |
|---------------------------|----------------------|--------------|------------|------------------------|------------|------|----|-----|---|
| Trichloroethene | 23.22 | 2.0 | µg/L | 20 | 0 | 116 | 81 | 119 | 0 |
| 1,2-Dichloropropane | 21.68 | 2.0 | µg/L | 20 | 0 | 108 | 79 | 120 | 0 |
| Bromodichloromethane | 21.01 | 2.0 | µg/L | 20 | 0 | 105 | 77 | 131 | 0 |
| Dibromomethane | 19.94 | 2.0 | µg/L | 20 | 0 | 99.7 | 76 | 128 | 0 |
| 4-Methyl-2-pentanone | 18.5 | 10 | µg/L | 20 | 0 | 92.5 | 51 | 141 | 0 |
| cis-1,3-Dichloropropene | 20.94 | 1.0 | µg/L | 20 | 0 | 105 | 76 | 120 | 0 |
| Toluene | 21.95 | 2.0 | µg/L | 20 | 0 | 110 | 83 | 119 | 0 |
| trans-1,3-Dichloropropene | 20.93 | 1.0 | µg/L | 20 | 0 | 105 | 66 | 128 | 0 |
| 1,1,2-Trichloroethane | 20.34 | 2.0 | µg/L | 20 | 0 | 102 | 74 | 123 | 0 |
| 1,2-Dibromoethane | 20.65 | 2.0 | µg/L | 20 | 0 | 103 | 72 | 128 | 0 |
| 2-Hexanone | 18.18 | 10 | µg/L | 20 | 0 | 90.9 | 31 | 148 | 0 |
| 1,3-Dichloropropane | 20.59 | 2.0 | µg/L | 20 | 0 | 103 | 76 | 122 | 0 |
| Tetrachloroethene | 22.29 | 2.0 | µg/L | 20 | 0 | 111 | 81 | 124 | 0 |
| Dibromochloromethane | 19.81 | 2.0 | µg/L | 20 | 0 | 99 | 63 | 126 | 0 |
| Chlorobenzene | 19.91 | 2.0 | µg/L | 20 | 0 | 99.6 | 84 | 113 | 0 |
| 1,1,1,2-Tetrachloroethane | 20.76 | 2.0 | µg/L | 20 | 0 | 104 | 73 | 124 | 0 |
| Ethylbenzene | 21.45 | 2.0 | µg/L | 20 | 0 | 107 | 83 | 118 | 0 |
| m,p-Xylene | 43.07 | 2.0 | µg/L | 40 | 0 | 108 | 85 | 116 | 0 |
| o-Xylene | 19.94 | 2.0 | µg/L | 20 | 0 | 99.7 | 84 | 115 | 0 |
| Styrene | 20.05 | 2.0 | µg/L | 20 | 0 | 100 | 81 | 118 | 0 |
| Bromoform | 17.31 | 2.0 | µg/L | 20 | 0 | 86.6 | 55 | 126 | 0 |
| Isopropylbenzene | 23.27 | 2.0 | µg/L | 20 | 0 | 116 | 77 | 125 | 0 |
| 1,1,2,2-Tetrachloroethane | 20.1 | 2.0 | µg/L | 20 | 0 | 100 | 62 | 134 | 0 |
| 1,2,3-Trichloropropane | 20 | 2.0 | µg/L | 20 | 0 | 100 | 62 | 132 | 0 |
| Bromobenzene | 19.36 | 2.0 | µg/L | 20 | 0 | 96.8 | 78 | 119 | 0 |
| n-Propylbenzene | 22.06 | 2.0 | µg/L | 20 | 0 | 110 | 77 | 127 | 0 |
| 2-Chlorotoluene | 21.4 | 2.0 | µg/L | 20 | 0 | 107 | 78 | 118 | 0 |
| 4-Chlorotoluene | 21.95 | 2.0 | µg/L | 20 | 0 | 110 | 77 | 119 | 0 |
| 1,3,5-Trimethylbenzene | 22.57 | 2.0 | µg/L | 20 | 0 | 113 | 80 | 120 | 0 |
| tert-Butylbenzene | 23.1 | 2.0 | µg/L | 20 | 0 | 116 | 81 | 120 | 0 |
| 1,2,4-Trimethylbenzene | 22.04 | 2.0 | µg/L | 20 | 0 | 110 | 80 | 118 | 0 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Laboratory Control Spike - Full List

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

| Compound | Concentration (µg/L) | Recovery (%) | Recovery (µg/L) | Spiked Amount (µg/L) | Recovery (%) | Recovery (µg/L) | Spiked Amount (µg/L) |
|-----------------------------|----------------------|--------------|-----------------|----------------------|--------------|-----------------|----------------------|
| sec-Butylbenzene | 19.73 | 2.0 | 20 | 20 | 98.6 | 82 | 123 |
| 4-Isopropyltoluene | 22.98 | 2.0 | 20 | 20 | 115 | 80 | 126 |
| 1,3-Dichlorobenzene | 20.22 | 2.0 | 20 | 20 | 101 | 84 | 115 |
| 1,4-Dichlorobenzene | 20.42 | 2.0 | 20 | 20 | 102 | 79 | 117 |
| n-Butylbenzene | 19.34 | 2.0 | 20 | 20 | 96.7 | 76 | 128 |
| 1,2-Dichlorobenzene | 19.92 | 2.0 | 20 | 20 | 99.6 | 81 | 117 |
| 1,2-Dibromo-3-chloropropane | 18.61 | 5.0 | 20 | 20 | 93 | 47 | 136 |
| 1,2,4-Trichlorobenzene | 20.38 | 2.0 | 20 | 20 | 102 | 73 | 126 |
| Hexachlorobutadiene | 20.28 | 2.0 | 20 | 20 | 101 | 77 | 134 |
| Naphthalene | 20.73 | 5.0 | 20 | 20 | 104 | 58 | 138 |
| 1,2,3-Trichlorobenzene | 20.62 | 2.0 | 20 | 20 | 103 | 76 | 124 |
| Surr: Dibromofluoromethane | 25.69 | 2.0 | 25 | 25 | 103 | 85 | 116 |
| Surr: 1,2-Dichloroethane-d4 | 27.39 | 2.0 | 25 | 25 | 110 | 77 | 127 |
| Surr: Toluene-d8 | 25.13 | 2.0 | 25 | 25 | 101 | 86 | 114 |
| Surr: 4-Bromofluorobenzene | 24.25 | 2.0 | 25 | 25 | 97 | 79 | 117 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Laboratory Control Spike - Full List

Sample ID: Icsf-02/22/07 Batch ID: R35964 Test Code: SW8260B Units: µg/L Analysis Date 2/22/2007 12:18:00 PM Prep Date: 2/22/2007
 Client ID: Run ID: V-1_070222A SeqNo: 596194

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | 16.24 | 5.0 | µg/L | 20 | 0 | 81.2 | 10 | 150 | 0 | 0 | | |
| Chloromethane | 16.21 | 5.0 | µg/L | 20 | 0 | 81 | 37 | 150 | 0 | 0 | | |
| Vinyl chloride | 17.6 | 2.0 | µg/L | 20 | 0 | 88 | 48 | 150 | 0 | 0 | | |
| Chloroethane | 18.68 | 5.0 | µg/L | 20 | 0 | 93.4 | 54 | 142 | 0 | 0 | | |
| Bromomethane | 18.53 | 2.0 | µg/L | 20 | 0 | 92.6 | 51 | 137 | 0 | 0 | | |
| Trichlorofluoromethane | 19.6 | 2.0 | µg/L | 20 | 0 | 98 | 62 | 141 | 0 | 0 | | |
| Diethyl ether | 20.64 | 5.0 | µg/L | 20 | 0 | 103 | 68 | 134 | 0 | 0 | | |
| Acetone | 20.99 | 10 | µg/L | 20 | 0 | 105 | 9 | 150 | 0 | 0 | | |
| 1,1-Dichloroethene | 17.2 | 1.0 | µg/L | 20 | 0 | 86 | 68 | 146 | 0 | 0 | | |
| Carbon disulfide | 18.52 | 2.0 | µg/L | 20 | 0 | 92.6 | 52 | 131 | 0 | 0 | | |
| Methylene chloride | 20.46 | 5.0 | µg/L | 20 | 0 | 102 | 67 | 138 | 0 | 0 | | |
| Methyl tert-butyl ether | 20.94 | 2.0 | µg/L | 20 | 0 | 105 | 63 | 139 | 0 | 0 | | |
| trans-1,2-Dichloroethene | 22.09 | 2.0 | µg/L | 20 | 0 | 110 | 81 | 126 | 0 | 0 | | |
| 1,1-Dichloroethane | 21.87 | 2.0 | µg/L | 20 | 0 | 109 | 78 | 124 | 0 | 0 | | |
| 2-Butanone | 21.48 | 10 | µg/L | 20 | 0 | 107 | 41 | 150 | 0 | 0 | | |
| 2,2-Dichloropropane | 25.31 | 2.0 | µg/L | 20 | 0 | 127 | 71 | 150 | 0 | 0 | | |
| cis-1,2-Dichloroethene | 18.83 | 2.0 | µg/L | 20 | 0 | 94.2 | 78 | 121 | 0 | 0 | | |
| Chloroform | 20.98 | 2.0 | µg/L | 20 | 0 | 105 | 82 | 123 | 0 | 0 | | |
| Tetrahydrofuran | 21.18 | 10 | µg/L | 20 | 0 | 106 | 51 | 146 | 0 | 0 | | |
| Bromochloromethane | 20.02 | 2.0 | µg/L | 20 | 0 | 100 | 77 | 131 | 0 | 0 | | |
| 1,1,1-Trichloroethane | 21.77 | 2.0 | µg/L | 20 | 0 | 109 | 81 | 127 | 0 | 0 | | |
| 1,1-Dichloropropene | 19.82 | 2.0 | µg/L | 20 | 0 | 99.1 | 76 | 119 | 0 | 0 | | |
| Carbon tetrachloride | 19.22 | 2.0 | µg/L | 20 | 0 | 96.1 | 76 | 129 | 0 | 0 | | |
| 1,2-Dichloroethane | 22.06 | 2.0 | µg/L | 20 | 0 | 110 | 76 | 127 | 0 | 0 | | |
| Benzene | 20.23 | 1.0 | µg/L | 20 | 0 | 101 | 81 | 118 | 0 | 0 | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Laboratory Control Spike - Full List

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

| Compound | Concentration (µg/L) | Recovery (%) | Method | QC Value |
|---------------------------|----------------------|--------------|--------|----------|
| Trichloroethene | 20.45 | 2.0 | µg/L | 119 |
| 1,2-Dichloropropane | 19.96 | 2.0 | µg/L | 120 |
| Bromodichloromethane | 20.35 | 2.0 | µg/L | 131 |
| Dibromomethane | 19.6 | 2.0 | µg/L | 128 |
| 4-Methyl-2-pentanone | 19.03 | 10 | µg/L | 141 |
| cis-1,3-Dichloropropene | 19.03 | 1.0 | µg/L | 120 |
| Toluene | 19.8 | 2.0 | µg/L | 119 |
| trans-1,3-Dichloropropene | 19.8 | 1.0 | µg/L | 128 |
| 1,1,2-Trichloroethane | 20.39 | 2.0 | µg/L | 123 |
| 1,2-Dibromoethane | 20.94 | 2.0 | µg/L | 128 |
| 2-Hexanone | 18.04 | 10 | µg/L | 148 |
| 1,3-Dichloropropane | 19.76 | 2.0 | µg/L | 122 |
| Tetrachloroethene | 20.03 | 2.0 | µg/L | 124 |
| Dibromochloromethane | 18.27 | 2.0 | µg/L | 126 |
| Chlorobenzene | 18.47 | 2.0 | µg/L | 113 |
| 1,1,1,2-Tetrachloroethane | 18.74 | 2.0 | µg/L | 124 |
| Ethylbenzene | 19.73 | 2.0 | µg/L | 118 |
| m,p-Xylene | 40.32 | 2.0 | µg/L | 116 |
| o-Xylene | 18.71 | 2.0 | µg/L | 115 |
| Styrene | 19.38 | 2.0 | µg/L | 118 |
| Bromoform | 17.07 | 2.0 | µg/L | 126 |
| Isopropylbenzene | 20.36 | 2.0 | µg/L | 125 |
| 1,1,1,2-Tetrachloroethane | 18.21 | 2.0 | µg/L | 134 |
| 1,2,3-Trichloropropane | 19.34 | 2.0 | µg/L | 132 |
| Bromobenzene | 18.84 | 2.0 | µg/L | 119 |
| n-Propylbenzene | 19.34 | 2.0 | µg/L | 127 |
| 2-Chlorotoluene | 19.29 | 2.0 | µg/L | 118 |
| 4-Chlorotoluene | 19.62 | 2.0 | µg/L | 119 |
| 1,3,5-Trimethylbenzene | 20.01 | 2.0 | µg/L | 120 |
| tert-Butylbenzene | 20.46 | 2.0 | µg/L | 120 |
| 1,2,4-Trimethylbenzene | 19.7 | 2.0 | µg/L | 118 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Laboratory Control Spike - Full List

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

Sample ID: **iesf-02/23/07** Batch ID: **R35993** Test Code: **SW8260B** Units: **µg/L** Analysis Date **2/23/2007 12:35:00 PM** Prep Date: **2/23/2007**
Client ID: **SeqNo: 596566** Run ID: **V-1_070223A**

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | 15.43 | 5.0 | µg/L | 20 | 0 | 77.2 | 10 | 150 | 0 | 0 | 150 | 0 |
| Chloromethane | 17.02 | 5.0 | µg/L | 20 | 0 | 85.1 | 37 | 150 | 0 | 0 | 150 | 0 |
| Vinyl chloride | 19.06 | 2.0 | µg/L | 20 | 0 | 95.3 | 48 | 150 | 0 | 0 | 150 | 0 |
| Chloroethane | 20.04 | 5.0 | µg/L | 20 | 0 | 100 | 54 | 142 | 0 | 0 | 142 | 0 |
| Bromomethane | 19.28 | 2.0 | µg/L | 20 | 0 | 96.4 | 51 | 137 | 0 | 0 | 137 | 0 |
| Trichlorofluoromethane | 21.68 | 2.0 | µg/L | 20 | 0 | 108 | 62 | 141 | 0 | 0 | 141 | 0 |
| Diethyl ether | 18.88 | 5.0 | µg/L | 20 | 0 | 94.4 | 68 | 134 | 0 | 0 | 134 | 0 |
| Acetone | 14.65 | 10 | µg/L | 20 | 0 | 73.2 | 9 | 150 | 0 | 0 | 150 | 0 |
| 1,1-Dichloroethene | 18.7 | 1.0 | µg/L | 20 | 0 | 93.5 | 68 | 146 | 0 | 0 | 146 | 0 |
| Carbon disulfide | 20.28 | 2.0 | µg/L | 20 | 0 | 101 | 52 | 131 | 0 | 0 | 131 | 0 |
| Methylene chloride | 21.88 | 5.0 | µg/L | 20 | 0 | 109 | 67 | 138 | 0 | 0 | 138 | 0 |
| Methyl tert-butyl ether | 21.55 | 2.0 | µg/L | 20 | 0 | 108 | 63 | 139 | 0 | 0 | 139 | 0 |
| trans-1,2-Dichloroethene | 22.93 | 2.0 | µg/L | 20 | 0 | 115 | 81 | 126 | 0 | 0 | 126 | 0 |
| 1,1-Dichloroethane | 22.97 | 2.0 | µg/L | 20 | 0 | 115 | 78 | 124 | 0 | 0 | 124 | 0 |
| 2-Butanone | 15.28 | 10 | µg/L | 20 | 0 | 76.4 | 41 | 150 | 0 | 0 | 150 | 0 |
| 2,2-Dichloropropane | 27.34 | 2.0 | µg/L | 20 | 0 | 137 | 71 | 150 | 0 | 0 | 150 | 0 |
| cis-1,2-Dichloroethene | 19.43 | 2.0 | µg/L | 20 | 0 | 97.2 | 78 | 121 | 0 | 0 | 121 | 0 |
| Chloroform | 21.77 | 2.0 | µg/L | 20 | 0 | 109 | 82 | 123 | 0 | 0 | 123 | 0 |
| Tetrahydrofuran | 16.64 | 10 | µg/L | 20 | 0 | 83.2 | 51 | 146 | 0 | 0 | 146 | 0 |
| Bromochloromethane | 21.29 | 2.0 | µg/L | 20 | 0 | 106 | 77 | 131 | 0 | 0 | 131 | 0 |
| 1,1,1-Trichloroethane | 23.4 | 2.0 | µg/L | 20 | 0 | 117 | 81 | 127 | 0 | 0 | 127 | 0 |
| 1,1-Dichloropropene | 20.56 | 2.0 | µg/L | 20 | 0 | 103 | 76 | 119 | 0 | 0 | 119 | 0 |
| Carbon tetrachloride | 21.12 | 2.0 | µg/L | 20 | 0 | 106 | 76 | 129 | 0 | 0 | 129 | 0 |
| 1,2-Dichloroethane | 21.42 | 2.0 | µg/L | 20 | 0 | 107 | 76 | 127 | 0 | 0 | 127 | 0 |
| Benzene | 21.03 | 1.0 | µg/L | 20 | 0 | 105 | 81 | 118 | 0 | 0 | 118 | 0 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Laboratory Control Spike - Full List

| Chemical Name | Reporting Limit | Concentration | Recovery | Acceptance | Spikes | Recovery | Acceptance | Spikes | Recovery | Acceptance | Spikes | Recovery | Acceptance | Spikes | Recovery | Acceptance | Spikes | Recovery | Acceptance | |
|---------------------------|-----------------|---------------|----------|------------|--------|----------|------------|--------|----------|------------|--------|----------|------------|--------|----------|------------|--------|----------|------------|---|
| Trichloroethene | 22.08 | 2.0 | µg/L | 20 | 0 | 110 | 81 | 119 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 1,2-Dichloropropane | 20.44 | 2.0 | µg/L | 20 | 0 | 102 | 79 | 120 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Bromodichloromethane | 20.44 | 2.0 | µg/L | 20 | 0 | 102 | 77 | 131 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Dibromomethane | 19.16 | 2.0 | µg/L | 20 | 0 | 95.8 | 76 | 128 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 4-Methyl-2-pentanone | 14.8 | 10 | µg/L | 20 | 0 | 74 | 51 | 141 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| cis-1,3-Dichloropropene | 20.12 | 1.0 | µg/L | 20 | 0 | 101 | 76 | 120 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Toluene | 21.6 | 2.0 | µg/L | 20 | 0 | 108 | 83 | 119 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| trans-1,3-Dichloropropene | 20 | 1.0 | µg/L | 20 | 0 | 100 | 66 | 128 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 1,1,2-Trichloroethane | 20 | 2.0 | µg/L | 20 | 0 | 100 | 74 | 123 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 1,2-Dibromoethane | 19.75 | 2.0 | µg/L | 20 | 0 | 98.8 | 72 | 128 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 2-Hexanone | 15.04 | 10 | µg/L | 20 | 0 | 75.2 | 31 | 148 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 1,3-Dichloropropane | 19.93 | 2.0 | µg/L | 20 | 0 | 99.7 | 76 | 122 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Tetrachloroethene | 21.64 | 2.0 | µg/L | 20 | 0 | 108 | 81 | 124 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Dibromochloromethane | 19.61 | 2.0 | µg/L | 20 | 0 | 98 | 63 | 126 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Chlorobenzene | 19.88 | 2.0 | µg/L | 20 | 0 | 99.4 | 84 | 113 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 1,1,1,2-Tetrachloroethane | 20.01 | 2.0 | µg/L | 20 | 0 | 100 | 73 | 124 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Ethylbenzene | 21.26 | 2.0 | µg/L | 20 | 0 | 106 | 83 | 118 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| m,p-Xylene | 42.16 | 2.0 | µg/L | 40 | 0 | 105 | 85 | 116 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| o-Xylene | 19.94 | 2.0 | µg/L | 20 | 0 | 99.7 | 84 | 115 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Styrene | 20.37 | 2.0 | µg/L | 20 | 0 | 102 | 81 | 118 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Bromoform | 16.71 | 2.0 | µg/L | 20 | 0 | 83.6 | 55 | 126 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Isopropylbenzene | 22.03 | 2.0 | µg/L | 20 | 0 | 110 | 77 | 125 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 1,1,2,2-Tetrachloroethane | 16.78 | 2.0 | µg/L | 20 | 0 | 83.9 | 62 | 134 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 1,2,3-Trichloropropane | 18.26 | 2.0 | µg/L | 20 | 0 | 91.3 | 62 | 132 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Bromobenzene | 18.81 | 2.0 | µg/L | 20 | 0 | 94.1 | 78 | 119 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| n-Propylbenzene | 21 | 2.0 | µg/L | 20 | 0 | 105 | 77 | 127 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 2-Chlorotoluene | 20.66 | 2.0 | µg/L | 20 | 0 | 103 | 78 | 118 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 4-Chlorotoluene | 20.8 | 2.0 | µg/L | 20 | 0 | 104 | 77 | 119 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 1,3,5-Trimethylbenzene | 21.4 | 2.0 | µg/L | 20 | 0 | 107 | 80 | 120 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| tert-Butylbenzene | 21.88 | 2.0 | µg/L | 20 | 0 | 109 | 81 | 120 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 1,2,4-Trimethylbenzene | 20.8 | 2.0 | µg/L | 20 | 0 | 104 | 80 | 118 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Laboratory Control Spike - Full List

| Compound | Concentration (µg/L) | Recovery (%) | Acceptance | Count | Count |
|-----------------------------|----------------------|--------------|------------|-------|-------|
| sec-Butylbenzene | 18.23 | 2.0 | µg/L | 20 | 82 |
| 4-Isopropyltoluene | 21.45 | 2.0 | µg/L | 20 | 80 |
| 1,3-Dichlorobenzene | 19.37 | 2.0 | µg/L | 20 | 84 |
| 1,4-Dichlorobenzene | 19.03 | 2.0 | µg/L | 20 | 79 |
| n-Butylbenzene | 17.29 | 2.0 | µg/L | 20 | 76 |
| 1,2-Dichlorobenzene | 18.93 | 2.0 | µg/L | 20 | 81 |
| 1,2-Dibromo-3-chloropropane | 15.41 | 5.0 | µg/L | 20 | 47 |
| 1,2,4-Trichlorobenzene | 17.9 | 2.0 | µg/L | 20 | 73 |
| Hexachlorobutadiene | 17.35 | 2.0 | µg/L | 20 | 77 |
| Naphthalene | 17.05 | 5.0 | µg/L | 20 | 58 |
| 1,2,3-Trichlorobenzene | 16.93 | 2.0 | µg/L | 20 | 76 |
| Surr: Dibromofluoromethane | 25.04 | 2.0 | µg/L | 25 | 85 |
| Surr: 1,2-Dichloroethane-d4 | 25.82 | 2.0 | µg/L | 25 | 77 |
| Surr: Toluene-d8 | 25.56 | 2.0 | µg/L | 25 | 86 |
| Surr: 4-Bromofluorobenzene | 24.62 | 2.0 | µg/L | 25 | 79 |

Qualifiers: ND - Not Detected at the Reporting Limit
 J - Analyte detected below quantitation limits
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.
 S - Spike Recovery outside accepted recovery limits
 R - RPD outside accepted recovery limits
 B - Analyte detected in the associated Method Blank
 NA - Not applicable where J values or ND results occur

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Laboratory Control Spike Duplicate - Full List

Sample ID: lcsdf-02/23/07 Batch ID: R35983 Test Code: SW8260B Units: µg/L Analysis Date 2/23/2007 1:09:00 PM Prep Date: 2/23/2007

Client ID: Run ID: V-1_070223A SeqNo: 596567

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|--------|----------|-----|
| Dichlorodifluoromethane | 16.78 | 5.0 | µg/L | 20 | 0 | 83.9 | 10 | 150 | 15.43 | 8.38 | 25 | |
| Chloromethane | 17.69 | 5.0 | µg/L | 20 | 0 | 88.4 | 37 | 150 | 17.02 | 3.86 | 25 | |
| Vinyl chloride | 19.34 | 2.0 | µg/L | 20 | 0 | 96.7 | 48 | 150 | 19.06 | 1.46 | 25 | |
| Chloroethane | 18.49 | 5.0 | µg/L | 20 | 0 | 92.5 | 54 | 142 | 20.04 | 8.05 | 25 | |
| Bromomethane | 19.3 | 2.0 | µg/L | 20 | 0 | 96.5 | 51 | 137 | 19.28 | 0.104 | 25 | |
| Trichlorofluoromethane | 21.62 | 2.0 | µg/L | 20 | 0 | 108 | 62 | 141 | 21.68 | 0.277 | 25 | |
| Acetone | 16.26 | 10 | µg/L | 20 | 0 | 81.3 | 9 | 150 | 14.65 | 10.4 | 25 | |
| 1,1-Dichloroethene | 18.55 | 1.0 | µg/L | 20 | 0 | 92.8 | 68 | 146 | 18.7 | 0.805 | 25 | |
| Carbon disulfide | 20.33 | 2.0 | µg/L | 20 | 0 | 102 | 52 | 131 | 20.28 | 0.246 | 25 | |
| Methylene chloride | 22.14 | 5.0 | µg/L | 20 | 0 | 111 | 67 | 138 | 21.88 | 1.18 | 25 | |
| Methyl tert-butyl ether | 21.45 | 2.0 | µg/L | 20 | 0 | 107 | 63 | 139 | 21.55 | 0.465 | 25 | |
| trans-1,2-Dichloroethene | 23.12 | 2.0 | µg/L | 20 | 0 | 116 | 81 | 126 | 22.93 | 0.825 | 25 | |
| 1,1-Dichloroethane | 23.22 | 2.0 | µg/L | 20 | 0 | 116 | 78 | 124 | 22.97 | 1.08 | 25 | |
| 2-Butanone | 15.15 | 10 | µg/L | 20 | 0 | 75.8 | 41 | 150 | 15.28 | 0.854 | 25 | |
| 2,2-Dichloropropane | 26.83 | 2.0 | µg/L | 20 | 0 | 134 | 71 | 150 | 27.34 | 1.88 | 25 | |
| cis-1,2-Dichloroethene | 19.61 | 2.0 | µg/L | 20 | 0 | 98 | 78 | 121 | 19.43 | 0.922 | 25 | |
| Chloroform | 21 | 2.0 | µg/L | 20 | 0 | 105 | 82 | 123 | 21.77 | 3.6 | 25 | |
| Bromochloromethane | 20.65 | 2.0 | µg/L | 20 | 0 | 103 | 77 | 131 | 21.29 | 3.05 | 25 | |
| 1,1,1-Trichloroethane | 23.22 | 2.0 | µg/L | 20 | 0 | 116 | 81 | 127 | 23.4 | 0.772 | 25 | |
| 1,1-Dichloropropene | 20.55 | 2.0 | µg/L | 20 | 0 | 103 | 76 | 119 | 20.56 | 0.0486 | 25 | |
| Carbon tetrachloride | 20.73 | 2.0 | µg/L | 20 | 0 | 104 | 76 | 129 | 21.12 | 1.86 | 25 | |
| 1,2-Dichloroethane | 21.76 | 2.0 | µg/L | 20 | 0 | 109 | 76 | 127 | 21.42 | 1.57 | 25 | |
| Benzene | 20.64 | 1.0 | µg/L | 20 | 0 | 103 | 81 | 118 | 21.03 | 1.87 | 25 | |
| Trichloroethene | 21.44 | 2.0 | µg/L | 20 | 0 | 107 | 81 | 119 | 22.08 | 2.94 | 25 | |
| 1,2-Dichloropropane | 20.98 | 2.0 | µg/L | 20 | 0 | 105 | 79 | 120 | 20.44 | 2.61 | 25 | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Laboratory Control Spike Duplicate - Full List

| Chemical Name | Reporting Limit | Concentration | Recovery | Method | Concentration | Method | Concentration | Method | Concentration | Method | |
|---------------------------|-----------------|---------------|----------|--------|---------------|--------|---------------|--------|---------------|--------|----|
| Bromodichloromethane | 20.49 | 2.0 | µg/L | 20 | 0 | 102 | 77 | 131 | 20.44 | 0.244 | 25 |
| Dibromomethane | 18.82 | 2.0 | µg/L | 20 | 0 | 94.1 | 76 | 128 | 19.16 | 1.79 | 25 |
| 4-Methyl-2-pentanone | 16.64 | 10 | µg/L | 20 | 0 | 83.2 | 51 | 141 | 14.8 | 11.7 | 25 |
| cis-1,3-Dichloropropene | 19.67 | 1.0 | µg/L | 20 | 0 | 98.4 | 76 | 120 | 20.12 | 2.26 | 25 |
| Toluene | 21.28 | 2.0 | µg/L | 20 | 0 | 106 | 83 | 119 | 21.6 | 1.49 | 25 |
| trans-1,3-Dichloropropene | 19.79 | 1.0 | µg/L | 20 | 0 | 99 | 66 | 128 | 20 | 1.06 | 25 |
| 1,1,2-Trichloroethane | 19.36 | 2.0 | µg/L | 20 | 0 | 96.8 | 74 | 123 | 20 | 3.25 | 25 |
| 1,2-Dibromoethane | 19.54 | 2.0 | µg/L | 20 | 0 | 97.7 | 72 | 128 | 19.75 | 1.07 | 25 |
| 2-Hexanone | 14.97 | 10 | µg/L | 20 | 0 | 74.8 | 31 | 148 | 15.04 | 0.467 | 25 |
| 1,3-Dichloropropane | 19.47 | 2.0 | µg/L | 20 | 0 | 97.4 | 76 | 122 | 19.93 | 2.34 | 25 |
| Tetrachloroethene | 20.82 | 2.0 | µg/L | 20 | 0 | 104 | 81 | 124 | 21.64 | 3.86 | 25 |
| Dibromochloromethane | 18.49 | 2.0 | µg/L | 20 | 0 | 92.5 | 63 | 126 | 19.61 | 5.88 | 25 |
| Chlorobenzene | 19.43 | 2.0 | µg/L | 20 | 0 | 97.2 | 84 | 113 | 19.88 | 2.29 | 25 |
| 1,1,1,2-Tetrachloroethane | 19.75 | 2.0 | µg/L | 20 | 0 | 98.8 | 73 | 124 | 20.01 | 1.31 | 25 |
| Ethylbenzene | 20.76 | 2.0 | µg/L | 20 | 0 | 104 | 83 | 118 | 21.26 | 2.38 | 25 |
| m,p-Xylene | 41.65 | 2.0 | µg/L | 40 | 0 | 104 | 85 | 116 | 42.16 | 1.22 | 25 |
| o-Xylene | 19.69 | 2.0 | µg/L | 20 | 0 | 98.4 | 84 | 115 | 19.94 | 1.26 | 25 |
| Styrene | 19.55 | 2.0 | µg/L | 20 | 0 | 97.8 | 81 | 118 | 20.37 | 4.11 | 25 |
| Bromoform | 15.81 | 2.0 | µg/L | 20 | 0 | 79 | 55 | 126 | 16.71 | 5.54 | 25 |
| Isopropylbenzene | 21.93 | 2.0 | µg/L | 20 | 0 | 110 | 77 | 125 | 22.03 | 0.455 | 25 |
| 1,1,2,2-Tetrachloroethane | 17.16 | 2.0 | µg/L | 20 | 0 | 85.8 | 62 | 134 | 16.78 | 2.24 | 25 |
| 1,2,3-Trichloropropane | 17.42 | 2.0 | µg/L | 20 | 0 | 87.1 | 62 | 132 | 18.26 | 4.71 | 25 |
| Bromobenzene | 18.77 | 2.0 | µg/L | 20 | 0 | 93.8 | 78 | 119 | 18.81 | 0.213 | 25 |
| n-Propylbenzene | 21.46 | 2.0 | µg/L | 20 | 0 | 107 | 77 | 127 | 21 | 2.17 | 25 |
| 2-Chlorotoluene | 20.62 | 2.0 | µg/L | 20 | 0 | 103 | 78 | 118 | 20.66 | 0.194 | 25 |
| 4-Chlorotoluene | 20.95 | 2.0 | µg/L | 20 | 0 | 105 | 77 | 119 | 20.8 | 0.719 | 25 |
| 1,3,5-Trimethylbenzene | 20.87 | 2.0 | µg/L | 20 | 0 | 104 | 80 | 120 | 21.4 | 2.51 | 25 |
| tert-Butylbenzene | 20.88 | 2.0 | µg/L | 20 | 0 | 104 | 81 | 120 | 21.88 | 4.68 | 25 |
| 1,2,4-Trimethylbenzene | 20.48 | 2.0 | µg/L | 20 | 0 | 102 | 80 | 118 | 20.8 | 1.55 | 25 |
| sec-Butylbenzene | 17.89 | 2.0 | µg/L | 20 | 0 | 89.4 | 82 | 123 | 18.23 | 1.88 | 25 |
| 4-Isopropyltoluene | 20.59 | 2.0 | µg/L | 20 | 0 | 103 | 80 | 126 | 21.45 | 4.09 | 25 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Laboratory Control Spike Duplicate - Full List

| | 19.17 | 2.0 | µg/L | 20 | 0 | 95.8 | 84 | 115 | 19.37 | 1.04 | 25 |
|-----------------------------|-------|-----|------|----|---|------|----|-----|-------|------|----|
| 1,3-Dichlorobenzene | 18.82 | 2.0 | µg/L | 20 | 0 | 94.1 | 79 | 117 | 19.03 | 1.11 | 25 |
| 1,4-Dichlorobenzene | 17.01 | 2.0 | µg/L | 20 | 0 | 85 | 76 | 128 | 17.29 | 1.63 | 25 |
| n-Butylbenzene | 18.01 | 2.0 | µg/L | 20 | 0 | 90 | 81 | 117 | 18.93 | 4.98 | 25 |
| 1,2-Dichlorobenzene | 15.7 | 5.0 | µg/L | 20 | 0 | 78.5 | 47 | 136 | 15.41 | 1.86 | 25 |
| 1,2-Dibromo-3-chloropropane | 16.55 | 2.0 | µg/L | 20 | 0 | 82.8 | 73 | 126 | 17.9 | 7.84 | 25 |
| 1,2,4-Trichlorobenzene | 16.55 | 2.0 | µg/L | 20 | 0 | 82.8 | 77 | 134 | 17.35 | 4.72 | 25 |
| Hexachlorobutadiene | 16.75 | 5.0 | µg/L | 20 | 0 | 83.8 | 58 | 138 | 17.05 | 1.78 | 25 |
| Naphthalene | 15.69 | 2.0 | µg/L | 20 | 0 | 78.5 | 76 | 124 | 16.93 | 7.6 | 25 |
| 1,2,3-Trichlorobenzene | 25.74 | 2.0 | µg/L | 25 | 0 | 103 | 85 | 116 | 0 | 0 | 0 |
| Surr: Dibromofluoromethane | 26.86 | 2.0 | µg/L | 25 | 0 | 107 | 77 | 127 | 0 | 0 | 0 |
| Surr: 1,2-Dichloroethane-d4 | 25.15 | 2.0 | µg/L | 25 | 0 | 101 | 86 | 114 | 0 | 0 | 0 |
| Surr: Toluene-d8 | 24.35 | 2.0 | µg/L | 25 | 0 | 97.4 | 79 | 117 | 0 | 0 | 0 |
| Surr: 4-Bromofluorobenzene | | | | | | | | | | | |

Qualifiers: ND - Not Detected at the Reporting Limit

J - Analyte detected below quantitation limits

RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

S - Spike Recovery outside accepted recovery limits

R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank

NA - Not applicable where J values or ND results occur

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike - Full List

Sample ID: 0702072-15Amsf Batch ID: R35964 Test Code: SW8260B Units: µg/L Analysis Date 2/22/2007 10:05:00 PM Prep Date: 2/17/2007
 Client ID: MW-209D Run ID: V-1_070222A SeqNo: 596251

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | 179.5 | 50 | µg/L | 200 | 0 | 89.8 | 16 | 150 | 0 | | | |
| Chloromethane | 183.6 | 50 | µg/L | 200 | 0 | 91.8 | 35 | 150 | 0 | | | |
| Vinyl chloride | 210.2 | 20 | µg/L | 200 | 0 | 105 | 49 | 150 | 0 | | | |
| Chloroethane | 196.3 | 50 | µg/L | 200 | 0 | 98.2 | 58 | 147 | 0 | | | |
| Bromomethane | 188.6 | 20 | µg/L | 200 | 0 | 94.3 | 49 | 142 | 0 | | | |
| Trichlorofluoromethane | 215.4 | 20 | µg/L | 200 | 0 | 108 | 57 | 149 | 0 | | | |
| Diethyl ether | 176.1 | 50 | µg/L | 200 | 0 | 88 | 66 | 136 | 0 | | | |
| Acetone | 140.3 | 100 | µg/L | 200 | 0 | 70.2 | 16 | 150 | 0 | | | |
| 1,1-Dichloroethene | 184.2 | 10 | µg/L | 200 | 0 | 92.1 | 70 | 150 | 0 | | | |
| Carbon disulfide | 202.5 | 20 | µg/L | 200 | 0 | 101 | 47 | 135 | 0 | | | |
| Methylene chloride | 213.5 | 50 | µg/L | 200 | 0 | 107 | 66 | 142 | 0 | | | |
| Methyl tert-butyl ether | 210.6 | 20 | µg/L | 200 | 0 | 105 | 63 | 138 | 0 | | | |
| trans-1,2-Dichloroethene | 227.8 | 20 | µg/L | 200 | 0 | 114 | 78 | 135 | 0 | | | |
| 1,1-Dichloroethane | 221.6 | 20 | µg/L | 200 | 0 | 111 | 76 | 131 | 0 | | | |
| 2-Butanone | 123.5 | 100 | µg/L | 200 | 0 | 61.8 | 51 | 142 | 0 | | | |
| 2,2-Dichloropropane | 217.9 | 20 | µg/L | 200 | 0 | 109 | 60 | 149 | 0 | | | |
| cis-1,2-Dichloroethene | 196.5 | 20 | µg/L | 200 | 0 | 98.2 | 74 | 128 | 0 | | | |
| Chloroform | 194.4 | 20 | µg/L | 200 | 0 | 97.2 | 80 | 129 | 0 | | | |
| Tetrahydrofuran | 144.7 | 100 | µg/L | 200 | 0 | 72.4 | 53 | 145 | 0 | | | |
| Bromochloromethane | 177.2 | 20 | µg/L | 200 | 0 | 88.6 | 78 | 130 | 0 | | | |
| 1,1,1-Trichloroethane | 222.1 | 20 | µg/L | 200 | 0 | 111 | 77 | 139 | 0 | | | |
| 1,1-Dichloropropene | 185.4 | 20 | µg/L | 200 | 0 | 92.7 | 74 | 127 | 0 | | | |
| Carbon tetrachloride | 184.6 | 20 | µg/L | 200 | 0 | 92.3 | 73 | 138 | 0 | | | |
| 1,2-Dichloroethane | 205.2 | 20 | µg/L | 200 | 0 | 103 | 75 | 130 | 0 | | | |
| Benzene | 202.1 | 10 | µg/L | 200 | 0 | 101 | 79 | 123 | 0 | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit, defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike - Full List

| Compound | Concentration (µg/L) | Volume (mL) | Recovery (%) | Spikes | Matrix Spike |
|---------------------------|----------------------|-------------|--------------|--------|--------------|
| Trichloroethene | 316.2 | 20 | 110.5 | 103 | 79 |
| 1,2-Dichloropropane | 193.4 | 20 | 0 | 96.7 | 76 |
| Bromodichloromethane | 189.5 | 20 | 0 | 94.8 | 69 |
| Dibromomethane | 190.1 | 20 | 0 | 95 | 76 |
| 4-Methyl-2-pentanone | 146.8 | 100 | 0 | 73.4 | 53 |
| cis-1,3-Dichloropropene | 176.5 | 10 | 0 | 88.2 | 70 |
| Toluene | 205.1 | 20 | 0 | 103 | 82 |
| trans-1,3-Dichloropropene | 172.4 | 10 | 0 | 86.2 | 64 |
| 1,1,2-Trichloroethane | 184.4 | 20 | 0 | 92.2 | 73 |
| 1,2-Dibromoethane | 188.7 | 20 | 0 | 94.4 | 73 |
| 2-Hexanone | 130 | 100 | 0 | 65 | 37 |
| 1,3-Dichloropropane | 181.8 | 20 | 0 | 90.9 | 76 |
| Tetrachloroethene | 584.2 | 20 | 431.8 | 76.2 | 82 |
| Dibromochloromethane | 175.7 | 20 | 0 | 87.8 | 59 |
| Chlorobenzene | 187.5 | 20 | 0 | 93.8 | 80 |
| 1,1,1,2-Tetrachloroethane | 183.7 | 20 | 0 | 91.8 | 72 |
| Ethylbenzene | 199.4 | 20 | 0 | 99.7 | 83 |
| m,p-Xylene | 405.6 | 20 | 0 | 101 | 84 |
| o-Xylene | 188.3 | 20 | 0 | 94.2 | 83 |
| Styrene | 192.9 | 20 | 0 | 96.5 | 80 |
| Bromoform | 139.7 | 20 | 0 | 69.8 | 54 |
| Isopropylbenzene | 212.8 | 20 | 0 | 106 | 75 |
| 1,1,2,2-Tetrachloroethane | 150.8 | 20 | 0 | 75.4 | 61 |
| 1,2,3-Trichloropropane | 157.1 | 20 | 0 | 78.6 | 66 |
| Bromobenzene | 175.6 | 20 | 0 | 87.8 | 77 |
| n-Propylbenzene | 199.4 | 20 | 0 | 99.7 | 76 |
| 2-Chlorotoluene | 185.3 | 20 | 0 | 92.6 | 78 |
| 4-Chlorotoluene | 190.8 | 20 | 0 | 95.4 | 75 |
| 1,3,5-Trimethylbenzene | 200.4 | 20 | 0 | 100 | 79 |
| tert-Butylbenzene | 203.4 | 20 | 0 | 102 | 79 |
| 1,2,4-Trimethylbenzene | 193.7 | 20 | 0 | 96.8 | 77 |

Qualifiers: ND - Not Detected at the Reporting Limit
 J - Analyte detected below quantitation limits
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

S - Spike Recovery outside accepted recovery limits
 R - RPD outside accepted recovery limits
 NA - Not applicable where J values or ND results occur

B - Analyte detected in the associated Method Blank

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT

Matrix Spike - Full List

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

| Compound | 172.8 | 195 | 173.8 | 168.8 | 159 | 171.2 | 128 | 133.7 | 142.8 | 130.6 | 120.7 | 252 | 261.6 | 255.4 | 242.9 |
|-----------------------------|-------|------|-------|-------|------|-------|------|-------|-------|-------|-------|------|-------|-------|-------|
| sec-Butylbenzene | 20 | 20 | 20 | 20 | 20 | 20 | 50 | 20 | 20 | 50 | 20 | 20 | 20 | 20 | 20 |
| 4-Isopropyltoluene | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 250 | 250 | 250 | 250 |
| 1,3-Dichlorobenzene | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L |
| 1,4-Dichlorobenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| n-Butylbenzene | 86.4 | 97.5 | 86.9 | 84.4 | 79.5 | 85.6 | 64 | 66.8 | 71.4 | 65.3 | 60.4 | 101 | 105 | 102 | 97.2 |
| 1,2-Dichlorobenzene | 128 | 128 | 122 | 123 | 130 | 121 | 127 | 128 | 134 | 131 | 131 | 116 | 127 | 114 | 117 |
| 1,2-Dibromo-3-chloropropane | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 1,2,4-Trichlorobenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Hexachlorobutadiene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Naphthalene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 1,2,3-Trichlorobenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Surr: Dibromofluoromethane | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Surr: 1,2-Dichloroethane-d4 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Surr: Toluene-d8 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Surr: 4-Bromofluorobenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |

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Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Matrix Spike Duplicate - Full List

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

Sample ID: 0702072-15Amsdf Batch ID: R35964 Test Code: SW6260B Units: µg/L Analysis Date 2/22/2007 10:39:00 PM Prep Date: 2/17/2007
Client ID: MW-209D Run ID: V-1_070222A SeqNo: 596252

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | 198.3 | 50 | µg/L | 200 | 0 | 99.2 | 16 | 150 | 179.5 | 9.95 | 20 | |
| Chloromethane | 195.2 | 50 | µg/L | 200 | 0 | 97.6 | 35 | 150 | 183.6 | 6.12 | 20 | |
| Vinyl chloride | 222 | 20 | µg/L | 200 | 0 | 111 | 49 | 150 | 210.2 | 5.46 | 20 | |
| Chloroethane | 204.5 | 50 | µg/L | 200 | 0 | 102 | 58 | 147 | 196.3 | 4.09 | 20 | |
| Bromomethane | 203.4 | 20 | µg/L | 200 | 0 | 102 | 49 | 142 | 188.6 | 7.55 | 20 | |
| Trichlorofluoromethane | 234.4 | 20 | µg/L | 200 | 0 | 117 | 57 | 149 | 215.4 | 8.45 | 20 | |
| Diethyl ether | 192.4 | 50 | µg/L | 200 | 0 | 96.2 | 66 | 136 | 176.1 | 8.85 | 20 | |
| Acetone | 161.7 | 100 | µg/L | 200 | 0 | 80.8 | 16 | 150 | 140.3 | 14.2 | 20 | |
| 1,1-Dichloroethene | 209.7 | 10 | µg/L | 200 | 0 | 105 | 70 | 150 | 184.2 | 12.9 | 20 | |
| Carbon disulfide | 219.2 | 20 | µg/L | 200 | 0 | 110 | 47 | 135 | 202.5 | 7.92 | 20 | |
| Methylene chloride | 231.7 | 50 | µg/L | 200 | 0 | 116 | 66 | 142 | 213.5 | 8.18 | 20 | |
| Methyl tert-butyl ether | 228.8 | 20 | µg/L | 200 | 0 | 114 | 63 | 138 | 210.6 | 8.28 | 20 | |
| trans-1,2-Dichloroethene | 233.8 | 20 | µg/L | 200 | 0 | 117 | 78 | 135 | 227.8 | 2.6 | 20 | |
| 1,1-Dichloroethane | 239.2 | 20 | µg/L | 200 | 0 | 120 | 76 | 131 | 221.6 | 7.64 | 20 | |
| 2-Butanone | 145.3 | 100 | µg/L | 200 | 0 | 72.7 | 51 | 142 | 123.5 | 16.2 | 20 | |
| 2,2-Dichloropropane | 226.6 | 20 | µg/L | 200 | 0 | 113 | 60 | 149 | 217.9 | 3.91 | 20 | |
| cis-1,2-Dichloroethene | 203.8 | 20 | µg/L | 200 | 0 | 102 | 74 | 128 | 196.5 | 3.65 | 20 | |
| Chloroform | 208.9 | 20 | µg/L | 200 | 0 | 104 | 80 | 129 | 194.4 | 7.19 | 20 | |
| Tetrahydrofuran | 156.7 | 100 | µg/L | 200 | 0 | 78.4 | 53 | 145 | 144.7 | 7.96 | 20 | |
| Bromochloromethane | 190 | 20 | µg/L | 200 | 0 | 95 | 78 | 130 | 177.2 | 6.97 | 20 | |
| 1,1,1-Trichloroethane | 234.6 | 20 | µg/L | 200 | 0 | 117 | 77 | 139 | 222.1 | 5.47 | 20 | |
| 1,1-Dichloropropene | 192 | 20 | µg/L | 200 | 0 | 96 | 74 | 127 | 185.4 | 3.5 | 20 | |
| Carbon tetrachloride | 190 | 20 | µg/L | 200 | 0 | 95 | 73 | 138 | 184.6 | 2.88 | 20 | |
| 1,2-Dichloroethane | 218.7 | 20 | µg/L | 200 | 0 | 109 | 75 | 130 | 205.2 | 6.37 | 20 | |
| Benzene | 214.7 | 10 | µg/L | 200 | 0 | 107 | 79 | 123 | 202.1 | 6.05 | 20 | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Matrix Spike Duplicate - Full List

CLIENT: SHAW E & I, Inc.
Work Order: 0702072
Project: 101960 Textron Gorham

| | | | | | | | | | | | |
|---------------------------|-------|-----|------|-----|-------|------|----|-----|-------|-------|----|
| Trichloroethene | 331.8 | 20 | µg/L | 200 | 110.5 | 111 | 79 | 126 | 316.2 | 4.81 | 20 |
| 1,2-Dichloropropane | 216.5 | 20 | µg/L | 200 | 0 | 108 | 76 | 125 | 193.4 | 11.3 | 20 |
| Bromodichloromethane | 201.7 | 20 | µg/L | 200 | 0 | 101 | 69 | 119 | 189.5 | 6.24 | 20 |
| Dibromomethane | 191.7 | 20 | µg/L | 200 | 0 | 95.8 | 76 | 127 | 190.1 | 0.838 | 20 |
| 4-Methyl-2-pentanone | 154.6 | 100 | µg/L | 200 | 0 | 77.3 | 53 | 141 | 146.8 | 5.18 | 20 |
| cis-1,3-Dichloropropene | 189.6 | 10 | µg/L | 200 | 0 | 94.8 | 70 | 119 | 176.5 | 7.16 | 20 |
| Toluene | 217.9 | 20 | µg/L | 200 | 0 | 109 | 82 | 124 | 205.1 | 6.05 | 20 |
| trans-1,3-Dichloropropene | 185.3 | 10 | µg/L | 200 | 0 | 92.6 | 64 | 124 | 172.4 | 7.21 | 20 |
| 1,1,2-Trichloroethane | 197.5 | 20 | µg/L | 200 | 0 | 98.8 | 73 | 127 | 184.4 | 6.86 | 20 |
| 1,2-Dibromoethane | 198.4 | 20 | µg/L | 200 | 0 | 99.2 | 73 | 127 | 188.7 | 5.01 | 20 |
| 2-Hexanone | 140.8 | 100 | µg/L | 200 | 0 | 70.4 | 37 | 145 | 130 | 7.98 | 20 |
| 1,3-Dichloropropane | 194.1 | 20 | µg/L | 200 | 0 | 97 | 76 | 123 | 181.8 | 6.54 | 20 |
| Tetrachloroethene | 627.4 | 20 | µg/L | 200 | 431.8 | 97.8 | 82 | 129 | 584.2 | 7.13 | 20 |
| Dibromochloromethane | 181.7 | 20 | µg/L | 200 | 0 | 90.8 | 59 | 125 | 175.7 | 3.36 | 20 |
| Chlorobenzene | 195.1 | 20 | µg/L | 200 | 0 | 97.6 | 80 | 120 | 187.5 | 3.97 | 20 |
| 1,1,1,2-Tetrachloroethane | 192 | 20 | µg/L | 200 | 0 | 96 | 72 | 124 | 183.7 | 4.42 | 20 |
| Ethylbenzene | 215.1 | 20 | µg/L | 200 | 0 | 108 | 83 | 123 | 199.4 | 7.58 | 20 |
| m,p-Xylene | 431.4 | 20 | µg/L | 400 | 0 | 108 | 84 | 121 | 405.6 | 6.16 | 20 |
| o-Xylene | 200.3 | 20 | µg/L | 200 | 0 | 100 | 83 | 119 | 188.3 | 6.18 | 20 |
| Styrene | 207.6 | 20 | µg/L | 200 | 0 | 104 | 80 | 122 | 192.9 | 7.34 | 20 |
| Bromoform | 156.7 | 20 | µg/L | 200 | 0 | 78.4 | 54 | 119 | 139.7 | 11.5 | 20 |
| Isopropylbenzene | 235 | 20 | µg/L | 200 | 0 | 118 | 75 | 131 | 212.8 | 9.92 | 20 |
| 1,1,2,2-Tetrachloroethane | 168.7 | 20 | µg/L | 200 | 0 | 84.4 | 61 | 139 | 150.8 | 11.2 | 20 |
| 1,2,3-Trichloropropane | 175.7 | 20 | µg/L | 200 | 0 | 87.8 | 66 | 130 | 157.1 | 11.2 | 20 |
| Bromobenzene | 193.7 | 20 | µg/L | 200 | 0 | 96.8 | 77 | 124 | 175.6 | 9.8 | 20 |
| n-Propylbenzene | 220 | 20 | µg/L | 200 | 0 | 110 | 76 | 131 | 199.4 | 9.82 | 20 |
| 2-Chlorotoluene | 207.5 | 20 | µg/L | 200 | 0 | 104 | 78 | 125 | 185.3 | 11.3 | 20 |
| 4-Chlorotoluene | 206.4 | 20 | µg/L | 200 | 0 | 103 | 75 | 124 | 190.8 | 7.85 | 20 |
| 1,3,5-Trimethylbenzene | 216 | 20 | µg/L | 200 | 0 | 108 | 79 | 124 | 200.4 | 7.49 | 20 |
| tert-Butylbenzene | 222.7 | 20 | µg/L | 200 | 0 | 111 | 79 | 126 | 203.4 | 9.06 | 20 |
| 1,2,4-Trimethylbenzene | 207.7 | 20 | µg/L | 200 | 0 | 104 | 77 | 124 | 193.7 | 6.98 | 20 |

Qualifiers: ND - Not Detected at the Reporting Limit
J - Analyte detected below quantitation limits
RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

S - Spike Recovery outside accepted recovery limits
R - RPD outside accepted recovery limits
NA - Not applicable where J values or ND results occur

B - Analyte detected in the associated Method Blank

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike Duplicate - Full List

| Compound | Concentration (µg/L) | Volume (L) | Recovery (%) | Matrix Spike | Matrix Duplicate | Matrix Spike % Diff | Matrix Duplicate % Diff | Matrix Spike Conc (µg/L) | Matrix Duplicate Conc (µg/L) | Matrix Spike Conc (µg/L) | Matrix Duplicate Conc (µg/L) | Matrix Spike Conc (µg/L) | Matrix Duplicate Conc (µg/L) |
|-----------------------------|----------------------|------------|--------------|--------------|------------------|---------------------|-------------------------|--------------------------|------------------------------|--------------------------|------------------------------|--------------------------|------------------------------|
| sec-Butylbenzene | 188.4 | 20 | 0 | 94.2 | 82 | 128 | 172.8 | 8.64 | 20 | 200 | 200 | 200 | 200 |
| 4-Isopropyltoluene | 214.6 | 20 | 0 | 107 | 77 | 128 | 195 | 9.57 | 20 | 200 | 200 | 200 | 200 |
| 1,3-Dichlorobenzene | 186.2 | 20 | 0 | 93.1 | 80 | 122 | 173.8 | 6.89 | 20 | 200 | 200 | 200 | 200 |
| 1,4-Dichlorobenzene | 181 | 20 | 0 | 90.5 | 78 | 123 | 168.8 | 6.98 | 20 | 200 | 200 | 200 | 200 |
| n-Butylbenzene | 172.2 | 20 | 0 | 86.1 | 74 | 130 | 159 | 7.97 | 20 | 200 | 200 | 200 | 200 |
| 1,2-Dichlorobenzene | 182.8 | 20 | 0 | 91.4 | 78 | 121 | 171.2 | 6.55 | 20 | 200 | 200 | 200 | 200 |
| 1,2-Dibromo-3-chloropropane | 139.1 | 50 | 0 | 69.6 | 50 | 127 | 128 | 8.31 | 20 | 200 | 200 | 200 | 200 |
| 1,2,4-Trichlorobenzene | 162.5 | 20 | 0 | 81.2 | 67 | 128 | 133.7 | 19.4 | 20 | 200 | 200 | 200 | 200 |
| Hexachlorobutadiene | 166 | 20 | 0 | 83 | 74 | 134 | 142.8 | 15 | 20 | 200 | 200 | 200 | 200 |
| Naphthalene | 158.6 | 50 | 0 | 79.3 | 57 | 131 | 130.6 | 19.4 | 20 | 200 | 200 | 200 | 200 |
| 1,2,3-Trichlorobenzene | 147 | 20 | 0 | 73.5 | 64 | 131 | 120.7 | 19.6 | 20 | 200 | 200 | 200 | 200 |
| Surr: Dibromofluoromethane | 249.3 | 20 | 0 | 99.7 | 85 | 116 | 0 | 0 | 0 | 250 | 250 | 250 | 250 |
| Surr: 1,2-Dichloroethane-d4 | 271 | 20 | 0 | 108 | 77 | 127 | 0 | 0 | 0 | 250 | 250 | 250 | 250 |
| Surr: Toluene-d8 | 253.8 | 20 | 0 | 102 | 86 | 114 | 0 | 0 | 0 | 250 | 250 | 250 | 250 |
| Surr: 4-Bromofluorobenzene | 249.1 | 20 | 0 | 99.6 | 79 | 117 | 0 | 0 | 0 | 250 | 250 | 250 | 250 |

Qualifiers: ND - Not Detected at the Reporting Limit

J - Analyte detected below quantitation limits

RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

S - Spike Recovery outside accepted recovery limits

R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank

NA - Not applicable where J values or ND results occur

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT

Matrix Spike - Full List

CLIENT: SHAW E & I, Inc.
 Work Order: 0702072
 Project: 101960 Textron Gorham

Sample ID: 0702072-19Amsf Batch ID: R35993 Test Code: SW8260B Units: µg/L Analysis Date 2/23/2007 6:20:00 PM Prep Date: 2/17/2007
 Client ID: CW-1 Run ID: V-1_070223A SeqNo: 596559

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|--------------------------|------------------|-------|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | 1581 | 500 | µg/L | 2000 | 0 | 79 | 16 | 150 | 0 | | | |
| Chloromethane | 1769 | 500 | µg/L | 2000 | 0 | 88.4 | 35 | 150 | 0 | | | |
| Vinyl chloride | 1983 | 200 | µg/L | 2000 | 0 | 99.2 | 49 | 150 | 0 | | | |
| Chloroethane | 1976 | 500 | µg/L | 2000 | 0 | 98.8 | 58 | 147 | 0 | | | |
| Bromomethane | 1922 | 200 | µg/L | 2000 | 0 | 96.1 | 49 | 142 | 0 | | | |
| Trichlorofluoromethane | 2221 | 200 | µg/L | 2000 | 0 | 111 | 57 | 149 | 0 | | | |
| Diethyl ether | 1840 | 500 | µg/L | 2000 | 0 | 92 | 66 | 136 | 0 | | | |
| Acetone | 1338 | 1,000 | µg/L | 2000 | 0 | 66.9 | 16 | 150 | 0 | | | |
| 1,1-Dichloroethene | 1983 | 100 | µg/L | 2000 | 177 | 90.3 | 70 | 150 | 0 | | | |
| Carbon disulfide | 1891 | 200 | µg/L | 2000 | 0 | 94.6 | 47 | 135 | 0 | | | |
| Methylene chloride | 2150 | 500 | µg/L | 2000 | 0 | 108 | 66 | 142 | 0 | | | |
| Methyl tert-butyl ether | 2045 | 200 | µg/L | 2000 | 0 | 102 | 63 | 138 | 0 | | | |
| trans-1,2-Dichloroethene | 2206 | 200 | µg/L | 2000 | 0 | 110 | 78 | 135 | 0 | | | |
| 1,1-Dichloroethane | 2235 | 200 | µg/L | 2000 | 0 | 112 | 76 | 131 | 0 | | | |
| 2-Butanone | 1331 | 1,000 | µg/L | 2000 | 0 | 66.6 | 51 | 142 | 0 | | | |
| 2,2-Dichloropropane | 2473 | 200 | µg/L | 2000 | 0 | 124 | 60 | 149 | 0 | | | |
| cis-1,2-Dichloroethene | 2230 | 200 | µg/L | 2000 | 352 | 93.9 | 74 | 128 | 0 | | | |
| Chloroform | 2051 | 200 | µg/L | 2000 | 0 | 103 | 80 | 129 | 0 | | | |
| Tetrahydrofuran | 1451 | 1,000 | µg/L | 2000 | 0 | 72.6 | 53 | 145 | 0 | | | |
| Bromochloromethane | 1803 | 200 | µg/L | 2000 | 0 | 90.2 | 78 | 130 | 0 | | | |
| 1,1,1-Trichloroethane | 2216 | 200 | µg/L | 2000 | 0 | 111 | 77 | 139 | 0 | | | |
| 1,1-Dichloropropene | 1863 | 200 | µg/L | 2000 | 0 | 93.2 | 74 | 127 | 0 | | | |
| Carbon tetrachloride | 1901 | 200 | µg/L | 2000 | 0 | 95 | 73 | 138 | 0 | | | |
| 1,2-Dichloroethane | 2109 | 200 | µg/L | 2000 | 0 | 105 | 75 | 130 | 0 | | | |
| Benzene | 1961 | 100 | µg/L | 2000 | 0 | 98 | 79 | 123 | 0 | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike - Full List

| Compound | Reporting Limit | Concentration | Recovery | Acceptance | Recovery | Acceptance | Recovery | Acceptance | Recovery | Acceptance |
|---------------------------|-----------------|---------------|----------|------------|----------|------------|----------|------------|----------|------------|
| Trichloroethene | 7887 | 200 | µg/L | 2000 | 5939 | 97.4 | 79 | 126 | 0 | |
| 1,2-Dichloropropane | 1986 | 200 | µg/L | 2000 | 0 | 99.3 | 76 | 125 | 0 | |
| Bromodichloromethane | 1878 | 200 | µg/L | 2000 | 0 | 93.9 | 69 | 119 | 0 | |
| Dibromomethane | 1849 | 200 | µg/L | 2000 | 0 | 92.5 | 76 | 127 | 0 | |
| 4-Methyl-2-pentanone | 1296 | 1,000 | µg/L | 2000 | 0 | 64.8 | 53 | 141 | 0 | |
| cis-1,3-Dichloropropene | 1856 | 100 | µg/L | 2000 | 0 | 92.8 | 70 | 119 | 0 | |
| Toluene | 2050 | 200 | µg/L | 2000 | 0 | 103 | 82 | 124 | 0 | |
| trans-1,3-Dichloropropene | 1805 | 100 | µg/L | 2000 | 0 | 90.2 | 64 | 124 | 0 | |
| 1,1,2-Trichloroethane | 1905 | 200 | µg/L | 2000 | 0 | 95.2 | 73 | 127 | 0 | |
| 1,2-Dibromoethane | 1922 | 200 | µg/L | 2000 | 0 | 96.1 | 73 | 127 | 0 | |
| 2-Hexanone | 1311 | 1,000 | µg/L | 2000 | 0 | 65.6 | 37 | 145 | 0 | |
| 1,3-Dichloropropane | 1834 | 200 | µg/L | 2000 | 0 | 91.7 | 76 | 123 | 0 | |
| Tetrachloroethene | 2144 | 200 | µg/L | 2000 | 0 | 107 | 82 | 129 | 0 | |
| Dibromochloromethane | 1767 | 200 | µg/L | 2000 | 0 | 88.4 | 59 | 125 | 0 | |
| Chlorobenzene | 1835 | 200 | µg/L | 2000 | 0 | 91.8 | 80 | 120 | 0 | |
| 1,1,1,2-Tetrachloroethane | 1782 | 200 | µg/L | 2000 | 0 | 89.1 | 72 | 124 | 0 | |
| Ethylbenzene | 1944 | 200 | µg/L | 2000 | 0 | 97.2 | 83 | 123 | 0 | |
| m,p-Xylene | 3889 | 200 | µg/L | 4000 | 0 | 97.2 | 84 | 121 | 0 | |
| o-Xylene | 1877 | 200 | µg/L | 2000 | 0 | 93.8 | 83 | 119 | 0 | |
| Styrene | 1976 | 200 | µg/L | 2000 | 0 | 98.8 | 80 | 122 | 0 | |
| Bromoform | 1496 | 200 | µg/L | 2000 | 0 | 74.8 | 54 | 119 | 0 | |
| Isopropylbenzene | 2081 | 200 | µg/L | 2000 | 0 | 104 | 75 | 131 | 0 | |
| 1,1,2,2-Tetrachloroethane | 1564 | 200 | µg/L | 2000 | 0 | 78.2 | 61 | 139 | 0 | |
| 1,2,3-Trichloropropane | 1645 | 200 | µg/L | 2000 | 0 | 82.2 | 66 | 130 | 0 | |
| Bromobenzene | 1799 | 200 | µg/L | 2000 | 0 | 90 | 77 | 124 | 0 | |
| n-Propylbenzene | 1959 | 200 | µg/L | 2000 | 0 | 98 | 76 | 131 | 0 | |
| 2-Chlorotoluene | 1911 | 200 | µg/L | 2000 | 0 | 95.6 | 78 | 125 | 0 | |
| 4-Chlorotoluene | 1934 | 200 | µg/L | 2000 | 0 | 96.7 | 75 | 124 | 0 | |
| 1,3,5-Trimethylbenzene | 1940 | 200 | µg/L | 2000 | 0 | 97 | 79 | 124 | 0 | |
| tert-Butylbenzene | 2002 | 200 | µg/L | 2000 | 0 | 100 | 79 | 126 | 0 | |
| 1,2,4-Trimethylbenzene | 1947 | 200 | µg/L | 2000 | 0 | 97.4 | 77 | 124 | 0 | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
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AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike - Full List

| Compound | 1694 | 1920 | 1802 | 1781 | 1563 | 1749 | 1116 | 1397 | 1388 | 1305 | 1270 | 2548 | 2635 | 2563 | 2493 |
|-----------------------------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|
| sec-Butylbenzene | 200 | 200 | 200 | 200 | 200 | 200 | 500 | 200 | 200 | 500 | 200 | 200 | 200 | 200 | 200 |
| 4-Isopropyltoluene | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L |
| 1,3-Dichlorobenzene | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2500 | 2500 | 2500 | 2500 |
| 1,4-Dichlorobenzene | 82 | 77 | 80 | 78 | 74 | 78 | 50 | 67 | 74 | 57 | 64 | 85 | 77 | 86 | 79 |
| n-Butylbenzene | 0 | 96 | 90.1 | 89.1 | 78.2 | 87.5 | 55.8 | 69.8 | 69.4 | 65.2 | 63.5 | 102 | 105 | 103 | 99.7 |
| 1,2-Dichlorobenzene | 128 | 128 | 122 | 123 | 130 | 121 | 127 | 128 | 134 | 131 | 131 | 116 | 127 | 114 | 117 |
| 1,2-Dibromo-3-chloropropane | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 1,2,4-Trichlorobenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Hexachlorobutadiene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Naphthalene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 1,2,3-Trichlorobenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Surr: Dibromofluoromethane | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Surr: 1,2-Dichloroethane-d4 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Surr: Toluene-d8 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Surr: 4-Bromofluorobenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |

S

S

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike Duplicate - Full List

| Sample ID: 0702072-19Amsdf | Batch ID: R35993 | Test Code: SW8260B | Units: µg/L | Analysis Date 2/23/2007 6:54:00 PM | Prep Date: 2/17/2007 | | | | | | |
|----------------------------|---------------------|--------------------|-------------|------------------------------------|------------------------|-----------|----------|------|-------|----------|-----|
| Client ID: CW-1 | Run ID: V-1_070223A | SeqNo: 596561 | | | | | | | | | |
| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | HighLimit | LowLimit | %REC | %RPD | RPDLimit | Que |
| Dichlorodifluoromethane | 1697 | 500 | µg/L | 2000 | 0 | 150 | 16 | 84.8 | 7.08 | 20 | |
| Chloromethane | 1805 | 500 | µg/L | 2000 | 0 | 150 | 35 | 90.2 | 2.01 | 20 | |
| Vinyl chloride | 2084 | 200 | µg/L | 2000 | 0 | 150 | 49 | 104 | 4.97 | 20 | |
| Chloroethane | 1962 | 500 | µg/L | 2000 | 0 | 147 | 58 | 98.1 | 0.711 | 20 | |
| Bromomethane | 1909 | 200 | µg/L | 2000 | 0 | 142 | 49 | 95.4 | 0.679 | 20 | |
| Trichlorofluoromethane | 2260 | 200 | µg/L | 2000 | 0 | 149 | 57 | 113 | 1.74 | 20 | |
| Diethyl ether | 1883 | 500 | µg/L | 2000 | 0 | 136 | 66 | 94.2 | 2.31 | 20 | |
| Acetone | 1457 | 1,000 | µg/L | 2000 | 0 | 150 | 16 | 72.8 | 8.52 | 20 | |
| 1,1-Dichloroethene | 2145 | 100 | µg/L | 2000 | 177 | 150 | 70 | 98.4 | 7.85 | 20 | |
| Carbon disulfide | 2058 | 200 | µg/L | 2000 | 0 | 135 | 47 | 103 | 8.46 | 20 | |
| Methylene chloride | 2222 | 500 | µg/L | 2000 | 0 | 142 | 66 | 111 | 3.29 | 20 | |
| Methyl tert-butyl ether | 2128 | 200 | µg/L | 2000 | 0 | 138 | 63 | 106 | 3.98 | 20 | |
| trans-1,2-Dichloroethene | 2336 | 200 | µg/L | 2000 | 0 | 135 | 78 | 117 | 5.72 | 20 | |
| 1,1-Dichloroethane | 2437 | 200 | µg/L | 2000 | 0 | 131 | 76 | 122 | 8.65 | 20 | |
| 2-Butanone | 1456 | 1,000 | µg/L | 2000 | 0 | 142 | 51 | 72.8 | 8.97 | 20 | |
| 2,2-Dichloropropane | 2564 | 200 | µg/L | 2000 | 0 | 149 | 60 | 128 | 3.61 | 20 | |
| cis-1,2-Dichloroethene | 2418 | 200 | µg/L | 2000 | 352 | 128 | 74 | 103 | 8.09 | 20 | |
| Chloroform | 2097 | 200 | µg/L | 2000 | 0 | 129 | 80 | 105 | 2.22 | 20 | |
| Tetrahydrofuran | 1545 | 1,000 | µg/L | 2000 | 0 | 145 | 53 | 77.2 | 6.28 | 20 | |
| Bromochloromethane | 1974 | 200 | µg/L | 2000 | 0 | 130 | 78 | 98.7 | 9.05 | 20 | |
| 1,1,1-Trichloroethane | 2404 | 200 | µg/L | 2000 | 0 | 139 | 77 | 120 | 8.14 | 20 | |
| 1,1-Dichloropropene | 2026 | 200 | µg/L | 2000 | 0 | 127 | 74 | 101 | 8.38 | 20 | |
| Carbon tetrachloride | 2046 | 200 | µg/L | 2000 | 0 | 138 | 73 | 102 | 7.35 | 20 | |
| 1,2-Dichloroethane | 2155 | 200 | µg/L | 2000 | 0 | 130 | 75 | 108 | 2.16 | 20 | |
| Benzene | 2135 | 100 | µg/L | 2000 | 0 | 123 | 79 | 107 | 8.5 | 20 | |

Qualifiers: ND - Not Detected at the Reporting Limit
 S - Spike Recovery outside accepted recovery limits
 B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits
 R - RPD outside accepted recovery limits
 NA - Not applicable where J values or ND results occur
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AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike Duplicate - Full List

| Compound | 8359 | 2115 | 2039 | 1926 | 1428 | 1954 | 2172 | 1947 | 1965 | 1874 | 1327 | 1892 | 2137 | 1761 | 1941 | 1885 | 2089 | 4285 | 1965 | 2049 | 1495 | 2258 | 1580 | 1638 | 1842 | 2123 | 1984 | 2033 | 2116 | 2140 | 2050 | | | |
|---------------------------|------|------|------|------|-------|------|------|------|------|------|-------|------|-------|------|------|------|------|------|------|------|--------|------|------|-------|------|------|------|------|------|------|------|--|--|--|
| Trichloroethene | 200 | 200 | 200 | 200 | 1,000 | 100 | 200 | 100 | 200 | 200 | 1,000 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 | | | |
| 1,2-Dichloropropane | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | | | |
| Bromodichloromethane | 5939 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | | | |
| Dibromomethane | 121 | 106 | 102 | 96.3 | 71.4 | 97.7 | 109 | 97.4 | 98.2 | 93.7 | 66.4 | 94.6 | 107 | 88 | 97 | 94.2 | 104 | 107 | 98.2 | 102 | 74.8 | 113 | 79 | 81.9 | 92.1 | 106 | 99.2 | 102 | 106 | 107 | 103 | | | |
| 4-Methyl-2-pentanone | 79 | 76 | 69 | 76 | 53 | 70 | 82 | 64 | 73 | 73 | 37 | 76 | 82 | 59 | 80 | 72 | 83 | 84 | 83 | 80 | 54 | 75 | 61 | 66 | 77 | 76 | 78 | 75 | 79 | 79 | 77 | | | |
| cis-1,3-Dichloropropene | 126 | 125 | 119 | 127 | 141 | 119 | 124 | 124 | 127 | 127 | 145 | 123 | 129 | 125 | 120 | 124 | 123 | 121 | 119 | 122 | 119 | 131 | 139 | 130 | 124 | 131 | 125 | 124 | 124 | 126 | 124 | | | |
| Toluene | 7887 | 5.81 | 1986 | 1878 | 1849 | 1296 | 1856 | 2050 | 1805 | 1905 | 1311 | 1834 | 2144 | 1767 | 1835 | 1782 | 1944 | 3889 | 1877 | 1976 | 1496 | 2081 | 1564 | 1645 | 1799 | 1959 | 1911 | 1934 | 1940 | 2002 | 1947 | | | |
| trans-1,3-Dichloropropene | 5.81 | 6.29 | 8.22 | 4.08 | 9.69 | 5.14 | 5.78 | 7.57 | 3.1 | 2.53 | 1.21 | 3.11 | 0.327 | 0.34 | 5.61 | 5.62 | 7.19 | 9.69 | 4.58 | 3.63 | 0.0669 | 8.16 | 1.02 | 0.426 | 2.36 | 8.04 | 3.75 | 4.99 | 8.68 | 6.66 | 5.15 | | | |
| 1,1,2-Trichloroethane | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 1,2-Dibromoethane | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 2-Hexanone | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 1,3-Dichloropropane | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Tetrachloroethene | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Dibromochloromethane | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Chlorobenzene | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 1,1,1,2-Tetrachloroethane | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Ethylbenzene | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| m,p-Xylene | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| o-Xylene | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Styrene | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Bromoform | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Isopropylbenzene | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 1,1,2,2-Tetrachloroethane | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 1,2,3-Trichloropropane | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Bromobenzene | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| n-Propylbenzene | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 2-Chlorotoluene | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 4-Chlorotoluene | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 1,3,5-Trimethylbenzene | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| tert-Butylbenzene | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| 1,2,4-Trimethylbenzene | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Matrix Spike Duplicate - Full List

| Compound | 1827 | 2041 | 1812 | 1791 | 1707 | 1818 | 1385 | 1521 | 1641 | 1416 | 1410 | 2560 | 2632 | 2568 | 2451 |
|-----------------------------|------|------|-------|------|------|------|------|------|------|------|------|------|------|------|------|
| sec-Butylbenzene | 200 | 200 | 200 | 200 | 200 | 200 | 500 | 200 | 200 | 500 | 200 | 200 | 200 | 200 | 200 |
| 4-Isopropyltoluene | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2000 | 2500 | 2500 | 2500 | 2500 |
| 1,3-Dichlorobenzene | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L |
| 1,4-Dichlorobenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| n-Butylbenzene | 91.4 | 102 | 90.6 | 89.6 | 85.4 | 90.9 | 69.2 | 76 | 82 | 70.8 | 70.5 | 102 | 105 | 103 | 98 |
| 1,2-Dichlorobenzene | 82 | 77 | 80 | 78 | 74 | 78 | 50 | 67 | 74 | 57 | 64 | 85 | 77 | 86 | 79 |
| 1,2-Dibromo-3-chloropropane | 128 | 128 | 122 | 123 | 130 | 121 | 127 | 128 | 134 | 131 | 131 | 116 | 127 | 114 | 117 |
| 1,2,4-Trichlorobenzene | 1694 | 1920 | 1802 | 1781 | 1563 | 1749 | 1116 | 1397 | 1388 | 1305 | 1270 | 0 | 0 | 0 | 0 |
| Hexachlorobutadiene | 7.55 | 6.11 | 0.553 | 0.56 | 8.81 | 3.87 | 21.5 | 8.5 | 16.7 | 8.16 | 10.4 | 0 | 0 | 0 | 0 |
| Naphthalene | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 0 | 0 | 0 | 0 |
| 1,2,3-Trichlorobenzene | R | | | | | | | | | | | | | | |
| Surr: Dibromofluoromethane | | | | | | | | | | | | | | | |
| Surr: 1,2-Dichloroethane-d4 | | | | | | | | | | | | | | | |
| Surr: Toluene-d8 | | | | | | | | | | | | | | | |
| Surr: 4-Bromofluorobenzene | | | | | | | | | | | | | | | |

Qualifiers: ND - Not Detected at the Reporting Limit
 J - Analyte detected below quantitation limits
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

S - Spike Recovery outside accepted recovery limits
 R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank
 NA - Not applicable where J values or ND results occur

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

| | | | |
|-------------------|-----------------------|--------------------------|--------------------|
| CLIENT: | SHAW E & I, Inc. | Client Sample ID: | CW-6 |
| Lab Order: | 0702072 | Tag Number: | |
| Project: | 101960 Textron Gorham | Collection Date: | 2/17/07 1:45:00 PM |
| Lab ID: | 0702072-23A | Matrix: | GROUNDWATER |

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|---------------------------------------|--------|----------------|------|-------|----|--------------------|
| TPH BY GC/FID (MODIFIED 8015B) | | SW8015B | | | | Analyst: FQ |
| Gasoline | ND | 0.051 | | mg/L | 1 | 2/23/07 6:08:00 PM |
| Mineral Spirits | ND | 0.051 | | mg/L | 1 | 2/23/07 6:08:00 PM |
| Kerosene | ND | 0.051 | | mg/L | 1 | 2/23/07 6:08:00 PM |
| Diesel Fuel/Fuel Oil #2 | ND | 0.051 | | mg/L | 1 | 2/23/07 6:08:00 PM |
| Motor Oil/Hydraulic Oil | ND | 0.10 | | mg/L | 1 | 2/23/07 6:08:00 PM |
| Unidentified Hydrocarbons | 8.5 | 0.20 | | mg/L | 1 | 2/23/07 6:08:00 PM |
| Surr: o-Terphenyl | 63.5 | 31-131 | | %REC | 1 | 2/23/07 6:08:00 PM |

Gasoline cannot be accurately determined by this method. Purge and trap sample introduction into a GC or GCMS is the recommended approach for gasoline. Due to the physical, chemical, and biological processes which affect the chemical composition of fuel mixtures exposed to the environment, the qualitative identity of a hydrocarbon mixture as a fuel product is not always conclusive by this method due to the method's reliance on chromatographic pattern recognition. A result provided for a specific fuel indicates that the mixture present in the sample has a chromatographic pattern similar to the laboratory's reference standard for that fuel mixture under specific GC operating conditions utilized at the time of analysis. A result identified as Unidentified Hydrocarbons is based upon the detector response obtained for the laboratory's Fuel Oil#2 reference standard and includes the entire chromatographic response for the sample between n-Alkanes of carbon numbers C9 to C36.

Qualifiers:

| | |
|---|---|
| ND - Not Detected at the Reporting Limit | S - Spike Recovery outside accepted recovery limits |
| J - Analyte detected below quantitation limits | R - RPD outside accepted recovery limits |
| B - Analyte detected in the associated Method Blank | E - Value above quantitation range |
| H - Method prescribed holding time exceeded. | # - See Case Narrative |
| RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate. | |

AMRO Environmental Laboratories Corp.

Date: 01-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0702072
Project: 101960 Textron Gorham
Lab ID: 0702072-24A

Client Sample ID: CW-6 DUP
Tag Number:
Collection Date: 2/17/07 1:45:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|---------------------------------------|--------|----------------|------|-------|----|--------------------|
| TPH BY GC/FID (MODIFIED 8015B) | | SW8015B | | | | Analyst: FQ |
| Gasoline | ND | 0.052 | | mg/L | 1 | 2/23/07 6:44:00 PM |
| Mineral Spirits | ND | 0.052 | | mg/L | 1 | 2/23/07 6:44:00 PM |
| Kerosene | ND | 0.052 | | mg/L | 1 | 2/23/07 6:44:00 PM |
| Diesel Fuel/Fuel Oil #2 | ND | 0.052 | | mg/L | 1 | 2/23/07 6:44:00 PM |
| Motor Oil/Hydraulic Oil | ND | 0.10 | | mg/L | 1 | 2/23/07 6:44:00 PM |
| Unidentified Hydrocarbons | 8.6 | 0.21 | | mg/L | 1 | 2/23/07 6:44:00 PM |
| Surr: o-Terphenyl | 66.3 | 31-131 | | %REC | 1 | 2/23/07 6:44:00 PM |

Gasoline cannot be accurately determined by this method. Purge and trap sample introduction into a GC or GCMS is the recommended approach for gasoline. Due to the physical, chemical, and biological processes which affect the chemical composition of fuel mixtures exposed to the environment, the qualitative identity of a hydrocarbon mixture as a fuel product is not always conclusive by this method due to the method's reliance on chromatographic pattern recognition. A result provided for a specific fuel indicates that the mixture present in the sample has a chromatographic pattern similar to the laboratory's reference standard for that fuel mixture under specific GC operating conditions utilized at the time of analysis. A result identified as Unidentified Hydrocarbons is based upon the detector response obtained for the laboratory's Fuel Oil#2 reference standard and includes the entire chromatographic response for the sample between n-Alkanes of carbon numbers C9 to C36.

- Qualifiers:**
- ND - Not Detected at the Reporting Limit
 - J - Analyte detected below quantitation limits
 - B - Analyte detected in the associated Method Blank
 - H - Method prescribed holding time exceeded.
 - RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.
 - S - Spike Recovery outside accepted recovery limits
 - R - RPD outside accepted recovery limits
 - E - Value above quantitation range
 - # - See Case Narrative

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

QC SUMMARY REPORT
Method Blank

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

Sample ID: MB-16813 Batch ID: 16813 Test Code: SW8015B Units: mg/L Analysis Date 2/23/2007 4:18:00 PM Prep Date: 2/23/2007
 Client ID: Run ID: GC-FING1_070223A SeqNo: 596574

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|---------------------------|------------------|-------|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Gasoline | ND | 0.050 | mg/L | | | | | | | | | |
| Mineral Spirits | ND | 0.050 | mg/L | | | | | | | | | |
| Kerosene | ND | 0.050 | mg/L | | | | | | | | | |
| Diesel Fuel/Fuel Oil #2 | ND | 0.050 | mg/L | | | | | | | | | |
| Motor Oil/Hydraulic Oil | ND | 0.10 | mg/L | | | | | | | | | |
| Unidentified Hydrocarbons | ND | 0.20 | mg/L | | | | | | | | | |
| Surr: o-Terphenyl | 0.06958 | 0 | mg/L | 0.1 | 0 | 69.6 | 31 | 131 | 0 | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 28-Feb-07

CLIENT: SHAW E & I, Inc.

Work Order: 0702072

Project: 101960 Textron Gorham

QC SUMMARY REPORT

Laboratory Control Spike

Sample ID: LCS-16813 Batch ID: 16813 Test Code: SW8015B Units: mg/L Analysis Date 2/23/2007 4:55:00 PM Prep Date: 2/23/2007
 Client ID: Run ID: GC-FING1_070223A SeqNo: 596575

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|-------------------------|------------------|-------|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Diesel Fuel/Fuel Oil #2 | 1.236 | 0.050 | mg/L | 2 | 0 | 61.8 | 42 | 119 | 0 | | | |
| Surr: o-Terphenyl | 0.07476 | 0 | mg/L | 0.1 | 0 | 74.8 | 31 | 131 | 0 | | | |

Sample ID: LCSD-16813 Batch ID: 16813 Test Code: SW8015B Units: mg/L Analysis Date 2/23/2007 5:31:00 PM Prep Date: 2/23/2007
 Client ID: Run ID: GC-FING1_070223A SeqNo: 596576

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Que |
|-------------------------|------------------|-------|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Diesel Fuel/Fuel Oil #2 | 1.184 | 0.050 | mg/L | 2 | 0 | 59.2 | 42 | 119 | 1.236 | 4.29 | 40 | |
| Surr: o-Terphenyl | 0.06378 | 0 | mg/L | 0.1 | 0 | 63.8 | 31 | 131 | 0 | 0 | 0 | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.



111 Herrick Street, Merrimack, NH 03054
TEL: (603) 424-2022 • FAX: (603) 429-8496
www.amrolabs.com

March 22, 2007

ANALYTICAL TEST RESULTS

Ed VanDoren
SHAW E & I, Inc.
11 Northeastern Boulevard
Salem, NH 030791953
TEL: (603) 870-4500
FAX: (603) 870-4501

Subject: 101960 Textron

Workorder No.: 0703003

Dear Ed VanDoren:

AMRO Environmental Laboratories Corp. received 5 samples on 3/2/07 for the analyses presented in the following report.

AMRO is accredited in accordance with NELAC and certifies that these test results meet all the requirements of NELAC, where applicable, unless otherwise noted in the case narrative.

The enclosed Sample Receipt Checklist details the condition of your sample(s) upon receipt. Please be advised that any unused sample volume and sample extracts will be stored for a period of 60 days from sample receipt date (90 days for samples from New York). After this time, AMRO will properly dispose of the remaining sample(s). If you require further analysis, or need the samples held for a longer period, please contact us immediately.

This report consists of a total of 28 pages. This letter is an integral part of your data report. All results in this project relate only to the sample(s) as received by the laboratory and documented in the Chain-of-Custody. This report shall not be reproduced except in full, without the written approval of the laboratory. If you have any questions regarding this project in the future, please refer to the Workorder Number above.

Sincerely,

Nancy Stewart
Vice President

State Certifications: NH (NELAC): 1001, MA: M-NH012, CT: PH-0758, NY: 11278 (NELAC), ME: NH012 and 1001, NJ: NH125, RI: 00105, U.S. Army Corps of Engineers (USACE), Naval Facilities Engineering Service Center (NFESC).

Hard copy of the State Certification is available upon request.



CLIENT: SHAW E & I, Inc.
Project: 101960 Textron
Lab Order: 0703003
Date Received: 3/2/07

Work Order Sample Summary

| Lab Sample ID | Client Sample ID | Collection Date | Collection Time |
|----------------------|-------------------------|------------------------|------------------------|
| 0703003-01A | MW-220S | 3/1/07 | 11:00 AM |
| 0703003-02A | MW-221S | 3/1/07 | 11:30 AM |
| 0703003-02B | MW-221S | 3/1/07 | 11:30 AM |
| 0703003-03A | MW-109D | 3/1/07 | 1:00 PM |
| 0703003-04A | GZA-3 | 3/1/07 | 1:50 PM |
| 0703003-05A | Trip Blank | 3/1/07 | 12:00 AM |

DATES REPORT

Lab Order: 0703003
Client: SHAW E & I, Inc.
Project: 101960 Textron

| Sample ID | Client Sample ID | Collection Date | Matrix | Analytical Test Name | Prep Date | Analysis Date | TCLP Date |
|-------------|------------------|--------------------|-------------|---|-----------|------------------|-----------|
| | | | | Preparatory Test Name | | Batch ID | |
| 0703003-01A | MW-220S | 3/1/07 11:00:00 AM | Groundwater | EPA 8260B VOLATILES by GC/MS EPA 5030B | 3/1/07 | 3/6/07 R36052 | |
| 0703003-02A | MW-221S | 3/1/07 11:30:00 AM | | EPA 8260B VOLATILES by GC/MS | 3/1/07 | 3/9/07 R36118 | |
| 0703003-02B | | | | TPH by GC/FID (modified 8015B) AQPREP SEP FUNNEL: FING | 3/6/07 | 3/10/07 16854 | |
| 0703003-03A | MW-109D | 3/1/07 1:00:00 PM | | EPA 8260B VOLATILES by GC/MS EPA 5030B | 3/1/07 | 3/9/07 R36118 | |
| 0703003-04A | GZA-3 | 3/1/07 1:50:00 PM | | EPA 6010B ICP METALS, DISSOLVED EPA 3010 AQPREP TOTAL METALS: ICP/GFAA | 3/5/07 | 3/9/07 16844 | |
| 0703003-05A | Trip Blank | 3/1/07 | Trip Blank | EPA 8260B VOLATILES by GC/MS EPA 5030B | 3/1/07 | 3/6/07 R36052 | |

| | | | | | |
|--|---------------------------------|-----------------------------|--|--|-------------------------------------|
| Project No.: 101960 | Project Name: TEXTROL | Project State: RI | Project Manager: <i>[Signature]</i> | Samples (Signature): <i>[Signature]</i> | AMRO Project No.: 0703003 |
| P.O.#: | Results Needed by: | Total # of Cont. & Size | REQUESTED ANALYSES | | |
| QUOTE #: | Seal Intact? Yes No N/A | Matrix | Remarks | | |
| Sample ID: | Date/Time Sampled | Comp. Grab | Pb Sample Field Filtered 0.45 MICRON FILTER | | |
| MW-2205 | 3/1/07 @ 1100 | GW | X | | |
| MW-2215 | 3/1/07 @ 1130 | GW | X | | |
| MW-1041D | 3/1/07 @ 1300 | GW | X | | |
| GZA-3 | 3/1/07 @ 1350 | GW | X | | |
| TRIP BUBBLR | | | | | |
| Preservative: Cl-HCl, MeOH, N-HN03, S-H2SO4, Na-NaOH, O- Other | | | | | |
| Send Results To: SHAN ENV. ED VAN DUSEN 11 N. EASTERN BLVD SALUN, MA 03079 PHONE # 603 810 4530 E-mail: | | | | | |
| PRIORITY TURNAROUND TIME AUTHORIZATION Before submitting samples for expedited TAT, you must have a coded AUTHORIZATION NUMBER AUTHORIZATION No.: BY: | | | | | |
| METALS 8 RCRA <input type="checkbox"/> 13 PP <input type="checkbox"/> 23 TAL <input type="checkbox"/> 14 MCP <input type="checkbox"/> Method: 6010 <input type="checkbox"/> 200.7 <input type="checkbox"/> Other Metals: | | | | | |
| Dissolved Metals Field Filtered? YES <input type="checkbox"/> NO <input type="checkbox"/> MCP Presumptive Certainty Required? YES <input type="checkbox"/> NO <input type="checkbox"/> Received By: <i>[Signature]</i> | | | | | |
| Required Reporting Limits: S-1 <input type="checkbox"/> GW-1 <input type="checkbox"/> S-2 <input type="checkbox"/> GW-2 <input type="checkbox"/> S-3 <input type="checkbox"/> GW-3 <input type="checkbox"/> Other: | | | | | |
| AMRO policy requires notification in writing to the laboratory in cases where the samples were collected from highly contaminated sites. | | | | | |
| KNOWN SITE CONTAMINATION: | | | | | |
| SHEET 1 OF 1 AMROCC2004, Rev.3 08/18/04 | | | | | |

CLIENT: SHAW E & I, Inc.
Project: 101960 Textron
Lab Order: 0703003

CASE NARRATIVE

VOLATILES:

No QC deviations were observed.

TPH GC/FID:

No QC deviations were observed.

METALS:

No QC deviations were observed.

DATA COMMENT PAGE

Organic Data Qualifiers

| | |
|----|--|
| ND | Indicates compound was analyzed for, but not detected at or above the reporting limit. |
| J | Indicates an estimated value. This flag is used either when estimating a concentration for tentatively identified compounds where a 1:1 response is assumed, or when the data indicates the presence of a compound that meets the identification criteria but the result is less than the sample quantitation limit but greater than the method detection limit. |
| H | Method prescribed holding time exceeded. |
| E | This flag identifies compounds whose concentrations exceed the calibration range of the instrument for that specific analysis. |
| B | This flag is used when the analyte is found in the associated blank as well as in the sample. |
| R | RPD outside accepted recovery limits |
| RL | Reporting limit; defined as the lowest concentration the laboratory can accurately quantitate. |
| S | Spike Recovery outside accepted recovery limits. |
| # | See Case Narrative |

Micro Data Qualifiers

TNTC Too numerous to count

Inorganic Data Qualifiers

| | |
|---------|---|
| ND or U | Indicates element was analyzed for, but not detected at or above the reporting limit. |
| J | Indicates a value greater than or equal to the method detection limit, but less than the quantitation limit. |
| H | Indicates analytical holding time exceedance. |
| B | Indicates that the analyte is found in the associated blank, as well as in the sample. |
| MSA | Indicates value determined by the Method of Standard Addition |
| E | This flag identifies compounds whose concentrations exceed the calibration range of the instrument for that specific analysis. |
| R | RPD outside accepted recovery limits |
| RL | Reporting limit; defined as the lowest concentration the laboratory can accurately quantitate. |
| S | Spike Recovery outside accepted recovery limits. |
| W | Post-digestion spike for Furnace AA analysis is out of control limits (85-115), while sample absorbance is less than 50% of spike absorbance. |
| * | Duplicate analysis not within control limits. |
| + | Indicates the correlation coefficient for the Method of Standard Addition is less than 0.995 |
| # | See Case Narrative |

Report Comments:

1. Soil, sediment and sludge sample results are reported on a "dry weight" basis.
2. Reporting limits are adjusted for sample size used, dilutions and moisture content, if applicable.

AMRO Environmental Laboratories Corp.

Date: 15-Mar-07

CLIENT: SHAW E & I, Inc.
 Lab Order: 0703003
 Project: 101960 Textron
 Lab ID: 0703003-01A

Client Sample ID: MW-220S
 Collection Date: 3/1/07 11:00:00 AM
 Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|-------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 50 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Chloromethane | ND | 50 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Vinyl chloride | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Chloroethane | ND | 50 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Bromomethane | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Trichlorofluoromethane | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Diethyl ether | ND | 50 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Acetone | ND | 100 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,1-Dichloroethene | 11 | 10 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Carbon disulfide | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Methylene chloride | ND | 50 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Methyl tert-butyl ether | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| trans-1,2-Dichloroethene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,1-Dichloroethane | 290 | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 2-Butanone | ND | 100 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 2,2-Dichloropropane | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| cis-1,2-Dichloroethene | 240 | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Chloroform | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Tetrahydrofuran | ND | 100 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Bromochloromethane | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,1,1-Trichloroethane | 210 | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,1-Dichloropropene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Carbon tetrachloride | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,2-Dichloroethane | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Benzene | ND | 10 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Trichloroethene | 49 | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,2-Dichloropropane | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Bromodichloromethane | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Dibromomethane | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 4-Methyl-2-pentanone | ND | 100 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| cis-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Toluene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| trans-1,3-Dichloropropene | ND | 10 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,1,2-Trichloroethane | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,2-Dibromoethane | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 2-Hexanone | ND | 100 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,3-Dichloropropane | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Tetrachloroethene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Dibromochloromethane | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |

AMRO Environmental Laboratories Corp.

Date: 15-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0703003
Project: 101960 Textron
Lab ID: 0703003-01A

Client Sample ID: MW-220S
Collection Date: 3/1/07 11:00:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|-------------------|
| Chlorobenzene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Ethylbenzene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| m,p-Xylene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| o-Xylene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Styrene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Bromoform | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Isopropylbenzene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,2,3-Trichloropropane | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Bromobenzene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| n-Propylbenzene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 2-Chlorotoluene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 4-Chlorotoluene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,3,5-Trimethylbenzene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| tert-Butylbenzene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,2,4-Trimethylbenzene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| sec-Butylbenzene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 4-Isopropyltoluene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,3-Dichlorobenzene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,4-Dichlorobenzene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| n-Butylbenzene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,2-Dichlorobenzene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 50 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,2,4-Trichlorobenzene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Hexachlorobutadiene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Naphthalene | ND | 50 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| 1,2,3-Trichlorobenzene | ND | 20 | | µg/L | 10 | 3/6/07 5:00:00 PM |
| Surr: Dibromofluoromethane | 107 | 85-116 | | %REC | 10 | 3/6/07 5:00:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 110 | 77-127 | | %REC | 10 | 3/6/07 5:00:00 PM |
| Surr: Toluene-d8 | 104 | 86-114 | | %REC | 10 | 3/6/07 5:00:00 PM |
| Surr: 4-Bromofluorobenzene | 99.3 | 79-117 | | %REC | 10 | 3/6/07 5:00:00 PM |

AMRO Environmental Laboratories Corp.

Date: 15-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0703003
Project: 101960 Textron
Lab ID: 0703003-02A

Client Sample ID: MW-221S
Collection Date: 3/1/07 11:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|--------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: KT |
| Dichlorodifluoromethane | 5.1 | 5.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Vinyl chloride | 22 | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Chloroethane | 26 | 5.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Trichlorofluoromethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,1-Dichloroethene | 1.2 | 1.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,1-Dichloroethane | 140 | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| cis-1,2-Dichloroethene | 8.5 | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Chloroform | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Tetrahydrofuran | 29 | 10 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,1,1-Trichloroethane | 25 | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Benzene | 2.6 | 1.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Trichloroethene | 12 | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Bromodichloromethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Toluene | 3.1 | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Tetrachloroethene | 4.5 | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |

AMRO Environmental Laboratories Corp.

Date: 15-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0703003
Project: 101960 Textron
Lab ID: 0703003-02A

Client Sample ID: MW-221S
Collection Date: 3/1/07 11:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|-------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Ethylbenzene | 6.7 | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| m,p-Xylene | 14 | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| o-Xylene | 10 | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| n-Propylbenzene | 2.4 | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,3,5-Trimethylbenzene | 6.1 | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,2,4-Trimethylbenzene | 15 | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Naphthalene | 9.0 | 5.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:22:00 PM |
| Surr: Dibromofluoromethane | 101 | 85-116 | | %REC | 1 | 3/9/07 2:22:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 106 | 77-127 | | %REC | 1 | 3/9/07 2:22:00 PM |
| Surr: Toluene-d8 | 101 | 86-114 | | %REC | 1 | 3/9/07 2:22:00 PM |
| Surr: 4-Bromofluorobenzene | 99.5 | 79-117 | | %REC | 1 | 3/9/07 2:22:00 PM |

AMRO Environmental Laboratories Corp.

Date: 15-Mar-07

CLIENT: SHAW E & I, Inc.
 Lab Order: 0703003
 Project: 101960 Textron
 Lab ID: 0703003-03A

Client Sample ID: MW-109D
 Collection Date: 3/1/07 1:00:00 PM
 Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|-------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: KT |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Trichlorofluoromethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| cis-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Chloroform | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Trichloroethene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Bromodichloromethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Toluene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Tetrachloroethene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |

AMRO Environmental Laboratories Corp.

Date: 15-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0703003
Project: 101960 Textron
Lab ID: 0703003-03A

Client Sample ID: MW-109D
Collection Date: 3/1/07 1:00:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|-------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Ethylbenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| m,p-Xylene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| o-Xylene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,3,5-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,2,4-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Naphthalene | ND | 5.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 3/9/07 2:57:00 PM |
| Surr: Dibromofluoromethane | 104 | 85-116 | | %REC | 1 | 3/9/07 2:57:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 107 | 77-127 | | %REC | 1 | 3/9/07 2:57:00 PM |
| Surr: Toluene-d8 | 101 | 86-114 | | %REC | 1 | 3/9/07 2:57:00 PM |
| Surr: 4-Bromofluorobenzene | 97.0 | 79-117 | | %REC | 1 | 3/9/07 2:57:00 PM |

AMRO Environmental Laboratories Corp.

Date: 15-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0703003
Project: 101960 Textron
Lab ID: 0703003-05A

Client Sample ID: Trip Blank
Collection Date: 3/1/07
Matrix: TRIP BLANK

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-------------------------------------|--------|----------------|------|-------|----|-------------------|
| EPA 8260B VOLATILES BY GC/MS | | | | | | |
| | | SW8260B | | | | Analyst: SK |
| Dichlorodifluoromethane | ND | 5.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Chloromethane | ND | 5.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Vinyl chloride | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Chloroethane | ND | 5.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Bromomethane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Trichlorofluoromethane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Diethyl ether | ND | 5.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Acetone | ND | 10 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,1-Dichloroethene | ND | 1.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Carbon disulfide | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Methylene chloride | ND | 5.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Methyl tert-butyl ether | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| trans-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,1-Dichloroethane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 2-Butanone | ND | 10 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 2,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| cis-1,2-Dichloroethene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Chloroform | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Tetrahydrofuran | ND | 10 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Bromochloromethane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,1,1-Trichloroethane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,1-Dichloropropene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Carbon tetrachloride | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,2-Dichloroethane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Benzene | ND | 1.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Trichloroethene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,2-Dichloropropane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Bromodichloromethane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Dibromomethane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 4-Methyl-2-pentanone | ND | 10 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| cis-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Toluene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| trans-1,3-Dichloropropene | ND | 1.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,1,2-Trichloroethane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,2-Dibromoethane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 2-Hexanone | ND | 10 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,3-Dichloropropane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Tetrachloroethene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Dibromochloromethane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |

AMRO Environmental Laboratories Corp.

Date: 15-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0703003
Project: 101960 Textron
Lab ID: 0703003-05A

Client Sample ID: Trip Blank
Collection Date: 3/1/07
Matrix: TRIP BLANK

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|--------|------|-------|----|-------------------|
| Chlorobenzene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Ethylbenzene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| m,p-Xylene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| o-Xylene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Styrene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Bromoform | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Isopropylbenzene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,2,3-Trichloropropane | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Bromobenzene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| n-Propylbenzene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 2-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 4-Chlorotoluene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,3,5-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| tert-Butylbenzene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,2,4-Trimethylbenzene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| sec-Butylbenzene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 4-Isopropyltoluene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,3-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,4-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| n-Butylbenzene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,2-Dichlorobenzene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,2,4-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Hexachlorobutadiene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Naphthalene | ND | 5.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| 1,2,3-Trichlorobenzene | ND | 2.0 | | µg/L | 1 | 3/6/07 4:25:00 PM |
| Surr: Dibromofluoromethane | 102 | 85-116 | | %REC | 1 | 3/6/07 4:25:00 PM |
| Surr: 1,2-Dichloroethane-d4 | 102 | 77-127 | | %REC | 1 | 3/6/07 4:25:00 PM |
| Surr: Toluene-d8 | 102 | 86-114 | | %REC | 1 | 3/6/07 4:25:00 PM |
| Surr: 4-Bromofluorobenzene | 95.2 | 79-117 | | %REC | 1 | 3/6/07 4:25:00 PM |

AMRO Environmental Laboratories Corp.

Date: 21-Mar-07

QC SUMMARY REPORT
Method Blank

CLIENT: SHAW E & I, Inc.
Work Order: 0703003
Project: 101960 Textron

Sample ID mb-03/06/07 Batch ID: R36052 Test Code: SW6260B Units: µg/L Analysis Date 3/6/07 11:14:00 AM Prep Date 3/6/07
Client ID: Run ID: V-3_070306A SeqNo: 597671

| Analyte | QC Sample | | QC Spike | | Original Sample | | %RPD | RPDLimit | Qua |
|--------------------------|-----------|-----|----------|--------|-----------------|--------------|------|----------|-----|
| | Result | RL | Amount | Result | HighLimit | or MS Result | | | |
| Dichlorodifluoromethane | ND | 5.0 | µg/L | | | | | | |
| Chloromethane | ND | 5.0 | µg/L | | | | | | |
| Vinyl chloride | ND | 2.0 | µg/L | | | | | | |
| Chloroethane | ND | 5.0 | µg/L | | | | | | |
| Bromomethane | ND | 2.0 | µg/L | | | | | | |
| Trichlorofluoromethane | ND | 2.0 | µg/L | | | | | | |
| Diethyl ether | ND | 5.0 | µg/L | | | | | | |
| Acetone | ND | 10 | µg/L | | | | | | |
| 1,1-Dichloroethene | ND | 1.0 | µg/L | | | | | | |
| Carbon disulfide | ND | 2.0 | µg/L | | | | | | |
| Methylene chloride | ND | 5.0 | µg/L | | | | | | |
| Methyl tert-butyl ether | ND | 2.0 | µg/L | | | | | | |
| trans-1,2-Dichloroethene | ND | 2.0 | µg/L | | | | | | |
| 1,1-Dichloroethane | ND | 2.0 | µg/L | | | | | | |
| 2-Butanone | ND | 10 | µg/L | | | | | | |
| 2,2-Dichloropropane | ND | 2.0 | µg/L | | | | | | |
| cis-1,2-Dichloroethene | ND | 2.0 | µg/L | | | | | | |
| Chloroform | ND | 2.0 | µg/L | | | | | | |
| Tetrahydrofuran | ND | 10 | µg/L | | | | | | |
| Bromochloromethane | ND | 2.0 | µg/L | | | | | | |
| 1,1,1-Trichloroethane | ND | 2.0 | µg/L | | | | | | |
| 1,1-Dichloropropene | ND | 2.0 | µg/L | | | | | | |
| Carbon tetrachloride | ND | 2.0 | µg/L | | | | | | |
| 1,2-Dichloroethane | ND | 2.0 | µg/L | | | | | | |
| Benzene | ND | 1.0 | µg/L | | | | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 21-Mar-07

QC SUMMARY REPORT
Method Blank

CLIENT: SHAW E & I, Inc.
Work Order: 0703003
Project: 101960 Textron

| | | | |
|---------------------------|----|-----|------|
| Trichloroethene | ND | 2.0 | µg/L |
| 1,2-Dichloropropane | ND | 2.0 | µg/L |
| Bromodichloromethane | ND | 2.0 | µg/L |
| Dibromomethane | ND | 2.0 | µg/L |
| 4-Methyl-2-pentanone | ND | 10 | µg/L |
| cis-1,3-Dichloropropene | ND | 1.0 | µg/L |
| Toluene | ND | 2.0 | µg/L |
| trans-1,3-Dichloropropene | ND | 1.0 | µg/L |
| 1,1,2-Trichloroethane | ND | 2.0 | µg/L |
| 1,2-Dibromoethane | ND | 2.0 | µg/L |
| 2-Hexanone | ND | 10 | µg/L |
| 1,3-Dichloropropane | ND | 2.0 | µg/L |
| Tetrachloroethene | ND | 2.0 | µg/L |
| Dibromochloromethane | ND | 2.0 | µg/L |
| Chlorobenzene | ND | 2.0 | µg/L |
| 1,1,1,2-Tetrachloroethane | ND | 2.0 | µg/L |
| Ethylbenzene | ND | 2.0 | µg/L |
| m,p-Xylene | ND | 2.0 | µg/L |
| o-Xylene | ND | 2.0 | µg/L |
| Styrene | ND | 2.0 | µg/L |
| Bromoform | ND | 2.0 | µg/L |
| Isopropylbenzene | ND | 2.0 | µg/L |
| 1,1,2,2-Tetrachloroethane | ND | 2.0 | µg/L |
| 1,2,3-Trichloropropane | ND | 2.0 | µg/L |
| Bromobenzene | ND | 2.0 | µg/L |
| n-Propylbenzene | ND | 2.0 | µg/L |
| 2-Chlorotoluene | ND | 2.0 | µg/L |
| 4-Chlorotoluene | ND | 2.0 | µg/L |
| 1,3,5-Trimethylbenzene | ND | 2.0 | µg/L |
| tert-Butylbenzene | ND | 2.0 | µg/L |
| 1,2,4-Trimethylbenzene | ND | 2.0 | µg/L |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 21-Mar-07

QC SUMMARY REPORT
Method Blank

CLIENT: SHAW E & I, Inc.
Work Order: 0703003
Project: 101960 Textron

| Compound | Reporting Limit | Concentration | Recovery | Recovery Limit | Concentration | Recovery | Recovery Limit |
|-----------------------------|-----------------|---------------|----------|----------------|---------------|----------|----------------|
| sec-Butylbenzene | ND | 2.0 | µg/L | | | | |
| 4-Isopropyltoluene | ND | 2.0 | µg/L | | | | |
| 1,3-Dichlorobenzene | ND | 2.0 | µg/L | | | | |
| 1,4-Dichlorobenzene | ND | 2.0 | µg/L | | | | |
| n-Butylbenzene | ND | 2.0 | µg/L | | | | |
| 1,2-Dichlorobenzene | ND | 2.0 | µg/L | | | | |
| 1,2-Dibromo-3-chloropropane | ND | 5.0 | µg/L | | | | |
| 1,2,4-Trichlorobenzene | ND | 2.0 | µg/L | | | | |
| Hexachlorobutadiene | ND | 2.0 | µg/L | | | | |
| Naphthalene | ND | 5.0 | µg/L | | | | |
| 1,2,3-Trichlorobenzene | ND | 2.0 | µg/L | | | | |
| Surr: Dibromofluoromethane | 25.3 | 2.0 | µg/L | 25 | 0 | 101 | 85 |
| Surr: 1,2-Dichloroethane-d4 | 26.77 | 2.0 | µg/L | 25 | 0 | 107 | 77 |
| Surr: Toluene-d8 | 26.03 | 2.0 | µg/L | 25 | 0 | 104 | 86 |
| Surr: 4-Bromofluorobenzene | 25.2 | 2.0 | µg/L | 25 | 0 | 101 | 79 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 21-Mar-07

CLIENT: SHAW E & I, Inc.
 Work Order: 0703003
 Project: 101960 Textron

QC SUMMARY REPORT
 Laboratory Control Spike - Full List

Sample ID Icsf-03/06/07 Batch ID: R36052 Test Code: SW6260B Units: µg/L Analysis Date 3/6/07 9:31:00 AM Prep Date 3/6/07
 Client ID: Run ID: V-3_070306A SeqNo: 597672

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Qua |
|--------------------------|------------------|-----|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Dichlorodifluoromethane | 25.64 | 5.0 | µg/L | 20 | 0 | 128 | 10 | 150 | 0 | 0 | 0 | 0 |
| Chloromethane | 22.39 | 5.0 | µg/L | 20 | 0 | 112 | 37 | 150 | 0 | 0 | 0 | 0 |
| Vinyl chloride | 22.78 | 2.0 | µg/L | 20 | 0 | 114 | 48 | 150 | 0 | 0 | 0 | 0 |
| Chloroethane | 20.86 | 5.0 | µg/L | 20 | 0 | 104 | 54 | 142 | 0 | 0 | 0 | 0 |
| Bromomethane | 21.97 | 2.0 | µg/L | 20 | 0 | 110 | 51 | 137 | 0 | 0 | 0 | 0 |
| Trichlorofluoromethane | 23.38 | 2.0 | µg/L | 20 | 0 | 117 | 62 | 141 | 0 | 0 | 0 | 0 |
| Diethyl ether | 20.11 | 5.0 | µg/L | 20 | 0 | 101 | 68 | 134 | 0 | 0 | 0 | 0 |
| Acetone | 21.04 | 10 | µg/L | 20 | 0 | 105 | 9 | 150 | 0 | 0 | 0 | 0 |
| 1,1-Dichloroethene | 20.98 | 1.0 | µg/L | 20 | 0 | 105 | 68 | 146 | 0 | 0 | 0 | 0 |
| Carbon disulfide | 23.28 | 2.0 | µg/L | 20 | 0 | 116 | 52 | 131 | 0 | 0 | 0 | 0 |
| Methylene chloride | 23.51 | 5.0 | µg/L | 20 | 0 | 118 | 67 | 138 | 0 | 0 | 0 | 0 |
| Methyl tert-butyl ether | 23 | 2.0 | µg/L | 20 | 0 | 115 | 63 | 139 | 0 | 0 | 0 | 0 |
| trans-1,2-Dichloroethene | 25.03 | 2.0 | µg/L | 20 | 0 | 125 | 81 | 126 | 0 | 0 | 0 | 0 |
| 1,1-Dichloroethane | 24.36 | 2.0 | µg/L | 20 | 0 | 122 | 78 | 124 | 0 | 0 | 0 | 0 |
| 2-Butanone | 17.58 | 10 | µg/L | 20 | 0 | 87.9 | 41 | 150 | 0 | 0 | 0 | 0 |
| 2,2-Dichloropropane | 29.07 | 2.0 | µg/L | 20 | 0 | 145 | 71 | 150 | 0 | 0 | 0 | 0 |
| cis-1,2-Dichloroethene | 20.5 | 2.0 | µg/L | 20 | 0 | 103 | 78 | 121 | 0 | 0 | 0 | 0 |
| Chloroform | 21.78 | 2.0 | µg/L | 20 | 0 | 109 | 82 | 123 | 0 | 0 | 0 | 0 |
| Tetrahydrofuran | 19.31 | 10 | µg/L | 20 | 0 | 96.6 | 51 | 146 | 0 | 0 | 0 | 0 |
| Bromochloromethane | 20.93 | 2.0 | µg/L | 20 | 0 | 105 | 77 | 131 | 0 | 0 | 0 | 0 |
| 1,1,1-Trichloroethane | 23.32 | 2.0 | µg/L | 20 | 0 | 117 | 81 | 127 | 0 | 0 | 0 | 0 |
| 1,1-Dichloropropene | 20.94 | 2.0 | µg/L | 20 | 0 | 105 | 76 | 119 | 0 | 0 | 0 | 0 |
| Carbon tetrachloride | 21.73 | 2.0 | µg/L | 20 | 0 | 109 | 76 | 129 | 0 | 0 | 0 | 0 |
| 1,2-Dichloroethane | 22.88 | 2.0 | µg/L | 20 | 0 | 114 | 76 | 127 | 0 | 0 | 0 | 0 |
| Benzene | 22.13 | 1.0 | µg/L | 20 | 0 | 111 | 81 | 118 | 0 | 0 | 0 | 0 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 21-Mar-07

QC SUMMARY REPORT

Laboratory Control Spike - Full List

CLIENT: SHAW E & I, Inc.
 Work Order: 0703003
 Project: 101960 Textron

| Compound | Reporting Limit | Concentration | Recovery | Accepted Recovery Limits | Recovery | Accepted Recovery Limits | Reporting Limit | Concentration | Recovery | Accepted Recovery Limits |
|---------------------------|-----------------|---------------|----------|--------------------------|----------|--------------------------|-----------------|---------------|----------|--------------------------|
| Trichloroethene | 22.95 | 2.0 | µg/L | 20 | 0 | 115 | 81 | 119 | 0 | 0 |
| 1,2-Dichloropropane | 21.53 | 2.0 | µg/L | 20 | 0 | 108 | 79 | 120 | 0 | 0 |
| Bromodichloromethane | 22.24 | 2.0 | µg/L | 20 | 0 | 111 | 77 | 131 | 0 | 0 |
| Dibromomethane | 20.92 | 2.0 | µg/L | 20 | 0 | 105 | 76 | 128 | 0 | 0 |
| 4-Methyl-2-pentanone | 17.75 | 10 | µg/L | 20 | 0 | 88.8 | 51 | 141 | 0 | 0 |
| cis-1,3-Dichloropropene | 21.39 | 1.0 | µg/L | 20 | 0 | 107 | 76 | 120 | 0 | 0 |
| Toluene | 22.24 | 2.0 | µg/L | 20 | 0 | 111 | 83 | 119 | 0 | 0 |
| trans-1,3-Dichloropropene | 21.05 | 1.0 | µg/L | 20 | 0 | 105 | 66 | 128 | 0 | 0 |
| 1,1,2-Trichloroethane | 21.36 | 2.0 | µg/L | 20 | 0 | 107 | 74 | 123 | 0 | 0 |
| 1,2-Dibromoethane | 21.49 | 2.0 | µg/L | 20 | 0 | 107 | 72 | 128 | 0 | 0 |
| 2-Hexanone | 16.68 | 10 | µg/L | 20 | 0 | 83.4 | 31 | 148 | 0 | 0 |
| 1,3-Dichloropropane | 20.94 | 2.0 | µg/L | 20 | 0 | 105 | 76 | 122 | 0 | 0 |
| Tetrachloroethene | 21.42 | 2.0 | µg/L | 20 | 0 | 107 | 81 | 124 | 0 | 0 |
| Dibromochloromethane | 20.17 | 2.0 | µg/L | 20 | 0 | 101 | 63 | 126 | 0 | 0 |
| Chlorobenzene | 20.59 | 2.0 | µg/L | 20 | 0 | 103 | 84 | 113 | 0 | 0 |
| 1,1,1,2-Tetrachloroethane | 20.91 | 2.0 | µg/L | 20 | 0 | 105 | 73 | 124 | 0 | 0 |
| Ethylbenzene | 21.54 | 2.0 | µg/L | 20 | 0 | 108 | 83 | 118 | 0 | 0 |
| m,p-Xylene | 43.65 | 2.0 | µg/L | 40 | 0 | 109 | 85 | 116 | 0 | 0 |
| o-Xylene | 20.55 | 2.0 | µg/L | 20 | 0 | 103 | 84 | 115 | 0 | 0 |
| Styrene | 21.46 | 2.0 | µg/L | 20 | 0 | 107 | 81 | 118 | 0 | 0 |
| Bromoform | 18.27 | 2.0 | µg/L | 20 | 0 | 91.4 | 55 | 126 | 0 | 0 |
| Isopropylbenzene | 22.33 | 2.0 | µg/L | 20 | 0 | 112 | 77 | 125 | 0 | 0 |
| 1,1,2,2-Tetrachloroethane | 17.8 | 2.0 | µg/L | 20 | 0 | 89 | 62 | 134 | 0 | 0 |
| 1,2,3-Trichloropropane | 18.79 | 2.0 | µg/L | 20 | 0 | 94 | 62 | 132 | 0 | 0 |
| Bromobenzene | 19.05 | 2.0 | µg/L | 20 | 0 | 95.2 | 78 | 119 | 0 | 0 |
| n-Propylbenzene | 21.49 | 2.0 | µg/L | 20 | 0 | 107 | 77 | 127 | 0 | 0 |
| 2-Chlorotoluene | 20.09 | 2.0 | µg/L | 20 | 0 | 100 | 78 | 118 | 0 | 0 |
| 4-Chlorotoluene | 20.81 | 2.0 | µg/L | 20 | 0 | 104 | 77 | 119 | 0 | 0 |
| 1,3,5-Trimethylbenzene | 21.18 | 2.0 | µg/L | 20 | 0 | 106 | 80 | 120 | 0 | 0 |
| tert-Butylbenzene | 21.55 | 2.0 | µg/L | 20 | 0 | 108 | 81 | 120 | 0 | 0 |
| 1,2,4-Trimethylbenzene | 20.84 | 2.0 | µg/L | 20 | 0 | 104 | 80 | 118 | 0 | 0 |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 21-Mar-07

CLIENT: SHAW E & I, Inc.
 Work Order: 0703003
 Project: 101960 Textron

QC SUMMARY REPORT
 Laboratory Control Spike - Full List

| Compound | 18.27 | 21.2 | 19.43 | 18.91 | 17.72 | 18.68 | 17.56 | 17.72 | 17.35 | 18.27 | 17.55 | 25.03 | 26.05 | 25.49 | 25.24 |
|-----------------------------|-------|------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|
| sec-Butylbenzene | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 5.0 | 2.0 | 2.0 | 5.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 4-Isopropyltoluene | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L | μg/L |
| 1,3-Dichlorobenzene | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 25 | 25 | 25 | 25 | 25 |
| 1,4-Dichlorobenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| n-Butylbenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 1,2-Dichlorobenzene | 91.4 | 106 | 97.2 | 94.6 | 88.6 | 93.4 | 87.8 | 88.6 | 86.8 | 91.4 | 87.8 | 100 | 104 | 102 | 101 |
| 1,2-Dibromo-3-chloropropane | 82 | 80 | 84 | 79 | 76 | 81 | 47 | 73 | 77 | 58 | 76 | 85 | 77 | 86 | 79 |
| 1,2,4-Trichlorobenzene | 123 | 126 | 115 | 117 | 128 | 117 | 136 | 126 | 134 | 138 | 124 | 116 | 127 | 114 | 117 |
| Hexachlorobutadiene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Naphthalene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| 1,2,3-Trichlorobenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Surr: Dibromofluoromethane | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Surr: 1,2-Dichloroethane-d4 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Surr: Toluene-d8 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |
| Surr: 4-Bromofluorobenzene | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 | 0 |

Qualifiers: ND - Not Detected at the Reporting Limit
 J - Analyte detected below quantitation limits
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

S - Spike Recovery outside accepted recovery limits
 R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank
 NA - Not applicable where J values or ND results occur

AMRO Environmental Laboratories Corp.

Date: 15-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0703003
Project: 101960 Textron
Lab ID: 0703003-02B

Client Sample ID: MW-221S
Tag Number:
Collection Date: 3/1/07 11:30:00 AM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|---------------------------------------|--------|----------------|------|-------|----|--------------------|
| TPH BY GC/FID (MODIFIED 8015B) | | SW8015B | | | | Analyst: FQ |
| Gasoline | ND | 0.50 | | mg/L | 10 | 3/10/07 3:29:00 AM |
| Mineral Spirits | ND | 0.50 | | mg/L | 10 | 3/10/07 3:29:00 AM |
| Kerosene | ND | 0.50 | | mg/L | 10 | 3/10/07 3:29:00 AM |
| Diesel Fuel/Fuel Oil #2 | ND | 0.50 | | mg/L | 10 | 3/10/07 3:29:00 AM |
| Motor Oil/Hydraulic Oil | ND | 1.0 | | mg/L | 10 | 3/10/07 3:29:00 AM |
| Unidentified Hydrocarbons | 35 | 2.0 | | mg/L | 10 | 3/10/07 3:29:00 AM |
| Surr: o-Terphenyl | 53.2 | 31-131 | | %REC | 10 | 3/10/07 3:29:00 AM |

Gasoline cannot be accurately determined by this method. Purge and trap sample introduction into a GC or GCMS is the recommended approach for gasoline. Due to the physical, chemical, and biological processes which affect the chemical composition of fuel mixtures exposed to the environment, the qualitative identity of a hydrocarbon mixture as a fuel product is not always conclusive by this method due to the method's reliance on chromatographic pattern recognition. A result provided for a specific fuel indicates that the mixture present in the sample has a chromatographic pattern similar to the laboratory's reference standard for that fuel mixture under specific GC operating conditions utilized at the time of analysis. A result identified as Unidentified Hydrocarbons is based upon the detector response obtained for the laboratory's Fuel Oil#2 reference standard and includes the entire chromatographic response for the sample between n-Alkanes of carbon numbers C9 to C36.

- Qualifiers:**
- ND - Not Detected at the Reporting Limit
 - J - Analyte detected below quantitation limits
 - B - Analyte detected in the associated Method Blank
 - H - Method prescribed holding time exceeded.
 - RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.
 - S - Spike Recovery outside accepted recovery limits
 - R - RPD outside accepted recovery limits
 - E - Value above quantitation range
 - # - See Case Narrative

Date: 21-Mar-07

AMRO Environmental Laboratories Corp.

CLIENT: SHAW E & I, Inc.
 Work Order: 0703003
 Project: 101960 Textron
QC SUMMARY REPORT
 Method Blank

Sample ID MB-16854 Batch ID: 16854 Test Code: SW6015B Units: mg/L Analysis Date 3/10/07 1:07:00 AM Prep Date 3/6/07
 Client ID: Run ID: GC-FING1_070309A SeqNo: 598887

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Qua |
|---------------------------|------------------|-------|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Gasoline | ND | 0.050 | mg/L | | | | | | | | | |
| Mineral Spirits | ND | 0.050 | mg/L | | | | | | | | | |
| Kerosene | ND | 0.050 | mg/L | | | | | | | | | |
| Diesel Fuel/Fuel Oil #2 | ND | 0.050 | mg/L | | | | | | | | | |
| Motor Oil/Hydraulic Oil | ND | 0.10 | mg/L | | | | | | | | | |
| Unidentified Hydrocarbons | ND | 0.20 | mg/L | | | | | | | | | |
| Surr. o-Terphenyl | 0.09015 | 0 | mg/L | 0.1 | 0 | 90.2 | 31 | 131 | 0 | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit, defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 21-Mar-07

QC SUMMARY REPORT
Laboratory Control Spike

CLIENT: SHAW E & I, Inc.
Work Order: 0703003
Project: 101960 Textron

Sample ID LCS-16854 Batch ID: 16854 Test Code: SW6015B Units: mg/L Analysis Date 3/10/07 1:43:00 AM Prep Date 3/6/07
Client ID: Run ID: GC-FING1_070309A SeqNo: 598888

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Qua |
|-------------------------|------------------|-------|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Diesel Fuel/Fuel Oil #2 | 1.407 | 0.050 | mg/L | 2 | 0 | 70.4 | 42 | 119 | 0 | | | |
| Surr: o-Terphenyl | 0.07872 | 0 | mg/L | 0.1 | 0 | 78.7 | 31 | 131 | 0 | | | |

Sample ID LCSD-16854 Batch ID: 16854 Test Code: SW6015B Units: mg/L Analysis Date 3/10/07 2:18:00 AM Prep Date 3/6/07
Client ID: Run ID: GC-FING1_070309A SeqNo: 598889

| Analyte | QC Sample Result | RL | Units | QC Spike Amount | Original Sample Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Qua |
|-------------------------|------------------|-------|-------|-----------------|------------------------|------|----------|-----------|------------------------------|------|----------|-----|
| Diesel Fuel/Fuel Oil #2 | 1.029 | 0.050 | mg/L | 2 | 0 | 51.5 | 42 | 119 | 1.407 | 31 | 40 | |
| Surr: o-Terphenyl | 0.06007 | 0 | mg/L | 0.1 | 0 | 60.1 | 31 | 131 | 0 | 0 | 0 | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 15-Mar-07

CLIENT: SHAW E & I, Inc.
Lab Order: 0703003
Project: 101960 Textron
Lab ID: 0703003-04A

Client Sample ID: GZA-3
Collection Date: 3/1/07 1:50:00 PM
Matrix: GROUNDWATER

| Analyses | Result | RL | Qual | Units | DF | Date Analyzed |
|-----------------------------|--------|------|---------|-------|----|--------------------|
| ICP METALS DISSOLVED SW-846 | | | SW6010B | | | Analyst: RK |
| Lead | ND | 12.0 | | µg/L | 1 | 3/9/07 12:29:25 PM |

AMRO Environmental Laboratories Corp.

Date: 19-Mar-07

QC SUMMARY REPORT
Method Blank

CLIENT: SHAW E & I, Inc.
Work Order: 0703003
Project: 101960 Textron

Sample ID MB-16884 Batch ID: 16844 Test Code: SW6010B Units: µg/L Analysis Date 3/9/07 11:14:52 AM Prep Date
Client ID: Run ID: ICP-OPTIMA_070309A SeqNo: 598186

| Analyte | QC Sample Result | RL | Units | QC Spike Original Sample Amount | Result | %REC | LowLimit | HighLimit | Original Sample or MS Result | %RPD | RPDLimit | Qua |
|---------|------------------|----|-------|---------------------------------|--------|------|----------|-----------|------------------------------|------|----------|-----|
| Lead | ND | 12 | µg/L | | | | | | | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.

AMRO Environmental Laboratories Corp.

Date: 19-Mar-07

QC SUMMARY REPORT
Laboratory Control Spike

CLIENT: SHAW E & I, Inc.
Work Order: 0703003
Project: 101960 Textron

| | | | | | | | | | | | |
|------------|-----------|------------------|--------------------|-----------------|---------|------------------------|------|---------------|--------------------|-----------|----------|
| Sample ID | LCS-16884 | Batch ID: | 16844 | Test Code: | SW6010B | Units: | µg/L | Analysis Date | 3/9/07 11:19:04 AM | Prep Date | |
| Client ID: | | Run ID: | ICP-OPTIMA_070309A | QC Spike Amount | 1998 | Original Sample Result | 0 | SeqNo: | 598187 | | |
| Analyte | | QC Sample Result | 1994 | RL | 12 | Units | µg/L | LowLimit | 80 | %RPD | RPDLimit |
| Lead | | | | | | | | HighLimit | 120 | %RPD | RPDLimit |
| | | | | | | | | | | | Qua |
| | | | | | | | | | | | |

| | | | | | | | | | | | |
|------------|------------|------------------|--------------------|-----------------|---------|------------------------|------|---------------|--------------------|-----------|----------|
| Sample ID | LCS-16884D | Batch ID: | 16844 | Test Code: | SW6010B | Units: | µg/L | Analysis Date | 3/9/07 11:24:34 AM | Prep Date | |
| Client ID: | | Run ID: | ICP-OPTIMA_070309A | QC Spike Amount | 1998 | Original Sample Result | 0 | SeqNo: | 598188 | | |
| Analyte | | QC Sample Result | 1975 | RL | 12 | Units | µg/L | LowLimit | 80 | %RPD | RPDLimit |
| Lead | | | | | | | | HighLimit | 120 | %RPD | RPDLimit |
| | | | | | | | | | | | Qua |
| | | | | | | | | | | | |

Qualifiers: ND - Not Detected at the Reporting Limit S - Spike Recovery outside accepted recovery limits B - Analyte detected in the associated Method Blank
 J - Analyte detected below quantitation limits R - RPD outside accepted recovery limits NA - Not applicable where J values or ND results occur
 RL - Reporting Limit; defined as the lowest concentration the laboratory can accurately quantitate.