

Periodic Groundwater Report For Site 15 Hancock Air National Guard Base Draft Final

Site:

Hancock Air National Guard Base
Syracuse, New York

Prepared for:

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List of Acronyms and Abbreviations

ANG	Air National Guard
ARAR	Applicable or Relevant and Appropriate Requirements
AWQS	Ambient Water Quality Standards
BEX	Benzene, Ethylbenzene, Xylenes
bgs	Below Ground Surface
cfm	Cubic feet per minute
COC	Compounds of Concern
DO	Dissolved Oxygen
DOD	Department of Defense
ELAP	Environmental Laboratory Accreditation Program
ERM	Environmental Resources Management
ERP	Environmental Restoration Program
FFS	Focused Feasibility Study
FSP	Field Sampling Plan
ft	Feet
HANGB	Hancock Air National Guard Base
IRM	Interim Remedial Measure
JP	Jet Propulsion
lbs	Pounds
mg/L	Milligram per liter
µg/L	Micrograms per Liter
MNA	Monitored Natural Attenuation
MW	Monitoring Well
NGB	National Guard Bureau
NYS	New York State
NYSDEC	New York State Department of Environmental Conservation
ORP	Oxidation Reduction Potential
PCB	Polychlorinated Biphenyls
pH	Pondus Hydrogenii
PVC	Polyvinyl Chloride
QAPP	Quality Assurance Project Plan
RAO	Remedial Action Objective
RAWP	Remedial Action Work Plan
ROD	Record of Decision
ROI	Radius of Influence
STARS	Spills Technology and Remediation Series
USEPA	United States Environmental Protection Agency
VOC	Volatile Organic Compound

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1 Introduction

This Periodic Groundwater Report summarizes the progress of the remedial activities implemented to address petroleum hydrocarbon groundwater contamination associated with Environmental Restoration Program (ERP) Site 15 at the 174th Fighter Wing, New York Air National Guard (ANG), Hancock Air National Guard Base (HANGB), Syracuse, New York. A Final Record of Decision (ROD) was issued by the ANG in April 2011 for the remediation of groundwater at Site 15. The remedy selected in the ROD consists of the injection of calcium peroxide to enhance aerobic biodegradation of the dissolved phase petroleum hydrocarbon groundwater plume, coupled with long-term monitoring.

The remedial activities conducted to date at ERP Site 15 under this contract consist of the installation of a biosparging system to address residual contamination in soil in October 2011; and an initial injection of calcium peroxide substrate conducted in October 2012. A baseline groundwater monitoring event was conducted in September 2012 and the first post-injection groundwater monitoring event was conducted in January 2013. This report provides a comparison of pre- and post-injection groundwater conditions at ERP Site 15 including groundwater monitoring results from June and August 2013.

The activities summarized in this report were conducted in accordance with the *Final Remedial Action Work Plan* (WP) submitted to the New York State Department of Environmental Conservation (NYSDEC) in September 2011 and approved by NYSDEC in October 2011. The WP was prepared in accordance with NYSDEC Program Policy, DER-10; Chapter 5, Section 5.3.

1.1 Site Description

HANGB is located in Syracuse, New York and ERP Site 15 comprises approximately 2.5 acres in the southeastern portion of the HANGB. Site 15 was formerly used as a pump house and fuel storage facility where numerous spills of jet propulsion (JP)-4 and JP-8 military aviation fuels and polychlorinated biphenyls (PCBs) occurred during active use. The pump house, tanks, and associated piping and structures have been demolished and the majority of the Site is now a large open field. Site 15 was listed as a Class 2 site on the NYS Inactive Hazardous Waste Disposal Site Registry in 1994 as Site Number 734054. PCBs identified at the site have been successfully remediated. A map illustrating the location of the HANGB and ERP Site 15 is provided as Figure 1.

The recurrent releases of JP-4 and JP-8 at ERP Site 15 resulted in contamination of soil and a dissolved-phase plume of petroleum hydrocarbons with maximum historic dimensions of approximately 1,000 feet (ft) along the axis in the north-south direction and a width of approximately 150 ft. The plume is aligned with groundwater flow at Site 15, which is to the south-southeast towards Ley Creek and Onondaga Lake. The source area is located at the northern end of the plume in the vicinity of monitoring well MW-101; the down-gradient edge of the plume is at the boundary of the General Electric Property. The contaminants of concern (COCs) within the plume are benzene, ethylbenzene, and xylenes (BEX). The historic maximum detection of BEX occurred at MW-19 in September 2005 at a concentration of 1,500 micrograms per liter ($\mu\text{g/L}$). The first post injection monitoring event was conducted in January 2013, BEX compounds exceeded cleanup criteria in only three (3) of 19 monitoring wells sampled, and in each instance the total BEX concentration was less than 50 $\mu\text{g/L}$. The most recent sampling event was conducted in August 2013 and there was only one detection of Benzene in excess of the cleanup criteria. ERP Site 15 monitoring well locations are depicted in Figure 2 along with groundwater elevation contours and flow directions. Historic plume dimensions are depicted in Figure 3.

1.1.1 Site Geology

The surficial geology at Site 15 consists of glaciofluvial sediments deposited by glacial melt water underlain by poorly sorted till deposited directly by glaciers. The glaciofluvial sediments include clayey silts, sands and gravels with thickness ranging from 45 to 55 ft. The underlying till consists of gravel, cobbles, and boulders entrained in a clayey silt matrix and ranges in thickness from 30 to 100 ft (Lockheed 1997).

Bedrock is encountered at depths ranging from 75 to 109 ft below ground surface (bgs), and is part of the Upper Silurian Vernon Formation. This formation consists of thinly bedded soft red shale with thin beds of green shale, gypsum, halite and dolomite. Competence varies from soft and crumbly to dense and hard. The degree of competence appears to be proportional to the density of the fractures in the shale. The shale is characterized by enlarged fractures, joints and bedding planes (Lockheed 1997). Significant portions of ERP Site 15 have been re-graded and filled due to previous construction and demolition activities.

1.1.2 Site Hydrogeology

The overburden at Site 15 consists of fine-grained sediments. The subgrade soils are fairly uniform with the upper 10 to 15 ft of the soil characterized by relatively soft, dark yellowish-brown silt and clayey silt. Towards the southeast, the interval thins to approximately 5 ft. Beneath the clayey silt are yellowish brown to dark brown fine to medium-grained sands with silt and trace amounts of clay down to a depth of approximately 20 ft. Underlying these silty sands is a lens of stiff clayey silts (often called glacial till). Till has been encountered as much as 15 ft thick (Lockheed 1997).

Groundwater is generally encountered at depths of 1.5 to 13 ft bgs and, as previously discussed, flows in a south to southeasterly direction towards Ley Creek and eventually into Onondaga Lake. Groundwater contours based on elevation data gathered during the September 2013 performance monitoring event are depicted in Figure 2.

1.2 Previous Remedial Activities

Based on an evaluation of the site conditions, the compounds of concern (COC), and an analysis of applicable or relevant and appropriate requirements (ARARs), the following remedial action objectives (RAOs) were developed for groundwater contaminated with BEX at ERP Site 15:

- Prevent exposure to contaminated groundwater containing BEX concentrations above the NYSDEC Ambient Water Quality Standards (AWQS) and Guidance Values;
- Prevent or minimize further off-site migration of the contaminant plume (plume containment);
- Prevent or minimize further migration of contaminants from source materials to groundwater (source control); and
- Enhance the natural process for the attenuation of BEX compounds on-site and off-site.

Achievement of RAOs will be quantitatively measured by the achievement of NYSDEC AWQS included in *NYS Division of Water Technical and Operational Guidance Series (1.1.1) 1998*. The AWQS for BEX are included on Table 1.

Environmental studies performed from 1990 to 2009 identified Site 15 and down-gradient off-site areas as having soil and groundwater impacted with petroleum hydrocarbons. A Focused Feasibility Study (FFS) recommended excavation and off-site disposal of the source areas (these were completed in 2003 and 2008) and focused enhanced aerobic bioremediation with monitored natural attenuation (MNA). Two vadose zone source area removals have occurred as interim remedial measures (IRMs):

- **Year 2003:** removal of 5,360 tons of petroleum-impacted soil from the vadose zone, steel tanks and associated piping; and
- **Year 2008:** excavation of 2,890 tons of petroleum-impacted soil from the vadose zone source area followed by application of 4,800 pounds of an oxygen-releasing product (calcium peroxide) within the bottom of the excavation areas.

The location of the IRM soil excavations is presented in Figure 3. Calcium peroxide was applied directly to the bottom of the 2008 excavations prior to backfilling with crushed concrete. Post excavation confirmation soil sampling and groundwater monitoring results from wells adjacent to the excavations indicated that there was an area of residual soil contamination as well as a lack of oxygen available in this portion of the aquifer, potentially limiting biodegradation in the residual petroleum impacted soil. Biosparging was selected for accelerated biodegradation of the COCs in the source area as the presence of petroleum hydrocarbons adsorbed to soil in the saturated region can act as a continuous source of dissolved phase contamination; limiting the effectiveness of calcium peroxide injections in achieving RAOs in a reasonable timeframe.

A biosparge system was designed and installed in October 2011 to inject air into the saturated zone to stimulate aerobic biodegradation of residual source area smear zone impacts. Fifteen one-inch diameter polyvinyl chloride (PVC) biosparge wells screened from 18 to 20 ft bgs were installed at Site 15 in accordance with the WP. The biosparge well locations are depicted in Figure 3. The biosparging system was operated from November 2011 through December 2012, when it was shutdown to monitor for contaminant rebound. During operation, filtered atmospheric air was delivered to each well at a rate of 1 to 2 cubic feet per minute (cfm). The radius of influence (ROI) for sparge wells was estimated at up to 30 feet, based on dissolved oxygen measurements taken at ERP Site 15 monitoring wells. The temporary electrical drop to the biosparge system has been removed and the system will be shipped off site at the end of November 2013. The biosparge points will be decommissioned along with the monitoring wells after the site is closed.

2 Calcium Peroxide Injections

The prescribed remedy for remediation of BEX impacted groundwater is enhanced aerobic biodegradation via calcium peroxide injection. The injection of calcium peroxide provides and extended release of oxygen into the subsurface to maintain an aerobic environment; enhancing microbial activity which results in the degradation of petroleum compounds. Delivery of calcium peroxide to the aquifer is accomplished by pressure injection, which ensures a uniform distribution across the injection interval and allows for delivery of the substrate at a faster rate. Distribution of calcium within the aquifer is by advection along the natural hydraulic gradient.

An initial calcium peroxide injection event was conducted from October 2-5, 2012. A total of 2,200 pounds of calcium peroxide were injected at 44 injection locations. Each injection location received 50 pounds of calcium peroxide, which was injected as a 40-percent slurry comprised of 9 gallons of water and 50 pounds of calcium peroxide. Substrate was injected using direct-push technology (DPT) techniques. Substrate was injected using a 'bottom up' approach across a 15-foot injection interval at each location. Injection intervals ranged from 3- to 18-feet bgs to 15- to 30-feet bgs, dependent on the water table elevation in monitoring wells located adjacent to the injection locations. The injection locations are depicted in Figure 4.

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3 Groundwater Monitoring

3.1 Groundwater Monitoring Program Overview

3.1.1 Rationale

The groundwater monitoring program at ERP Site 15 was developed to provide baseline characterization of the BEX plume by sampling the 30 existing monitoring wells; and for performance monitoring to assess the effectiveness of biosparging and calcium peroxide injection in remediating BEX contamination. Performance monitoring will consist of sampling up to 30 monitoring wells on a quarterly basis for two years after the initial injection event, followed by semi-annual sampling of up to 20 monitoring wells for two additional years. Samples are analyzed for VOCs and geochemical data is also collected to assess the performance of the remedial action. After four quarters of non-detect or detections below the NYSDEC AGWQS at any particular well, it will be proposed that the well be removed from the sampling network.

3.2 Sampling Methodology

3.2.1 Sample Collection

Groundwater samples are collected using low-flow (minimal draw down) sampling techniques. Wells are gauged with an oil-water interface probe prior to sampling to determine the water level, total well depth and the presence/absence of non-aqueous phase liquids (NAPL). Water level data is utilized to develop potentiometric surface maps. Following gauging, the wells are purged using a peristaltic pump and dedicated tubing. During purging, the depth to water is monitored as well as geochemical and physical parameters including temperature, pH, dissolved oxygen (DO), oxidation-reduction potential (ORP), specific conductivity and turbidity. The geochemical and physical parameters are utilized to determine well stabilization which indicates that the water being purged is from the formation surrounding the well and will provide a representative sample. VOC samples are collected in 40 milliliter (mL) vials preserved with hydrochloric acid provided by the analytical laboratory and shipped on ice to the laboratory for analysis. Geochemical and physical parameter readings are recorded on field sheets; and the geochemical data, particularly DO and ORP, are evaluated as indicators of the favorability of the aquifer for promoting aerobic biodegradation of BEX. In some instances, the hydraulic conductivity at a given well is insufficient to allow for low-flow sampling. In such cases, the well was purged dry, allowed to recharge and sampled with a disposable bailer. Additional information regarding sampling techniques can be found in the *Final Field Sampling Plan* (AECOM, 2011), which is included as an appendix of the WP.

3.2.2 Sample Analysis

Samples are analyzed for BEX using United States Environmental Protection Agency (USEPA) Method 8260. Samples are analyzed at a Spectrum Analytical, Inc. laboratories in North Kingstown, Rhode Island; a Department of Defense (DoD) Environmental Laboratory Accreditation Program (ELAP) certified facility. In addition to environmental samples, quality assurance/quality control (QA/QC) samples are also submitted to the laboratory. These include duplicates, matrix spike and matrix spike duplicate (MS/MSD) samples, trip blanks, and ambient blanks. Additional information on analytical procedures and QA/QC procedures can be found in the Quality Assurance Program Plan (QAPP), which is included as an appendix of the WP.

3.2.3 Sample Locations

Sample locations are selected from the existing 30 well monitoring network at ERP Site 15. All 30 wells were sampled during the October 2012 and January 13 events. Wells in which BEX was not detected in these events as well as previous events were dropped from future events as proposed in the April 2013 Periodic Groundwater Report. During the June and August 2013 sampling events 10 monitoring wells were sampled. These wells had detections of VOCs following the calcium peroxide injections. Future events will include fewer locations as wells are dropped from the network based on contaminant trends. The sampling locations and historic plume dimensions are depicted in Figure 5.

3.3 Groundwater Monitoring Results

The following sections present groundwater monitoring results from five sampling events. The October 2010 event was conducted by Environmental Resources Management (ERM) and predates the installation of the biosparge system and calcium peroxide injection. The second event, conducted by AECOM, occurred in September 2012, following 10 months of biosparging and prior to the initial calcium peroxide injection event. The third event presented in this report occurred in January 2013,

approximately three months after the initial calcium peroxide injection event. The fourth event discussed in this report was from a sampling event that occurred in June 2013, eight months after initial calcium peroxide injection, and the final event discussed in this report details the results from the most recent, August 2013, sampling event.

3.3.1 October 2010 Groundwater Monitoring Results

Thirty (30) monitoring wells were sampled by ERM in October 2010. COCs were detected in 10 of the 30 wells sampled; with detections at 6 wells exceeding the NYSDEC AWQS. Benzene was detected in samples from six (6) monitoring wells at concentrations ranging from 0.36J-µg/L to 26-µg/L. Five of the six detections of benzene exceeded the NYSDEC AWQS of 1-µg/L. The maximum detection of benzene occurred at MW-11. Ethylbenzene was detected in samples from six wells at concentrations ranging from 3.8-µg/L to 100-µg/L. Five of the six detections of ethylbenzene exceeded the NYSDEC AWQS of 5-µg/L. The maximum detection of ethylbenzene occurred at MW-19. Xylenes were detected in samples from eight wells at concentrations ranging from 0.72J-µg/L to 92-µg/L. Three of the eight detections of xylenes exceeded the NYSDEC AWQS of 5-µg/L. The maximum detection of xylenes occurred at MW-105. MW-112 was the furthest down-gradient monitoring well with detections of COCs in October 2010. Benzene, ethylbenzene, and xylenes were detected at MW-112 at concentrations of 0.36J-µg/L, 10-µg/L, and 2.7-µg/L, respectively. The detection of ethylbenzene at MW-112 exceeded the NYSDEC AWQS. Groundwater analytical data is summarized in Table 1 and the extent of the BEX plume during the October 2010 monitoring event is depicted in Figure 5.

Geochemical parameter monitoring conducted during sampling indicated that conditions were generally anaerobic (DO less than 1-mg/L) and reducing (ORP below 0 mV) throughout ERP Site 15. These conditions are typically of sites contaminated with petroleum hydrocarbons that have not undergone chemical or biological amendment. Geochemical data for the October 2010 monitoring event is summarized in Table 2.

3.3.2 September 2012 Groundwater Monitoring Results

Thirty (30) monitoring wells were sampled in September 2012. The intent of this monitoring event was to serve as a baseline for evaluating the effectiveness of the calcium peroxide injections to be conducted in October 2012. The biosparging system had been operational for approximately 10 months at the time of the September 2012 monitoring event. COCs were detected in samples from seven wells during the September 2012 monitoring event. Benzene was detected in samples from three monitoring wells at concentrations ranging from 1.4-µg/L to 5.9-µg/L. All three detections of benzene exceeded the NYSDEC AWQS of 1-µg/L. The maximum detection of benzene occurred at MW-103. Ethylbenzene was detected in samples from four wells at concentrations ranging from 1.7-µg/L to 53-µg/L. Three detections exceeded the NYSDEC AWQS of 5-µg/L. The maximum detection of ethylbenzene occurred at MW-112. Xylenes were detected at three monitoring wells at concentrations ranging from 26-µg/L to 67-µg/L; all of which exceeded the NYSDEC AWQS of 5-µg/L. The maximum detection of xylenes occurred at MW-112. As during the October 2010 event, MW-112 was the furthest down-gradient monitoring well with detections of COCs. Ethylbenzene was detected in the sample from MW-112 at 29-µg/L and xylenes were detected at 18-µg/L. Both of these detections exceed their respective NYSDEC AWQS of 5-µg/L. Groundwater analytical data is summarized in Table 1 and the extent of the BEX plume during the September 2012 monitoring event is depicted in Figure 5.

Geochemical parameter monitoring conducted during sampling indicated that dissolved-oxygen concentrations had increased in many areas since the October 2010 monitoring event, as would be expected following 10 months of biosparging. ORP readings were still within the reducing range throughout the majority of ERP Site 15. Geochemical data for the September 2012 monitoring event is summarized in Table 2.

3.3.3 January 2013 Groundwater Monitoring Results

Thirty (30) monitoring wells were sampled at ERP Site 15 in January 2013 to assess the performance of calcium peroxide injections and biosparging at reducing dissolved-phase BEX contamination. COCs were detected in samples from four monitoring wells, with at least one COC exceeding the NYSDEC AWQS at three wells. This represents a significant decrease from October 2010, when COCs were detected in ten wells and exceeded the NYSDEC AWQS at six wells. The frequency of COC detections and exceedances of the NYSDEC AWQS also decreased from the September 2012 monitoring event. Benzene was detected in samples from two wells at concentrations of 0.78J-µg/L and 4.7-µg/L. The detection of 4.7-µg/L, at MW-11, exceeded the NYSDEC AWQS of 1-µg/L. Ethylbenzene was detected in samples from three monitoring wells at concentrations of 0.63J-µg/L to 29-µg/L. The detection of ethylbenzene at 29-µg/L, which occurred at MW-112, exceeded the NYSDEC AWQS of 1-µg/L. Xylenes were detected at MW-112 and MW-101 at concentrations of 18-µg/L and 26-µg/L, respectively. Both detections of xylene exceeded the NYSDEC AWQS of 1-µg/L. As in the previous monitoring events, MW-112 was the furthest down-gradient well with BEX impacts. The data from the January 2013 monitoring indicate that there is no longer a contiguous BEX plume at ERP Site 15 but rather pockets of residual groundwater contamination.

DO and ORP measurements taken while sampling show increases in DO at numerous wells when compared to October 2010 and September 2012 data. While DO concentrations decreased at some wells from September 2012, the values were generally

higher than those in October 2010, prior to the implementation of biosparging and calcium peroxide injection. ORP data shows many wells still within the reducing range.

3.3.4 June 2013 Groundwater Monitoring Results

During the June 2013 sampling event, 10 monitoring wells were sampled (MW-14, MW-101, RW-01, MW-11, MW-15, MW-17, MW-19, MW-103, MW-105, MW-106, and MW-112). The other 19 wells showed no sign of contamination for four consecutive sampling events, therefore, were no longer sampled. This change to the monitoring network was proposed in the April 2013 Periodic Groundwater Report, Table 3. Out of the 11 wells sampled, COCs were detected in the sample from one well, MW-101. Contaminants found included Ethyl Benzene (5.7 µg/L) and Total Xylene (7.2 µg/L), both in exceedance of NYSDEC standard of 5 µg/L.

DO measurements taken during this sampling event showed both increases and decreases in many of the wells. Monitoring wells that showed a decrease in DO (MW-11, MW-15, MW-101, MW-112, and RW-1) dropped much more significantly than the monitoring wells that showed an increase (MW-14, MW-19, MW-103, MW-105, and MW-106,) went up in value. Wells with a decreasing DO dropped on an average of 2.775 mg/l, while monitoring wells that increased only rose 0.506 mg/l on average. ORP continued to fluctuate as it has in the past, but showed an overall increasing trend. The decreasing trend in DO is expected as the calcium peroxide is depleted.

3.3.5 August 2013 Groundwater Monitoring Results

A total of 10 monitoring wells were sampled on August 29, 2013 (MW-14, MW-101, RW-01, MW-11, MW-15, MW-19, MW-103, MW-105, MW-106, and MW-112). The sample from MW-11 contained benzene at a concentration of 9.6 µg/L which is above the NYSDEC standard of 1 µg/L. The result for the June 2013 sample from MW-11 was non-detect for benzene. The COCs ethyl-benzene (3.1 µg/L) and xylene (1.9 µg/L) were detected below NYSDEC standards in the sample from MW-101. These COCs had been detected above NYSDEC standards in the previous three quarterly samples. COCs were not detected in samples from the other 9 wells which were sampled during this event.

DO concentrations have stabilized below 1 mg/L in all of the wells sampling. This suggests that the calcium peroxide injected in the fall of 2012 has been depleted. The ORP measurements from this round of sampling continues to fluctuate but overall is showing a negative trend which is consistent with the depletion of the calcium peroxide.

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4 Conclusions and Recommendations

The post remedial groundwater monitoring data collected in September 2012, January 2013, June 2013, and August 2013 indicates that the addition of oxygen to the saturated zone at ERP Site 15 via bioparging and calcium peroxide injection has resulted in a significant decrease in BEX contamination within the aquifer, as evidenced by decreases in the frequency and magnitude of BEX detections over time. The current data indicates that the plume is no longer contiguous and that BEX impacts are isolated to one monitoring well, MW-11. The rebound seen in this well is typical and monitoring will be ongoing.

Calcium peroxide is typically effective in providing dissolved-oxygen to an aquifer for a period of 6 to 12 months, depending on groundwater flow rates, contaminant concentrations and utilization rates. The falling DO level in the monitoring wells suggests that the calcium peroxide injected in the fall of 2012 is nearing depletion. The next performance monitoring event is scheduled for December 2013. Should monitoring results indicate that BEX is still present at concentrations in excess of the NYSDEC AWQS, a second calcium peroxide injection will be conducted. This injection would possibly be augmented using a sodium-persulfate product, which degrades contaminants chemically; while the calcium peroxide would persist to stimulate biodegradation.

Quarterly groundwater monitoring should be continued until BEX concentrations are below the AWQS for four consecutive quarters as specified in DER-10 Section 6.4. The monitoring well network for December 2013 and subsequent events will include all wells where BEX compounds were detected in any of the three previous quarterly sampling events. The proposed monitoring well network is provided in Table 3.

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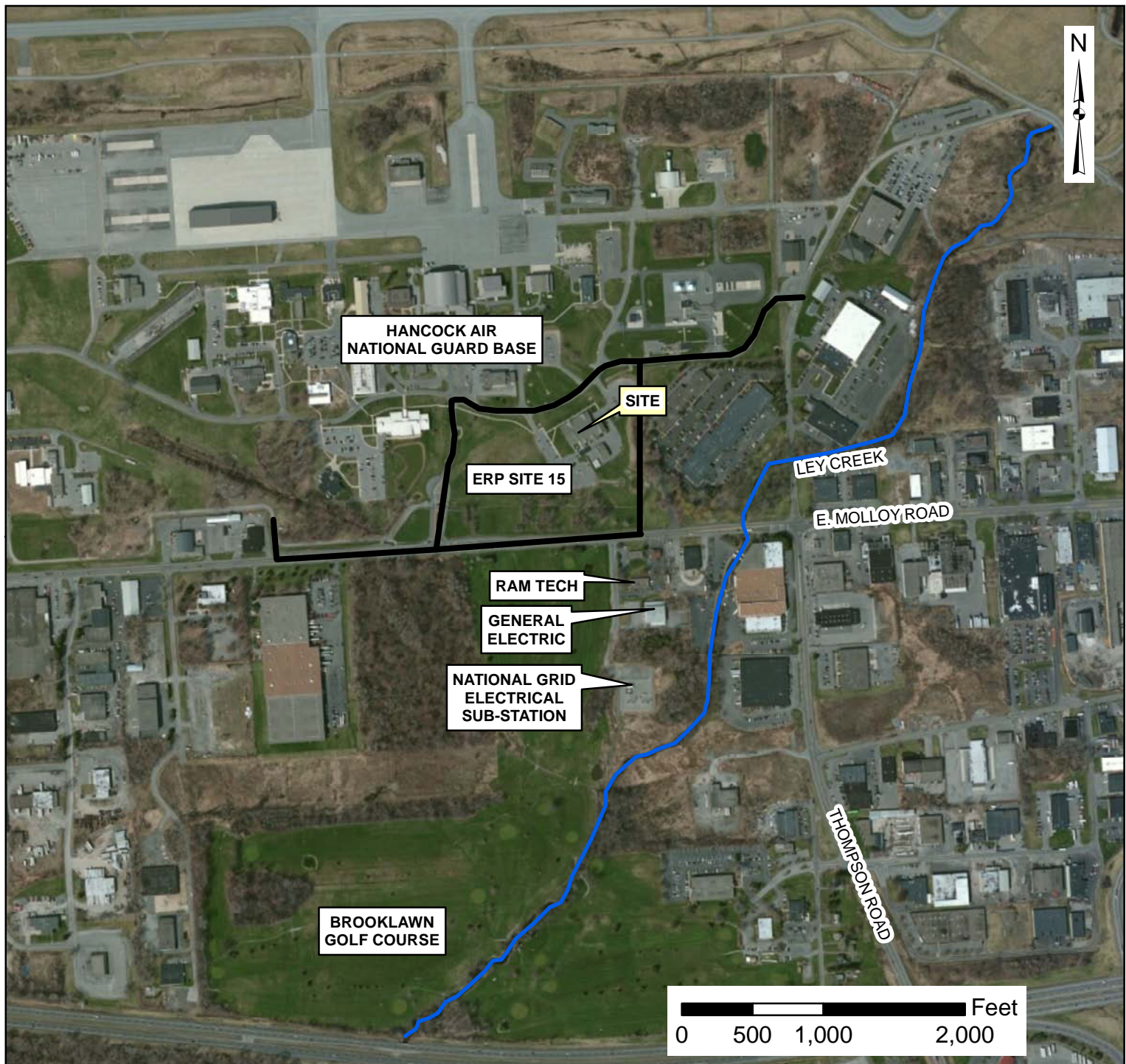
5 References

- AECOM, 2011. *Remedial Action Work Plan for Site 15* - Hancock Air National Guard Base - Syracuse, New York, September 2011.
- ERM, 2009. *Site 15 Final Construction Completion Report - Source Area Soil Removal*, 174th Fighter Wing - New York Air National Guard- Hancock Air National Guard Base - Syracuse, New York, January 2009.
- ERM, 2010. *Final Focused Feasibility Study* - 174th Fighter Wing - New York Air National Guard - Hancock Air National Guard Base - Syracuse, New York, March 2010.
- ERM, 2010. *Final Proposed Plan for Environmental Cleanup* - 174th Fighter Wing - New York Air National Guard- Hancock Air National Guard Base - Syracuse, New York, July 2010.
- ERM, 2011. *Environmental Restoration Program Final Record of Decision for Site 15- 74th Fighter Wing* - New York Air National Guard - Hancock Air National Guard Base - Syracuse, New York, April 2011.
- Lockheed, 1997. *Final Remedial Investigation Report for Petroleum, Oil and Lubricant Facility, Site 15*. Volumes I and II. Air National Guard Readiness Center, Andrews AFB, Maryland, July 1997.
- NYSDEC, 1998. Ambient Water Quality Standards and Guidance Values and Groundwater Effluent Limitations. *NYSDEC Division of Water Technical and Operational Guidance Series Memorandum Number 1.1.1.*, June 1998 (latest amendment April 2000).
- NYSDEC, 2010. DER-10, *Final Technical Guidance for Site Investigation and Remediation*, Division of Environmental Remediation, May 2010.

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Figures

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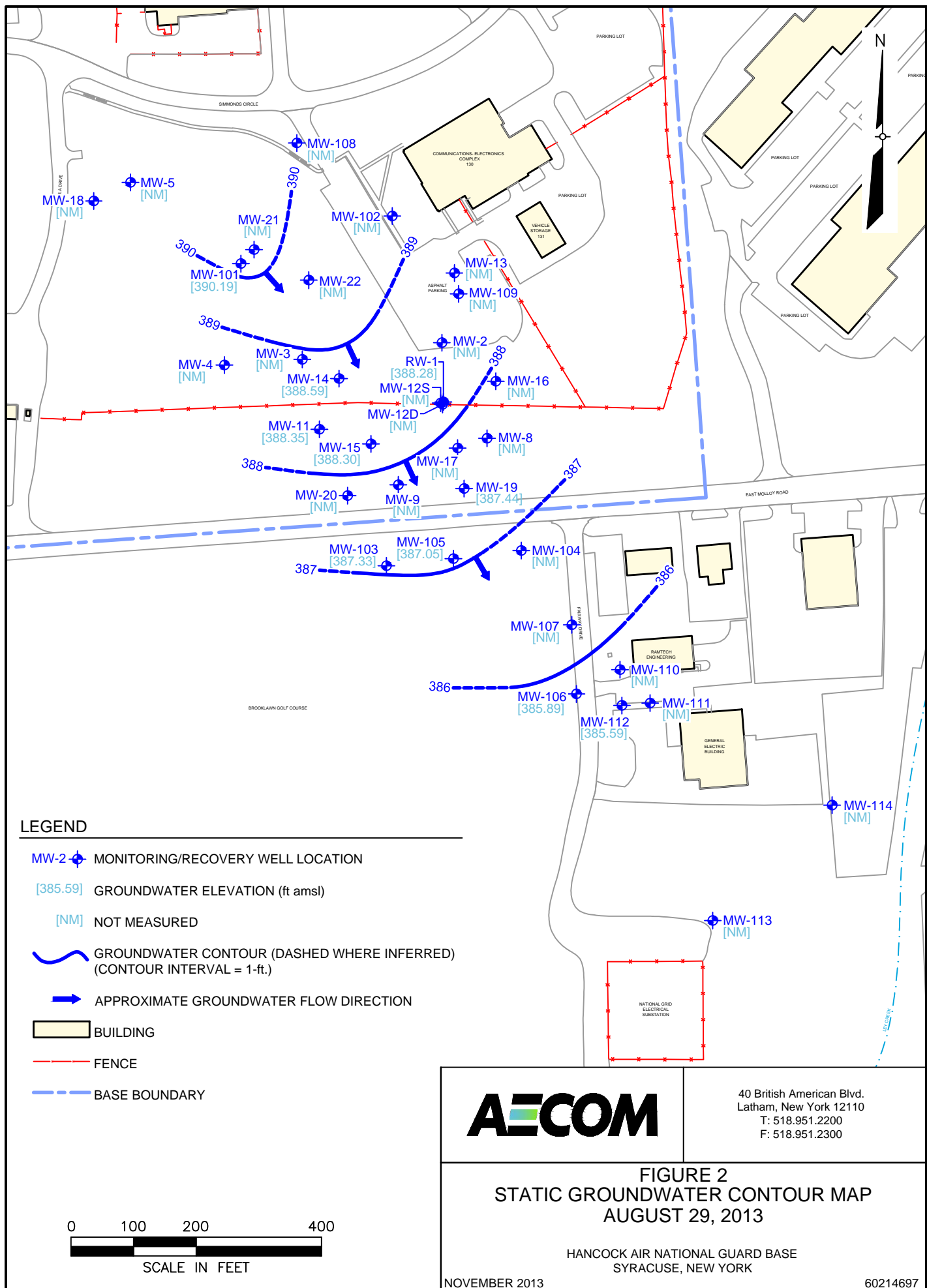


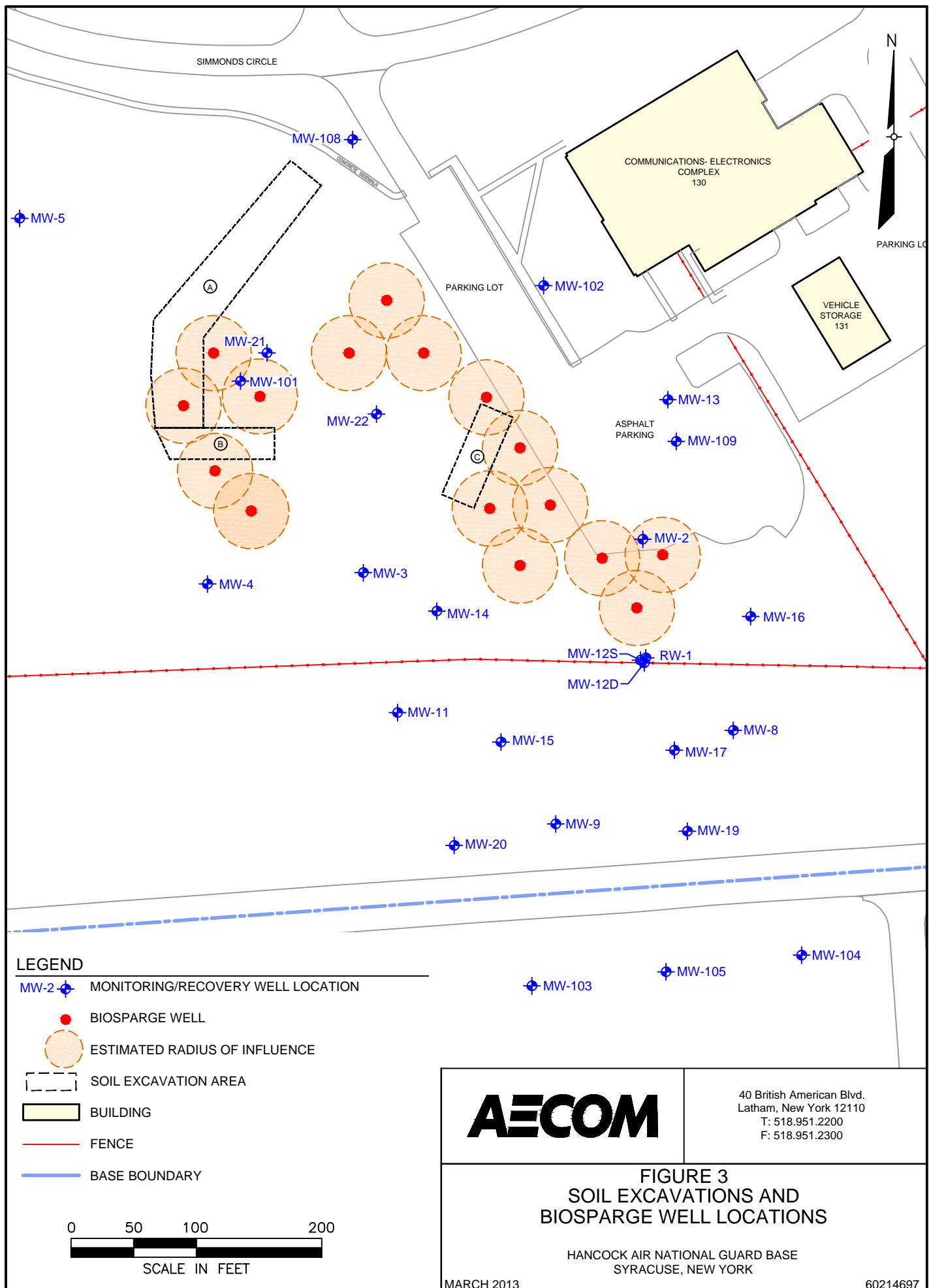
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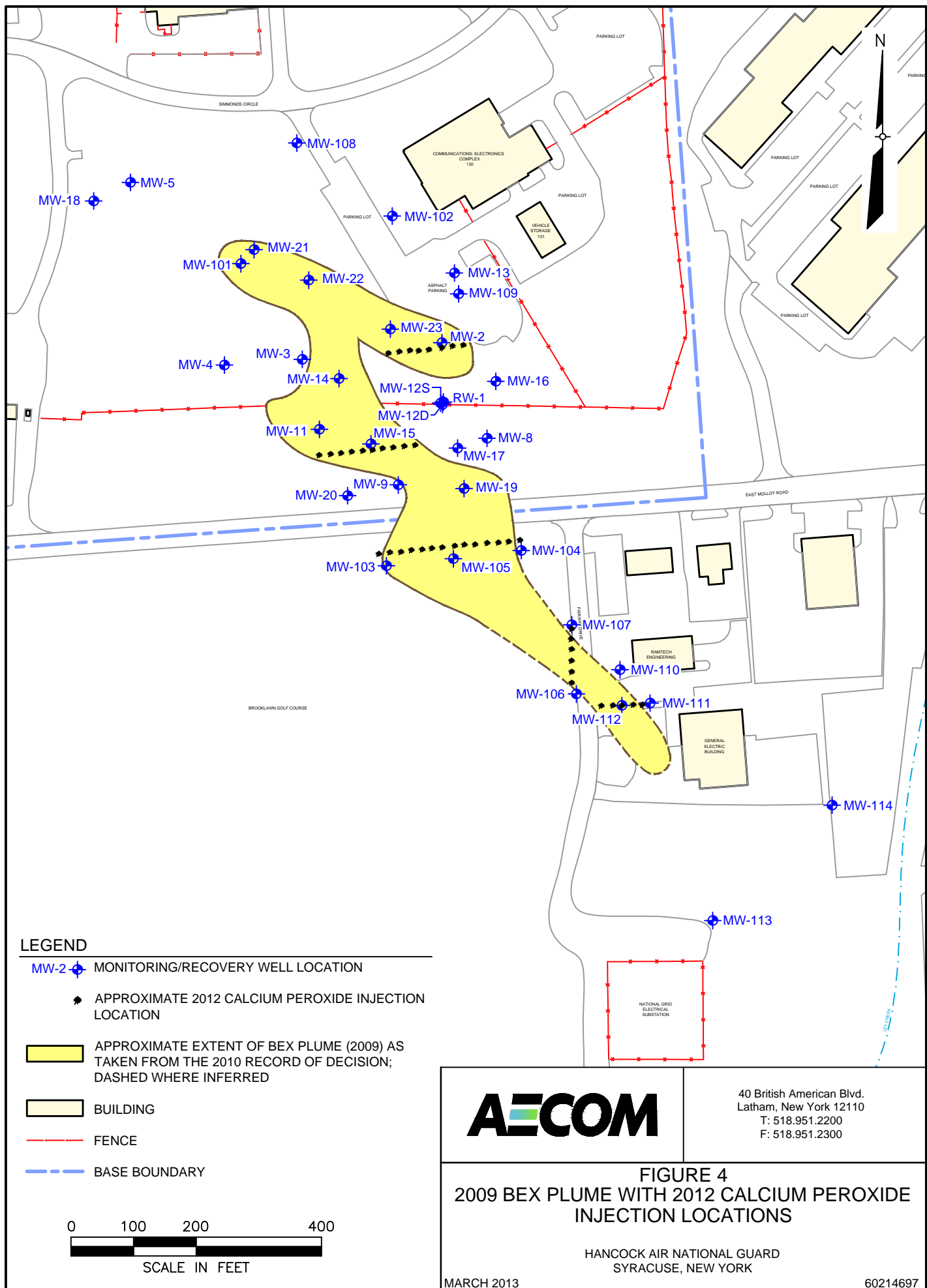
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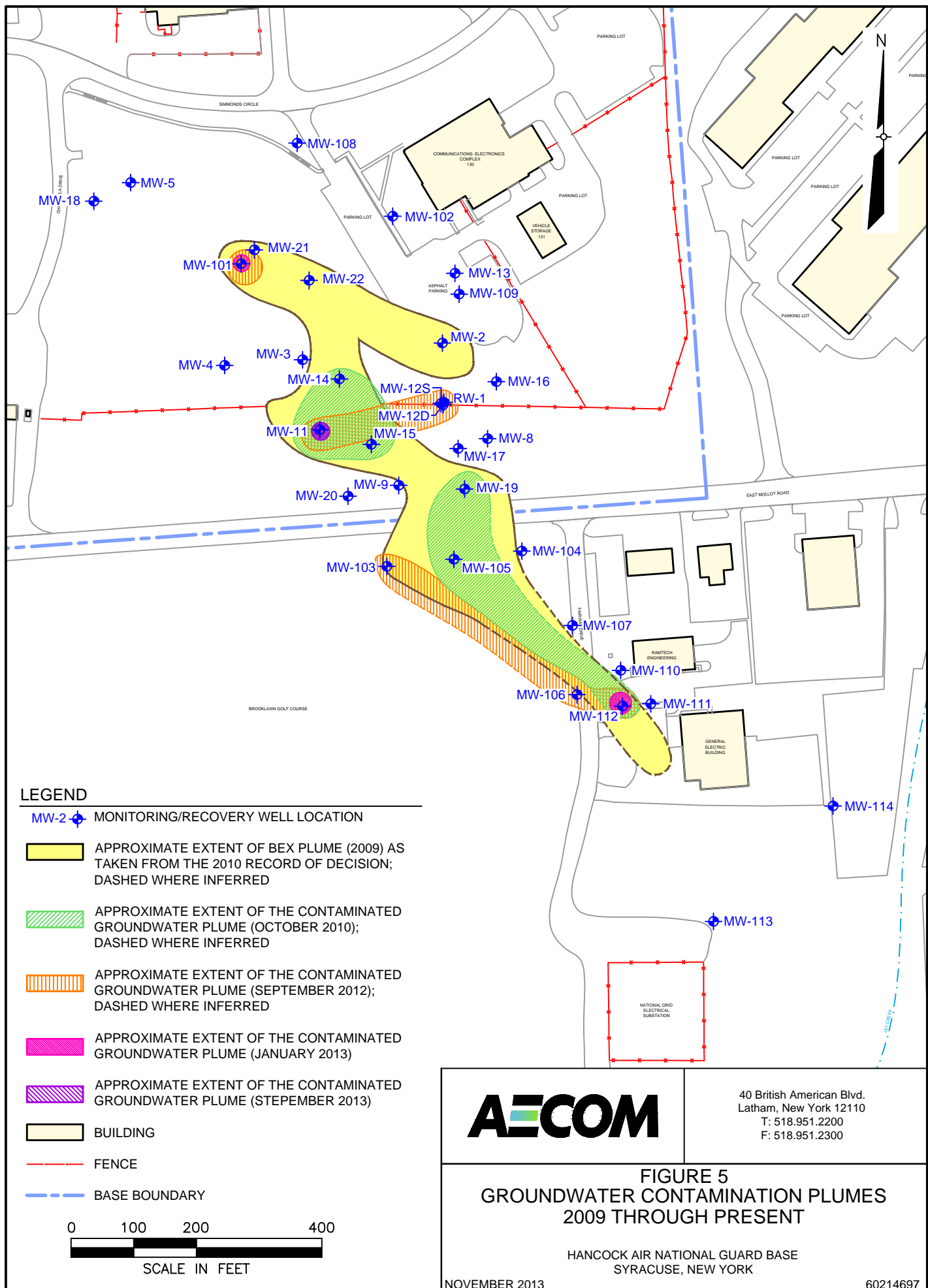
FIGURE 1
SITE LOCATION MAP

HANCOCK AIR NATIONAL GUARD
SYRACUSE, NEW YORK









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Tables

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Table 1
Groundwater Sampling Results
Hancock Air National Guard Base

Well ID	NYSDEC Std or GV	MW-112					MW-106					MW-19					MW-15					MW-11				
Sample Date		Oct-10	Sep-12	Jan-13	Jun-13	Aug-13	Oct-10	Sep-12	Jan-13	Jun-13	Aug-13	Oct-10	Sep-12	Jan-13	Jun-13	Aug-13	Oct-10	Sep-12	Jan-13	Jun-13	Aug-13	Oct-10	Sep-12	Jan-13	Jun-13	Aug-13
VOCs µg/L																										
Benzene	1	0.36 J	-	-	-	-	-	1.4	-	-	-	6.2	-	-	-	-	5.3	-	-	-	-	26	3.1	4.7	-	9.6
Ethyl Benzene	5	10	53	29	-	-	-	-	-	-	-	100	-	-	-	-	9.8	-	0.63 J	-	-	7.8	-	-	-	-
Total Xylene	5	2.7	67	18	-	-	-	-	-	-	-	18	-	-	-	-	0.72 J	-	-	-	-	9.5	-	-	-	-

Well ID	NYSDEC Std or GV	MW-103					MW-105					RW-1					MW-14					MW-101				
Sample Date		Oct-10	Sep-12	Jan-13	Jun-13	Aug-13	Oct-10	Sep-12	Jan-13	Jun-13	Aug-13	Oct-10	Sep-12	Jan-13	Jun-13	Aug-13	Oct-10	Sep-12	Jan-13	Jun-13	Aug-13	Oct-10	Sep-12	Jan-13	Jun-13	Aug-13
VOCs µg/L																										
Benzene	1	-	5.9	-	-	-	5.5	-	-	-	0.95 J	-	-	-	-	-	1.7	0.98 J	-	-	-	-	-	-	-	-
Ethyl Benzene	5	0.24 J	-	-	-	-	97	1.7	-	-	-	3.8	13	-	-	-	-	-	-	-	-	3.2	5.2	1.9	5.7	3.1
Total Xylene	5	-	-	-	-	-	92	-	-	-	-	2.8	27	-	-	-	-	-	-	-	-	2.0	26	26	7.2	1.9

Notes:
 NYSDEC GV or Std - Results compared to the New York State Department of Environmental Conservation (NYSDEC) Division of Water Technical and Operational Guidance Series (TOGS 1.1.1), 1998, standards (Std) and guidance values (GV)
 VOCs - Volatile organic compounds determined by United States Environmental Protection Agency (USEPA) Method 8260
 µg/L - Micrograms per liter
 -- - The compound was not detected at a concentration above the laboratory reporting limit
 Blank cells indicate that the compound was not analyzed for
 Shading indicates compounds detected above NYSDEC standards/guidance values
 J - The analyte was positively identified, the quantitation is an approximation

Table 2
Groundwater Field Parameters
Hancock Air National Guard Base

	MW-2		MW-3		MW-4		MW-5		MW-8		MW-9		MW-11		MW-14		MW-15		MW-16		MW-17		MW-18		MW-19		MW-20		MW-22		MW-101		MW-102		MW-103		MW-104		MW-105		MW-106		MW-107		MW-110		MW-111		MW-112		RW-1	
Date	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)	DO (mg/L)	ORP (mV)				
Oct-10	0.45	42.1	1.22	83.9	2.93	6.2	3.23	66	0.32	-97.4	0.85	43.8	0.36	-24.9	0.44	23.1	0.42	-7.1	0.84	114.8	0.58	-6.5	3.86	61.7	0.45	-66	0.61	8.9	1.87	-17.7	0.58	-5.5	0.45	30.7	0.51	3.3	1.75	25.9	0.68	2.6	0.68	30.3	2.57	34.2	0.5	25.3	0.69	-35.7	0.82	-106.9	0.46	-145.8
Nov-11	5.31	-318.8	2.32	-218.8	1.5	-210.2	5.83	-277.5	-	-	-	-	-	-	0.7	-268.6	-	-	-	-	-	-	4.86	-277.9	-	-	-	-	3.82	-249.2	-	-	0.14	-334.5	-	-	-	-	-	-	-	-	-	-	-	-	-	-				
Sep-12	6.3	-57	0.58	-79.5	0.66	-2.2	-	-	3.9	-88.2	4.34	-69.1	0.72	-122	1.04	-94.5	4.04	-142.9	1.04	214.4	0.38	-111	0.79	-93.6	0.34	-103.5	1.08	-94.2	9.31	9.1	1.23	-128.1	3.37	-89.3	2.75	-66.4	1.66	-38.9	4.09	-139.5	1.9	0.8	3.63	-124.4	-	-	-	-	0.4	-101.1		
Dec-12	5.11	-133.1	7.06	-93.8	2.91	-87.7	4	27.1	4.1	-10.8	4.81	47.2	3.73	-99.4	3.04	-46.5	2.96	-75.7	5.82	-14.2	8.02	-54.6	6.17	67.6	5.23	-67.1	3.4	-78.1	10.69	-24.9	4.13	-71.8	0.68	-52.8	4.1	-70.9	3.77	46.8	3.46	-112.4	2.04	-70.3	11.43	-53.8	-	-	-	-	3.08	-115.6		
Jan-13	1.34	-88	-0.03	-1.3	-	-	-	-	-	-	10.76	-8	4.75	-106.9	0.25	-52.9	2.39	-111.3	-	-	3.11	-96.3	-	-	-0.36	-47.4	-	-	3.88	168.8	2.65	-102	-	-	-0.1	-78.9	5.24	-15.4	0.01	-119.8	1.14	-68.4	5.09	14.8	0.24	70.9	1.78	-1.1	2.47	-146	2.62	6.2
Jun-13	-	-	-	-	-	-	-	-	-	-	-	-	0.23	-72.3	0.42	-9	0.9	-103.6	-	-	0.26	-134	-	-	0.65	-66.2	-	-	-	-	0.4	-172.4	-	-	0.17	-43.6	-	-	0.83	-91.2	1.61	-44.8	-	-	-	-	0.27	-82.7	0.57	-131.3		
Aug-13	-	-	-	-	-	-	-	-	-	-	-	-	0.18	-100	0.19	-43.6	0.31	-101.6	-	-	-	-	-	-	0.2	-80	-	-	-	-	0.13	-129.9	-	-	0.33	-83.8	-	-	0.11	-112.8	0.61	-35.9	-	-	-	-	0.19	-89.4	0.33	-107.9		

Table 3
Proposed Groundwater Monitoring Well Network
Hancock Air National Guard Base, Syracuse, NY

Previous Groundwater Monitoring Well Network	Proposed Groundwater Monitoring Well Network
MW-2	--
MW-3	--
MW-4	--
MW-5	--
MW-8	--
MW-9	--
MW-11	MW-11
MW-14	--
MW-15	--
MW-16	--
MW-17	--
MW-18	--
MW-19	--
MW-20	--
MW-22	--
MW-101	MW-101
MW-102	--
MW-103	MW-103
MW-104	--
MW-105	--
MW-106	MW-106
MW-107	--
MW-108	--
MW-109	--
MW-110	--
MW-111	--
MW-112	MW-112
MW-113	--
MW-114	--
RW-1	RW-1

Notes:

The proposed groundwater monitoring well network is based on the results of the baseline (September 2012) and performance monitoring (Jan, Jun, Sep 2013) groundwater sampling events. Wells that did not have any compounds of concern detected above laboratory reporting limits for four consecutive quarters (-) are proposed to be dropped from the sampling plan.

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Appendix A

Well Inspection Forms

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SITE NAME:

Hancock

SITE ID.:

60214697

INSPECTOR:

GW/HV

DATE/TIME:

11/30/11

WELL ID.:

mw-2

MONITORING WELL FIELD INSPECTION LOG

WELL VISIBLE? (If not, provide directions below)

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>

WELL COORDINATES? NYTM X _____ NYTM Y _____

PDOP Reading from Trimble Pathfinder: _____ Satellites: _____

GPS Method (circle) Trimble And/Or Magellan

WELL I.D. VISIBLE?

YES	NO
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>

WELL LOCATION MATCH SITE MAP? (if not, sketch actual location on back).....

WELL I.D. AS IT APPEARS ON PROTECTIVE CASING OR WELL:

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>

SURFACE SEAL PRESENT?

SURFACE SEAL COMPETENT? (If cracked, heaved etc., describe below)

PROTECTIVE CASING IN GOOD CONDITION? (If damaged, describe below)3 bolts missing

HEADSPACE READING (ppm) AND INSTRUMENT USED.....

TYPE OF PROTECTIVE CASING AND HEIGHT OF STICKUP IN FEET (If applicable)

PROTECTIVE CASING MATERIAL TYPE:

MEASURE PROTECTIVE CASING INSIDE DIAMETER (Inches):

NA
flushmount
steel

LOCK PRESENT?

LOCK FUNCTIONAL?

DID YOU REPLACE THE LOCK?

IS THERE EVIDENCE THAT THE WELL IS DOUBLE CASED? (If yes, describe below)

WELL MEASURING POINT VISIBLE?

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>

MEASURE WELL DEPTH FROM MEASURING POINT (Feet):

MEASURE DEPTH TO WATER FROM MEASURING POINT (Feet):

MEASURE WELL DIAMETER (Inches):

WELL CASING MATERIAL:

PHYSICAL CONDITION OF VISIBLE WELL CASING:

ATTACH ID MARKER (if well ID is confirmed) and IDENTIFY MARKER TYPE

PROXIMITY TO UNDERGROUND OR OVERHEAD UTILITIES.....

13.35
9.08
2"
steel
good
NA

DESCRIBE ACCESS TO WELL: (Include accessibility to truck mounted rig, natural obstructions, overhead power lines, proximity to permanent structures, etc.); ADD SKETCH OF LOCATION ON BACK, IF NECESSARY.

GoodDESCRIBE WELL SETTING (For example, located in a field, in a playground, on pavement, in a garden, etc.)
AND ASSESS THE TYPE OF RESTORATION REQUIRED.Open grassy field

IDENTIFY ANY NEARBY POTENTIAL SOURCES OF CONTAMINATION, IF PRESENT

(e.g. Gas station, salt pile, etc.):

REMARKS:

Temp 15.99
Spec Cond 1.521
Spec Cond 1.260

DO% 54.2
DO 5.32
Sketch

pH 6.69
ORP -318.8

SITE NAME:

Hancock

SITE ID.:

60214697

INSPECTOR:

C. C. / A. C.

DATE/TIME:

11/30

WELL ID.:

MW-3

MONITORING WELL FIELD INSPECTION LOG

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>

WELL VISIBLE? (If not, provide directions below)

WELL COORDINATES? NYTM X _____ NYTM Y _____

PDOP Reading from Trimble Pathfinder: _____ Satellites: _____

GPS Method (circle) Trimble And/Or Magellan

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>

WELL I.D. VISIBLE?

WELL LOCATION MATCH SITE MAP? (if not, sketch actual location on back).....

WELL I.D. AS IT APPEARS ON PROTECTIVE CASING OR WELL: MW-3

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>

SURFACE SEAL PRESENT?

SURFACE SEAL COMPETENT? (If cracked, heaved etc., describe below)

PROTECTIVE CASING IN GOOD CONDITION? (If damaged, describe below) 2 bolts missing

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>

HEADSPACE READING (ppm) AND INSTRUMENT USED.....

TYPE OF PROTECTIVE CASING AND HEIGHT OF STICKUP IN FEET (If applicable)

PROTECTIVE CASING MATERIAL TYPE:

MEASURE PROTECTIVE CASING INSIDE DIAMETER (Inches):

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>

LOCK PRESENT?

LOCK FUNCTIONAL?

DID YOU REPLACE THE LOCK?

IS THERE EVIDENCE THAT THE WELL IS DOUBLE CASED? (If yes, describe below) Steel (4")

WELL MEASURING POINT VISIBLE?

MEASURE WELL DEPTH FROM MEASURING POINT (Feet):

MEASURE DEPTH TO WATER FROM MEASURING POINT (Feet):

MEASURE WELL DIAMETER (Inches):

WELL CASING MATERIAL:

PHYSICAL CONDITION OF VISIBLE WELL CASING:

ATTACH ID MARKER (if well ID is confirmed) and IDENTIFY MARKER TYPE

PROXIMITY TO UNDERGROUND OR OVERHEAD UTILITIES.....

13.70
8.51
2"
Sched 40
good
NA

DESCRIBE ACCESS TO WELL: (Include accessibility to truck mounted rig, natural obstructions, overhead power lines, proximity to permanent structures, etc.); ADD SKETCH OF LOCATION ON BACK, IF NECESSARY.

good

DESCRIBE WELL SETTING (For example, located in a field, in a playground, on pavement, in a garden, etc.) AND ASSESS THE TYPE OF RESTORATION REQUIRED.

open field

IDENTIFY ANY NEARBY POTENTIAL SOURCES OF CONTAMINATION, IF PRESENT

(e.g. Gas station, salt pile, etc.):

NA

REMARKS:

2 Steel casings

Temp 13.06 DO% 22.1 ORP -218.8
Spec Cond .34 DO 2.32 pH 5.97
Spec Cond .242

Sketch

SITE NAME:

Hancock

SITE ID.:

60214617

INSPECTOR:

G. W. AL

DATE/TIME:

11/30/11

WELL ID.:

Mw-4

MONITORING WELL FIELD INSPECTION LOG

WELL VISIBLE? (If not, provide directions below)

YES	NO
X	

WELL COORDINATES? NYTM X _____ NYTM Y _____

PDOP Reading from Trimble Pathfinder: _____ Satellites: _____

GPS Method (circle) Trimble And/Or Magellan

WELL I.D. VISIBLE?

YES	NO
	X
X	

WELL LOCATION MATCH SITE MAP? (if not, sketch actual location on back).....

WELL I.D. AS IT APPEARS ON PROTECTIVE CASING OR WELL:

YES	NO
X	
X	
X	

SURFACE SEAL PRESENT?

SURFACE SEAL COMPETENT? (If cracked, heaved etc., describe below)

PROTECTIVE CASING IN GOOD CONDITION? (If damaged, describe below) 2 bulks missing

NA
flush mount

HEADSPACE READING (ppm) AND INSTRUMENT USED.....

TYPE OF PROTECTIVE CASING AND HEIGHT OF STICKUP IN FEET (If applicable)

PROTECTIVE CASING MATERIAL TYPE:

MEASURE PROTECTIVE CASING INSIDE DIAMETER (Inches):

YES	NO
X	
	X
	X
	X
X	X

LOCK PRESENT?

LOCK FUNCTIONAL?

DID YOU REPLACE THE LOCK?

IS THERE EVIDENCE THAT THE WELL IS DOUBLE CASED? (If yes, describe below)

WELL MEASURING POINT VISIBLE?

MEASURE WELL DEPTH FROM MEASURING POINT (Feet):

MEASURE DEPTH TO WATER FROM MEASURING POINT (Feet):

MEASURE WELL DIAMETER (Inches):

WELL CASING MATERIAL:

PHYSICAL CONDITION OF VISIBLE WELL CASING:

ATTACH ID MARKER (if well ID is confirmed) and IDENTIFY MARKER TYPE

PROXIMITY TO UNDERGROUND OR OVERHEAD UTILITIES.....

15.60
8.59
2"
4" steel
(crack)
NA
No

DESCRIBE ACCESS TO WELL: (Include accessibility to truck mounted rig, natural obstructions, overhead power lines, proximity to permanent structures, etc.); ADD SKETCH OF LOCATION ON BACK, IF NECESSARY.

Good → In open field

DESCRIBE WELL SETTING (For example, located in a field, in a playground, on pavement, in a garden, etc.)

AND ASSESS THE TYPE OF RESTORATION REQUIRED.

Open grassy field

IDENTIFY ANY NEARBY POTENTIAL SOURCES OF CONTAMINATION, IF PRESENT

(e.g. Gas station, salt pile, etc.):

REMARKS:

Double cased → steel casing Temp: 12.93°C Spec Cond: 0.434 ms/cm
DO: 14.5% DO: 1.50 mg/l 0.335 ms/cm

Sketch

pH: 6.58

CRP: -2102

SITE NAME:

Hancock

SITE ID.:

6021469-7

INSPECTOR:

GWS/BV

DATE/TIME:

11/30/11

WELL ID.:

NW-5

MONITORING WELL FIELD INSPECTION LOG

YES NO

WELL VISIBLE? (If not, provide directions below)

WELL COORDINATES? NYTM X _____ NYTM Y _____

PDOP Reading from Trimble Pathfinder: _____ Satellites: _____

GPS Method (circle) Trimble And/Or Magellan

YES NO

WELL I.D. VISIBLE?

WELL LOCATION MATCH SITE MAP? (if not, sketch actual location on back).....

WELL I.D. AS IT APPEARS ON PROTECTIVE CASING OR WELL:

YES NO

SURFACE SEAL PRESENT?

SURFACE SEAL COMPETENT? (If cracked, heaved etc., describe below)

PROTECTIVE CASING IN GOOD CONDITION? (If damaged, describe below)

HEADSPACE READING (ppm) AND INSTRUMENT USED:

TYPE OF PROTECTIVE CASING AND HEIGHT OF STICKUP IN FEET (If applicable)

PROTECTIVE CASING MATERIAL TYPE:

MEASURE PROTECTIVE CASING INSIDE DIAMETER (Inches):

YES NO

LOCK PRESENT?

LOCK FUNCTIONAL?

DID YOU REPLACE THE LOCK?

IS THERE EVIDENCE THAT THE WELL IS DOUBLE CASED? (If yes, describe below)

WELL MEASURING POINT VISIBLE?

MEASURE WELL DEPTH FROM MEASURING POINT (Feet):

MEASURE DEPTH TO WATER FROM MEASURING POINT (Feet):

MEASURE WELL DIAMETER (Inches):

WELL CASING MATERIAL:

PHYSICAL CONDITION OF VISIBLE WELL CASING:

ATTACH ID MARKER (if well ID is confirmed) and IDENTIFY MARKER TYPE

PROXIMITY TO UNDERGROUND OR OVERHEAD UTILITIES.....

15.22

5.75

2"

good steel

good

DESCRIBE ACCESS TO WELL: (Include accessibility to truck mounted rig, natural obstructions, overhead power lines, proximity to permanent structures, etc.); ADD SKETCH OF LOCATION ON BACK, IF NECESSARY.

Good

DESCRIBE WELL SETTING (For example, located in a field, in a playground, on pavement, in a garden, etc.) AND ASSESS THE TYPE OF RESTORATION REQUIRED.

Open, grassy field

IDENTIFY ANY NEARBY POTENTIAL SOURCES OF CONTAMINATION, IF PRESENT (e.g. Gas station, salt pile, etc.):

NA

REMARKS:

Temp 12.60

Spec Cond 1549

Spec Cond 1419

DO¹⁰ 50.7

DO 5.83

Sketch

pH 6.68
Torp -272.5

SITE NAME:

Harroct

SITE ID.:

60214697

INSPECTOR:

GWS/AV

DATE/TIME:

11/30/11

WELL ID.:

1110 11

MONITORING WELL FIELD INSPECTION LOG

WELL VISIBLE? (If not, provide directions below)

YES	NO
X	

WELL COORDINATES? NYTM X _____ NYTM Y _____

PDOP Reading from Trimble Pathfinder: _____ Satellites: _____

GPS Method (circle) Trimble And/Or Magellan

YES	NO
X	
X	

WELL I.D. VISIBLE?

WELL LOCATION MATCH SITE MAP? (if not, sketch actual location on back).....

WELL I.D. AS IT APPEARS ON PROTECTIVE CASING OR WELL:

YES	NO
X	
X	
X	

SURFACE SEAL PRESENT?

SURFACE SEAL COMPETENT? (If cracked, heaved etc., describe below)

PROTECTIVE CASING IN GOOD CONDITION? (If damaged, describe below)

YES	NO
X	
X	
X	

HEADSPACE READING (ppm) AND INSTRUMENT USED.....

TYPE OF PROTECTIVE CASING AND HEIGHT OF STICKUP IN FEET (If applicable) 3 ft

PROTECTIVE CASING MATERIAL TYPE:

MEASURE PROTECTIVE CASING INSIDE DIAMETER (Inches):

YES	NO
X	X
	X
	X
	X
	X

LOCK PRESENT?

LOCK FUNCTIONAL?

DID YOU REPLACE THE LOCK?

IS THERE EVIDENCE THAT THE WELL IS DOUBLE CASED? (If yes, describe below)

WELL MEASURING POINT VISIBLE?

17.74
11.90
211
Steel
Good
NA
Far near system
trenching

MEASURE WELL DEPTH FROM MEASURING POINT (Feet):

MEASURE DEPTH TO WATER FROM MEASURING POINT (Feet):

MEASURE WELL DIAMETER (Inches):

WELL CASING MATERIAL:

PHYSICAL CONDITION OF VISIBLE WELL CASING:

ATTACH ID MARKER (if well ID is confirmed) and IDENTIFY MARKER TYPE

PROXIMITY TO UNDERGROUND OR OVERHEAD UTILITIES.....

DESCRIBE ACCESS TO WELL: (Include accessibility to truck mounted rig, natural obstructions, overhead power lines, proximity to permanent structures, etc.); ADD SKETCH OF LOCATION ON BACK, IF NECESSARY.

Good → In open field

DESCRIBE WELL SETTING (For example, located in a field, in a playground, on pavement, in a garden, etc.)

AND ASSESS THE TYPE OF RESTORATION REQUIRED.

Open grassy field

IDENTIFY ANY NEARBY POTENTIAL SOURCES OF CONTAMINATION, IF PRESENT

(e.g. Gas station, salt pile, etc.):

REMARKS:

Temp 12.50 DO% 4.5
 Spec Cond 1.077 DO 0.47
 Spec Cond 0.820 Sketch

pH 6.51
 CRP -355.3

SITE NAME:

Hancock

SITE ID.:

6024697

INSPECTOR:

GUYAV

DATE/TIME:

11/30/11

WELL ID.:

ML2 L2D

MONITORING WELL FIELD INSPECTION LOG

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>

WELL VISIBLE? (If not, provide directions below)

WELL COORDINATES? NYTM X _____ NYTM Y _____

PDOP Reading from Trimble Pathfinder: _____ Satellites: _____

GPS Method (circle) Trimble And/Or Magellan

YES	NO
<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>

WELL I.D. VISIBLE?

WELL LOCATION MATCH SITE MAP? (if not, sketch actual location on back).....

WELL I.D. AS IT APPEARS ON PROTECTIVE CASING OR WELL:

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>

SURFACE SEAL PRESENT?

SURFACE SEAL COMPETENT? (If cracked, heaved etc., describe below) heaved off to

PROTECTIVE CASING IN GOOD CONDITION? (If damaged, describe below) ..well to the East ~ 1 ft.

HEADSPACE READING (ppm) AND INSTRUMENT USED.....

TYPE OF PROTECTIVE CASING AND HEIGHT OF STICKUP IN FEET (If applicable)

PROTECTIVE CASING MATERIAL TYPE:

MEASURE PROTECTIVE CASING INSIDE DIAMETER (Inches):

YES	NO
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>

LOCK PRESENT?

LOCK FUNCTIONAL?

DID YOU REPLACE THE LOCK?

IS THERE EVIDENCE THAT THE WELL IS DOUBLE CASED? (If yes, describe below)

WELL MEASURING POINT VISIBLE?

MEASURE WELL DEPTH FROM MEASURING POINT (Feet):

MEASURE DEPTH TO WATER FROM MEASURING POINT (Feet):

MEASURE WELL DIAMETER (Inches):

WELL CASING MATERIAL:

PHYSICAL CONDITION OF VISIBLE WELL CASING:

ATTACH ID MARKER (if well ID is confirmed) and IDENTIFY MARKER TYPE

PROXIMITY TO UNDERGROUND OR OVERHEAD UTILITIES.....

6.45
7.90
2"
Steel
Bad

DESCRIBE ACCESS TO WELL: (Include accessibility to truck mounted rig, natural obstructions, overhead power lines, proximity to permanent structures, etc.); ADD SKETCH OF LOCATION ON BACK, IF NECESSARY.

Goes

DESCRIBE WELL SETTING (For example, located in a field, in a playground, on pavement, in a garden, etc.)

AND ASSESS THE TYPE OF RESTORATION REQUIRED.

Open, grassy field

IDENTIFY ANY NEARBY POTENTIAL SOURCES OF CONTAMINATION, IF PRESENT

(e.g. Gas station, salt pile, etc.):

NA

REMARKS:

Well in circular shape.

Temp 11.50

Spec Cond 1.313

Sketch Spec Cond 0.781

DO% 19.7

DO 2.08

pH 6.49

ORP -341.4

SITE NAME:

Hancock

SITE ID:

60214697

INSPECTOR:

GW/AV

DATE/TIME:

11/30/11

WELL ID:

MW-125

MONITORING WELL FIELD INSPECTION LOG

WELL VISIBLE? (If not, provide directions below)

YES	NO
X	

WELL COORDINATES? NYTM X _____ NYTM Y _____

PDOP Reading from Trimble Pathfinder: _____ Satellites: _____

GPS Method (circle) Trimble And/Or Magellan

WELL I.D. VISIBLE?

YES	NO
	X
X	

WELL LOCATION MATCH SITE MAP? (if not, sketch actual location on back).....

WELL I.D. AS IT APPEARS ON PROTECTIVE CASING OR WELL:

YES	NO
X	
X	
	X

SURFACE SEAL PRESENT?

SURFACE SEAL COMPETENT? (If cracked, heaved etc., describe below)

PROTECTIVE CASING IN GOOD CONDITION? (If damaged, describe below)

Stickup casing is bent to be almost parallel to ground. Well rendered useless due to damage.

HEADSPACE READING (ppm) AND INSTRUMENT USED.....

TYPE OF PROTECTIVE CASING AND HEIGHT OF STICKUP IN FEET (If applicable)

Stick-up

PROTECTIVE CASING MATERIAL TYPE:

MEASURE PROTECTIVE CASING INSIDE DIAMETER (Inches):

YES	NO
X	
	X
	X
	X
	X

LOCK PRESENT?

LOCK FUNCTIONAL?

DID YOU REPLACE THE LOCK?

IS THERE EVIDENCE THAT THE WELL IS DOUBLE CASED? (If yes, describe below)

WELL MEASURING POINT VISIBLE?

MEASURE WELL DEPTH FROM MEASURING POINT (Feet):

MEASURE DEPTH TO WATER FROM MEASURING POINT (Feet):

MEASURE WELL DIAMETER (Inches):

WELL CASING MATERIAL:

PHYSICAL CONDITION OF VISIBLE WELL CASING:

ATTACH ID MARKER (if well ID is confirmed) and IDENTIFY MARKER TYPE

PROXIMITY TO UNDERGROUND OR OVERHEAD UTILITIES.....

DESCRIBE ACCESS TO WELL: (Include accessibility to truck mounted rig, natural obstructions, overhead power lines, proximity to permanent structures, etc.); ADD SKETCH OF LOCATION ON BACK, IF NECESSARY.

Good

DESCRIBE WELL SETTING (For example, located in a field, in a playground, on pavement, in a garden, etc.) AND ASSESS THE TYPE OF RESTORATION REQUIRED.

Open, grassy field

IDENTIFY ANY NEARBY POTENTIAL SOURCES OF CONTAMINATION, IF PRESENT (e.g. Gas station, salt pile, etc.):

NA

REMARKS:

Temp	DO %	pH
Spec Cond	DO	ORP
Spec Cond		

Sketch

SITE NAME:

Hancock

SITE ID.:

60214697

INSPECTOR:

GWAU

DATE/TIME:

11/30/11

WELL ID.:

M2-13

MONITORING WELL FIELD INSPECTION LOG

WELL VISIBLE? (If not, provide directions below)

YES	NO
X	

WELL COORDINATES? NYTM X _____ NYTM Y _____

PDOP Reading from Trimble Pathfinder: _____ Satellites: _____

GPS Method (circle) Trimble And/Or Magellan

WELL I.D. VISIBLE?

YES	NO
	X
X	

WELL LOCATION MATCH SITE MAP? (if not, sketch actual location on back).....

WELL I.D. AS IT APPEARS ON PROTECTIVE CASING OR WELL: Sewer Co.

SURFACE SEAL PRESENT?

SURFACE SEAL COMPETENT? (If cracked, heaved etc., describe below)

PROTECTIVE CASING IN GOOD CONDITION? (If damaged, describe below)

YES	NO
X	
X	
X	

HEADSPACE READING (ppm) AND INSTRUMENT USED.....

TYPE OF PROTECTIVE CASING AND HEIGHT OF STICKUP IN FEET (If applicable)

PROTECTIVE CASING MATERIAL TYPE:

MEASURE PROTECTIVE CASING INSIDE DIAMETER (Inches):

	NA
	flush

LOCK PRESENT?

LOCK FUNCTIONAL?

DID YOU REPLACE THE LOCK?

IS THERE EVIDENCE THAT THE WELL IS DOUBLE CASED? (If yes, describe below)

WELL MEASURING POINT VISIBLE?

YES	NO
X	
	X
	X
	X

MEASURE WELL DEPTH FROM MEASURING POINT (Feet):

MEASURE DEPTH TO WATER FROM MEASURING POINT (Feet):

MEASURE WELL DIAMETER (Inches):

WELL CASING MATERIAL:

PHYSICAL CONDITION OF VISIBLE WELL CASING:

ATTACH ID MARKER (if well ID is confirmed) and IDENTIFY MARKER TYPE

PROXIMITY TO UNDERGROUND OR OVERHEAD UTILITIES.....

	16.70
	8.96
	2"
	steel
	good
	NA

DESCRIBE ACCESS TO WELL: (Include accessibility to truck mounted rig, natural obstructions, overhead power lines, proximity to permanent structures, etc.); ADD SKETCH OF LOCATION ON BACK, IF NECESSARY.

Good

DESCRIBE WELL SETTING (For example, located in a field, in a playground, on pavement, in a garden, etc.)

AND ASSESS THE TYPE OF RESTORATION REQUIRED.

Open, grassy field

IDENTIFY ANY NEARBY POTENTIAL SOURCES OF CONTAMINATION, IF PRESENT

(e.g. Gas station, salt pile, etc.):

REMARKS:

Casing causes depression in pavement where H₂O pools over well

Temp

141.96

Spec cond

4.374

Spec Cond

3.555

DO%

Sketch 23-3

DO

2.33

pH 6.75

ORP -293.0

SITE NAME:

Hancock

SITE ID:

60214697

INSPECTOR:

GW/AV

DATE/TIME:

11/30/11

WELL ID:

MW-14

MONITORING WELL FIELD INSPECTION LOG

YES	NO
X	

WELL VISIBLE? (If not, provide directions below)

WELL COORDINATES? NYTM X _____ NYTM Y _____

PDOP Reading from Trimble Pathfinder: _____ Satellites: _____

GPS Method (circle) Trimble And/Or Magellan

YES	NO
X	
X	

WELL I.D. VISIBLE?

WELL LOCATION MATCH SITE MAP? (if not, sketch actual location on back).....

WELL I.D. AS IT APPEARS ON PROTECTIVE CASING OR WELL: MW-14

YES	NO
X	
X	
X	

SURFACE SEAL PRESENT?

SURFACE SEAL COMPETENT? (If cracked, heaved etc., describe below)

PROTECTIVE CASING IN GOOD CONDITION? (If damaged, describe below)

NA
3.5 ft
Steel

HEADSPACE READING (ppm) AND INSTRUMENT USED.....

TYPE OF PROTECTIVE CASING AND HEIGHT OF STICKUP IN FEET (If applicable)

PROTECTIVE CASING MATERIAL TYPE:

MEASURE PROTECTIVE CASING INSIDE DIAMETER (Inches):

YES	NO
X	
	X
	X
	X
X	

LOCK PRESENT?

LOCK FUNCTIONAL?

DID YOU REPLACE THE LOCK?

IS THERE EVIDENCE THAT THE WELL IS DOUBLE CASED? (If yes, describe below)

WELL MEASURING POINT VISIBLE?

15.09
12.44
2"
Steel
Good
NA

MEASURE WELL DEPTH FROM MEASURING POINT (Feet):

MEASURE DEPTH TO WATER FROM MEASURING POINT (Feet):

MEASURE WELL DIAMETER (Inches):

WELL CASING MATERIAL:

PHYSICAL CONDITION OF VISIBLE WELL CASING:

ATTACH ID MARKER (if well ID is confirmed) and IDENTIFY MARKER TYPE

PROXIMITY TO UNDERGROUND OR OVERHEAD UTILITIES.....

DESCRIBE ACCESS TO WELL: (Include accessibility to truck mounted rig, natural obstructions, overhead power lines, proximity to permanent structures, etc.); ADD SKETCH OF LOCATION ON BACK, IF NECESSARY.

Good

DESCRIBE WELL SETTING (For example, located in a field, in a playground, on pavement, in a garden, etc.)

AND ASSESS THE TYPE OF RESTORATION REQUIRED.

Open field

IDENTIFY ANY NEARBY POTENTIAL SOURCES OF CONTAMINATION, IF PRESENT

(e.g. Gas station, salt pile, etc.):

REMARKS:

Temp 12.63 Spec Cond 8.142 DO% 6.8 pH 6.60
 " 0.887 DO 0.70 BRP 208.6
 I - Plug sticks up so that casing does not sit snug

SITE NAME:

Hancock

SITE ID.:

60214697

INSPECTOR:

AV/GW

DATE/TIME:

11/30/11

WELL ID.:

MW-16

MONITORING WELL FIELD INSPECTION LOG

WELL VISIBLE? (If not, provide directions below)

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>

WELL COORDINATES? NYTM X _____ NYTM Y _____

PDOP Reading from Trimble Pathfinder: _____ Satellites: _____

GPS Method (circle) Trimble And/Or Magellan

WELL I.D. VISIBLE?

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>

WELL LOCATION MATCH SITE MAP? (if not, sketch actual location on back).....

WELL I.D. AS IT APPEARS ON PROTECTIVE CASING OR WELL:

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>

SURFACE SEAL PRESENT?

SURFACE SEAL COMPETENT? (If cracked, heaved etc., describe below)

PROTECTIVE CASING IN GOOD CONDITION? (If damaged, describe below)

HEADSPACE READING (ppm) AND INSTRUMENT USED.....

TYPE OF PROTECTIVE CASING AND HEIGHT OF STICKUP IN FEET (If applicable)

PROTECTIVE CASING MATERIAL TYPE:

MEASURE PROTECTIVE CASING INSIDE DIAMETER (Inches):

3'
Steel
8"

LOCK PRESENT?

LOCK FUNCTIONAL?

DID YOU REPLACE THE LOCK?

IS THERE EVIDENCE THAT THE WELL IS DOUBLE CASED? (If yes, describe below)

WELL MEASURING POINT VISIBLE?

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>

MEASURE WELL DEPTH FROM MEASURING POINT (Feet):

MEASURE DEPTH TO WATER FROM MEASURING POINT (Feet):

MEASURE WELL DIAMETER (Inches):

WELL CASING MATERIAL:

PHYSICAL CONDITION OF VISIBLE WELL CASING:

ATTACH ID MARKER (if well ID is confirmed) and IDENTIFY MARKER TYPE

PROXIMITY TO UNDERGROUND OR OVERHEAD UTILITIES.....

18.82
11.95
2"
Steel
Good
NO

DESCRIBE ACCESS TO WELL: (Include accessibility to truck mounted rig, natural obstructions, overhead power lines, proximity to permanent structures, etc.); ADD SKETCH OF LOCATION ON BACK, IF NECESSARY.

Good

DESCRIBE WELL SETTING (For example, located in a field, in a playground, on pavement, in a garden, etc.)

AND ASSESS THE TYPE OF RESTORATION REQUIRED.

Open, grassy field

IDENTIFY ANY NEARBY POTENTIAL SOURCES OF CONTAMINATION, IF PRESENT

(e.g. Gas station, salt pile, etc.):

REMARKS:

Temp 14.11

Spec Cond 1.939

DO% 2.9

pH 6.81

DO 0.30

ORP -301.0

Sketch

SITE NAME:

Hancock

SITE ID.:

60214697

INSPECTOR:

G.W./HV

DATE/TIME:

11/30/11

WELL ID.:

AW-156

MONITORING WELL FIELD INSPECTION LOG

WELL VISIBLE? (If not, provide directions below)

YES	NO
X	

WELL COORDINATES? NYTM X _____ NYTM Y _____

PDOP Reading from Trimble Pathfinder: _____ Satellites: _____

GPS Method (circle) Trimble And/Or Magellan

WELL I.D. VISIBLE?

YES	NO
	X
X	

WELL LOCATION MATCH SITE MAP? (if not, sketch actual location on back).....

WELL I.D. AS IT APPEARS ON PROTECTIVE CASING OR WELL:

YES	NO
X	
X	
X	

SURFACE SEAL PRESENT?

SURFACE SEAL COMPETENT? (If cracked, heaved etc., describe below)

PROTECTIVE CASING IN GOOD CONDITION? (If damaged, describe below)

Pushmount

HEADSPACE READING (ppm) AND INSTRUMENT USED.....

TYPE OF PROTECTIVE CASING AND HEIGHT OF STICKUP IN FEET (If applicable)

PROTECTIVE CASING MATERIAL TYPE:

MEASURE PROTECTIVE CASING INSIDE DIAMETER (Inches):

YES	NO
	X
	X
	X
X	

LOCK PRESENT?

LOCK FUNCTIONAL?

DID YOU REPLACE THE LOCK?

IS THERE EVIDENCE THAT THE WELL IS DOUBLE CASED? (If yes, describe below)

WELL MEASURING POINT VISIBLE?

MEASURE WELL DEPTH FROM MEASURING POINT (Feet):

MEASURE DEPTH TO WATER FROM MEASURING POINT (Feet):

MEASURE WELL DIAMETER (Inches):

WELL CASING MATERIAL:

PHYSICAL CONDITION OF VISIBLE WELL CASING:

ATTACH ID MARKER (if well ID is confirmed) and IDENTIFY MARKER TYPE

PROXIMITY TO UNDERGROUND OR OVERHEAD UTILITIES.....

15.30
5.20
3"
Good
few

DESCRIBE ACCESS TO WELL: (Include accessibility to truck mounted rig, natural obstructions, overhead power lines, proximity to permanent structures, etc.); ADD SKETCH OF LOCATION ON BACK, IF NECESSARY.

Good

DESCRIBE WELL SETTING (For example, located in a field, in a playground, on pavement, in a garden, etc.)

AND ASSESS THE TYPE OF RESTORATION REQUIRED.

Open, grassy field

IDENTIFY ANY NEARBY POTENTIAL SOURCES OF CONTAMINATION, IF PRESENT

(e.g. Gas station, salt pile, etc.):

NA

REMARKS:

Well depressed below grass.

2 10' N of sapling tree

Temp 13.30

Spec Cond 0.156

Sketch

Spec Cond 6.123

DO% 46.6 pH 6.57

DO 4.86 ORP -277.9

SITE NAME:

Hancock

SITE ID:

60214697

INSPECTOR:

GLW

DATE/TIME:

11/30/11

WELL ID:

MW-101

MONITORING WELL FIELD INSPECTION LOG

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>

WELL VISIBLE? (If not, provide directions below)

WELL COORDINATES? NYTM X _____ NYTM Y _____

PDOP Reading from Trimble Pathfinder: _____ Satelites: _____

GPS Method (circle) Trimble And/Or Magellan

YES	NO
<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>

WELL I.D. VISIBLE?

WELL LOCATION MATCH SITE MAP? (if not, sketch actual location on back).....

WELL I.D. AS IT APPEARS ON PROTECTIVE CASING OR WELL:

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>

SURFACE SEAL PRESENT?

SURFACE SEAL COMPETENT? (If cracked, heaved etc., describe below)

PROTECTIVE CASING IN GOOD CONDITION? (If damaged, describe below)

flushmount

HEADSPACE READING (ppm) AND INSTRUMENT USED.....

TYPE OF PROTECTIVE CASING AND HEIGHT OF STICKUP IN FEET (If applicable)

PROTECTIVE CASING MATERIAL TYPE:

MEASURE PROTECTIVE CASING INSIDE DIAMETER (Inches):

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>

LOCK PRESENT?

LOCK FUNCTIONAL?

DID YOU REPLACE THE LOCK?

IS THERE EVIDENCE THAT THE WELL IS DOUBLE CASED? (If yes, describe below)

WELL MEASURING POINT VISIBLE?

MEASURE WELL DEPTH FROM MEASURING POINT (Feet):

MEASURE DEPTH TO WATER FROM MEASURING POINT (Feet):

MEASURE WELL DIAMETER (Inches):

WELL CASING MATERIAL:

PHYSICAL CONDITION OF VISIBLE WELL CASING:

ATTACH ID MARKER (if well ID is confirmed) and IDENTIFY MARKER TYPE

PROXIMITY TO UNDERGROUND OR OVERHEAD UTILITIES.....

21.91
DTP: 5.98 DTW: 10.074
2"
sch d. 40
good
sharpie on plug
far, but
system trenching near

DESCRIBE ACCESS TO WELL: (Include accessibility to truck mounted rig, natural obstructions, overhead power lines, proximity to permanent structures, etc.); ADD SKETCH OF LOCATION ON BACK, IF NECESSARY.

in open field - good

DESCRIBE WELL SETTING (For example, located in a field, in a playground, on pavement, in a garden, etc.) AND ASSESS THE TYPE OF RESTORATION REQUIRED.

open grassy field

IDENTIFY ANY NEARBY POTENTIAL SOURCES OF CONTAMINATION, IF PRESENT

(e.g. Gas station, salt pile, etc.):

NA

REMARKS:

performed skim test on well today

Sketch

SITE NAME:

Hancock

SITE ID.:

6014697

INSPECTOR:

GW/AV

DATE/TIME:

11/30/11

WELL ID.:

M2-102

MONITORING WELL FIELD INSPECTION LOG

WELL VISIBLE? (If not, provide directions below)

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>

WELL COORDINATES? NYTM X _____ NYTM Y _____

PDOP Reading from Trimble Pathfinder: _____ Satellites: _____

GPS Method (circle) Trimble And/Or Magellan

WELL I.D. VISIBLE?

YES	NO
<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>

WELL LOCATION MATCH SITE MAP? (if not, sketch actual location on back).....

WELL I.D. AS IT APPEARS ON PROTECTIVE CASING OR WELL:

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input checked="" type="checkbox"/>	<input type="checkbox"/>

SURFACE SEAL PRESENT?

SURFACE SEAL COMPETENT? (If cracked, heaved etc., describe below)

PROTECTIVE CASING IN GOOD CONDITION? (If damaged, describe below)

HEADSPACE READING (ppm) AND INSTRUMENT USED.....

TYPE OF PROTECTIVE CASING AND HEIGHT OF STICKUP IN FEET (If applicable)

PROTECTIVE CASING MATERIAL TYPE:

MEASURE PROTECTIVE CASING INSIDE DIAMETER (Inches):

LOCK PRESENT?

LOCK FUNCTIONAL?

DID YOU REPLACE THE LOCK?

IS THERE EVIDENCE THAT THE WELL IS DOUBLE CASED? (If yes, describe below)

WELL MEASURING POINT VISIBLE?

YES	NO
<input checked="" type="checkbox"/>	<input type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>
<input type="checkbox"/>	<input checked="" type="checkbox"/>

MEASURE WELL DEPTH FROM MEASURING POINT (Feet):

MEASURE DEPTH TO WATER FROM MEASURING POINT (Feet):

MEASURE WELL DIAMETER (Inches):

WELL CASING MATERIAL:

PHYSICAL CONDITION OF VISIBLE WELL CASING:

ATTACH ID MARKER (if well ID is confirmed) and IDENTIFY MARKER TYPE

PROXIMITY TO UNDERGROUND OR OVERHEAD UTILITIES.....

22.49
8.57
2"
Sched 40
good
-
N25'

DESCRIBE ACCESS TO WELL: (Include accessibility to truck mounted rig, natural obstructions, overhead power lines, proximity to permanent structures, etc.); ADD SKETCH OF LOCATION ON BACK, IF NECESSARY.

good

DESCRIBE WELL SETTING (For example, located in a field, in a playground, on pavement, in a garden, etc.) AND ASSESS THE TYPE OF RESTORATION REQUIRED.

in grassy area

IDENTIFY ANY NEARBY POTENTIAL SOURCES OF CONTAMINATION, IF PRESENT

(e.g. Gas station, salt pile, etc.):

NA

REMARKS:

Temp 14.73

Spec Cond 2.793

DO% 1.5

pH 6.068

" " 2.243

DO 0.14

ORP -334.5

Sketch

SITE NAME:

Hanrock

SITE ID:

60214697

INSPECTOR:

GW/AV

DATE/TIME:

11/30/11

WELL ID:

MLW-108

MONITORING WELL FIELD INSPECTION LOG

WELL VISIBLE? (If not, provide directions below)

YES	NO
X	

WELL COORDINATES? NYTM X _____ NYTM Y _____

PDOP Reading from Trimble Pathfinder: _____ Satellites: _____

GPS Method (circle) Trimble And/Or Magellan

WELL I.D. VISIBLE?

YES	NO
	X
X	

WELL LOCATION MATCH SITE MAP? (if not, sketch actual location on back).....

WELL I.D. AS IT APPEARS ON PROTECTIVE CASING OR WELL:

YES	NO
X	
X	
X	

SURFACE SEAL PRESENT?

SURFACE SEAL COMPETENT? (If cracked, heaved etc., describe below)

PROTECTIVE CASING IN GOOD CONDITION? (If damaged, describe below) missing 3 bolts

HEADSPACE READING (ppm) AND INSTRUMENT USED.....

TYPE OF PROTECTIVE CASING AND HEIGHT OF STICKUP IN FEET (If applicable)

PROTECTIVE CASING MATERIAL TYPE:

MEASURE PROTECTIVE CASING INSIDE DIAMETER (Inches):

Pushmount
Steel

LOCK PRESENT?

LOCK FUNCTIONAL?

DID YOU REPLACE THE LOCK?

IS THERE EVIDENCE THAT THE WELL IS DOUBLE CASED? (If yes, describe below)

WELL MEASURING POINT VISIBLE?

YES	NO
	X
	X
	X
X	
	X

MEASURE WELL DEPTH FROM MEASURING POINT (Feet):

MEASURE DEPTH TO WATER FROM MEASURING POINT (Feet):

MEASURE WELL DIAMETER (Inches):

WELL CASING MATERIAL:

PHYSICAL CONDITION OF VISIBLE WELL CASING:

ATTACH ID MARKER (if well ID is confirmed) and IDENTIFY MARKER TYPE

PROXIMITY TO UNDERGROUND OR OVERHEAD UTILITIES.....

19.28
3.45
21.1
Steel
good

DESCRIBE ACCESS TO WELL: (Include accessibility to truck mounted rig, natural obstructions, overhead power lines, proximity to permanent structures, etc.); ADD SKETCH OF LOCATION ON BACK, IF NECESSARY.

Good

DESCRIBE WELL SETTING (For example, located in a field, in a playground, on pavement, in a garden, etc.)

AND ASSESS THE TYPE OF RESTORATION REQUIRED.

Open, grassy field

IDENTIFY ANY NEARBY POTENTIAL SOURCES OF CONTAMINATION, IF PRESENT

(e.g. Gas station, salt pile, etc.):

NA

REMARKS:

Casing filling w/ water

Sketch

SITE NAME:

Hancock

SITE ID.:

60214697

INSPECTOR:

GW/AV

DATE/TIME:

11/30/11

WELL ID.:

AW-13
NW-109

MONITORING WELL FIELD INSPECTION LOG

WELL VISIBLE? (If not, provide directions below)

YES	NO
X	

WELL COORDINATES? NYTM X _____ NYTM Y _____

PDOP Reading from Trimble Pathfinder: _____ Satellites: _____

GPS Method (circle) Trimble And/Or Magellan

WELL I.D. VISIBLE?

YES	NO
X	X

WELL LOCATION MATCH SITE MAP? (if not, sketch actual location on back).....

WELL I.D. AS IT APPEARS ON PROTECTIVE CASING OR WELL:

SURFACE SEAL PRESENT?

SURFACE SEAL COMPETENT? (If cracked, heaved etc., describe below)

PROTECTIVE CASING IN GOOD CONDITION? (If damaged, describe below)

YES	NO
X	
X	
X	

HEADSPACE READING (ppm) AND INSTRUMENT USED.....

TYPE OF PROTECTIVE CASING AND HEIGHT OF STICKUP IN FEET (If applicable)

PROTECTIVE CASING MATERIAL TYPE:

MEASURE PROTECTIVE CASING INSIDE DIAMETER (Inches):

NA
flush

LOCK PRESENT?

LOCK FUNCTIONAL?

DID YOU REPLACE THE LOCK?

IS THERE EVIDENCE THAT THE WELL IS DOUBLE CASED? (If yes, describe below)

WELL MEASURING POINT VISIBLE?

YES	NO

MEASURE WELL DEPTH FROM MEASURING POINT (Feet):

MEASURE DEPTH TO WATER FROM MEASURING POINT (Feet):

MEASURE WELL DIAMETER (Inches):

WELL CASING MATERIAL:

PHYSICAL CONDITION OF VISIBLE WELL CASING:

ATTACH ID MARKER (if well ID is confirmed) and IDENTIFY MARKER TYPE

PROXIMITY TO UNDERGROUND OR OVERHEAD UTILITIES.....

14.54
8.8 ft
21"
steel
good
tw

DESCRIBE ACCESS TO WELL: (Include accessibility to truck mounted rig, natural obstructions, overhead power lines, proximity to permanent structures, etc.); ADD SKETCH OF LOCATION ON BACK, IF NECESSARY.

Good

DESCRIBE WELL SETTING (For example, located in a field, in a playground, on pavement, in a garden, etc.)

AND ASSESS THE TYPE OF RESTORATION REQUIRED.

Open, grassy field

IDENTIFY ANY NEARBY POTENTIAL SOURCES OF CONTAMINATION, IF PRESENT

(e.g. Gas station, salt pile, etc.):

NA

REMARKS:

YST did not fit into well

Sketch

SITE NAME:

Hancock

SITE ID.:

60214697

INSPECTOR:

GW/AV

DATE/TIME:

11/30/11

WELL ID.:

RW-1

MONITORING WELL FIELD INSPECTION LOG

WELL VISIBLE? (If not, provide directions below)

YES	NO
X	

WELL COORDINATES? NYTM X _____ NYTM Y _____

PDOP Reading from Trimble Pathfinder: _____ Satellites: _____

GPS Method (circle) Trimble And/Or Magellan

WELL I.D. VISIBLE?

WELL LOCATION MATCH SITE MAP? (if not, sketch actual location on back).....

YES	NO
	X
X	

WELL I.D. AS IT APPEARS ON PROTECTIVE CASING OR WELL:

SURFACE SEAL PRESENT?

SURFACE SEAL COMPETENT? (If cracked, heaved etc., describe below)

PROTECTIVE CASING IN GOOD CONDITION? (If damaged, describe below)

YES	NO
X	
X	
X	

HEADSPACE READING (ppm) AND INSTRUMENT USED.....

TYPE OF PROTECTIVE CASING AND HEIGHT OF STICKUP IN FEET (If applicable) 2.5 ft

PROTECTIVE CASING MATERIAL TYPE:

MEASURE PROTECTIVE CASING INSIDE DIAMETER (Inches):

NA

~~Steel~~ stick-up

Steel

LOCK PRESENT?

LOCK FUNCTIONAL?

DID YOU REPLACE THE LOCK?

IS THERE EVIDENCE THAT THE WELL IS DOUBLE CASED? (If yes, describe below)

WELL MEASURING POINT VISIBLE?

YES	NO
	X
	X
	X
	X
X	

MEASURE WELL DEPTH FROM MEASURING POINT (Feet):

MEASURE DEPTH TO WATER FROM MEASURING POINT (Feet):

MEASURE WELL DIAMETER (Inches):

WELL CASING MATERIAL:

PHYSICAL CONDITION OF VISIBLE WELL CASING:

ATTACH ID MARKER (if well ID is confirmed) and IDENTIFY MARKER TYPE

PROXIMITY TO UNDERGROUND OR OVERHEAD UTILITIES.....

19.63
9.93
8"
Steel
Good
NA

DESCRIBE ACCESS TO WELL: (Include accessibility to truck mounted rig, natural obstructions, overhead power lines, proximity to permanent structures, etc.); ADD SKETCH OF LOCATION ON BACK, IF NECESSARY.

Good

DESCRIBE WELL SETTING (For example, located in a field, in a playground, on pavement, in a garden, etc.)

AND ASSESS THE TYPE OF RESTORATION REQUIRED.

Open field

IDENTIFY ANY NEARBY POTENTIAL SOURCES OF CONTAMINATION, IF PRESENT

(e.g. Gas station, salt pile, etc.):

REMARKS:

Temp 13.50 DO % 2.8 ORP -291.2

Spec Cond 1.108 DO 0.30 pH 6.71

Spec Cond 0.864 Sketch

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Appendix B

Groundwater Sampling Forms

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MONITORING WELL SAMPLE COLLECTION FORM HANCOCK ANGB - ERP SITE 15

LOCATION	Site: <u>Hancock</u>	LocID: <u>MW-11</u>	Page 1 of <u>1</u>
	Project Name: Hancock ANGB ERP Site 15	Project #: 60214697	Date: <u>6/4/13</u> Recorded By: <u>GLW</u> Checked By:
EQUIPMENT	H2O Quality Meter Type/ID #: <u>SI 600 XL</u>	Sampling Equipment: <u>Peri Pump</u>	PID Type/ID #: <u>NA</u>
	Water Level Indicator Type/ID #: <u>Geotech Interface Probe</u>	Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse	
WELL INFO	Casing I.D. (in): <u>2"</u>	Ambient PID (ppm): <u>NA</u>	Initial Depth to Water (ft): <u>10.55</u>
	Ground Condition of Well: <u>good</u>	Well Mouth PID (ppm): <u>NA</u>	Total Well Depth (ft): <u>~16.5</u>
	Approximate Pump Intake Depth (ft): <u>16'</u>		
	Remarks: <u>Begin purge @ 1535</u>		

Date (mm/dd/yy)	Time (24 hr)	Water Level (FTOC)	Volume Removed (Gals)	Pumping Rate (Lpm)	Temp. (C)	Specific Conduct. (mS/cm)	DO (mg/L)	pH	ORP (mv)	Turb. (NTU)	Pump Refill/Discharge (Seconds)	PSI (pump)	Note
6/4/13	1538	10.59	initial		12.67	1.042	1.80	7.00	-76.0	16.42			100%
	1543	10.59	0.40		10.42	0.977	0.47	6.95	-66.9	46.5			16.0
	1548	10.59	0.75		10.35	0.964	0.30	7.00	-64.7	13.9			4.3
	1553	10.59	1.0		10.44	0.964	0.23	7.02	-65.7	10.46			2.6
	1558	10.59	1.35		10.39	0.963	0.23	7.04	-71.0	10.23			2.1
	1603	10.59	1.70		10.42	0.963	0.23	7.03	-72.3	8.56			2.1
													2.0

Pumping Rate: ≤ 0.5 L/min Drawdown: < 0.33 ft Measurements: 3-5 min Stabilization: ± 0.5 C, $\pm 3\%$ conductivity, $\pm 10\%$ DO, ± 0.1 pH, ± 10 mv ORP, $\pm 10\%$ turb (≤ 10 NTU ideal) All for 3 consecutive readings

Sample ID #(s)/Time(s)	# of conds./Vol./Type	Preservative	Parameter(s)
MW11-060413 @ 1604	(3) 40 mL VOAs	HCl	VOCs by 8260B

HANCOCK AND B - ERP SITE 15

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HANCOCK ANGB - ERP SITE 15													
LOCATION		Site: <u>Hancock</u>			LocID: <u>MW-14</u>			Date: <u>6/5/13</u>			Page 1 of 1		
		Project Name: <u>Hancock ANGB ERP Site 15</u>			Project #: <u>60214697</u>			Recorded By: <u>GLW</u>			Checked By:		
EQUIPMENT		H2O Quality Meter Type/ID #: <u>YSI 600XL</u>			Sampling Equipment: <u>Peri Pump</u>			PID Type/ID #: <u>NA</u>					
		Water Level Indicator Type/ID #: <u>Geotech Interface</u>			Equipment Decon.: <u>Liquinox and Potable Wash/Potable Rinse/Distilled Rinse</u>								
WELL INFO		Casing I.D. (in): <u>2"</u>			Ambient PID (ppm): <u>NA</u>			Initial Depth to Water (ft): <u>13.60</u>					
		Ground Condition of Well: <u>good - stickup</u>			Well Mouth PID (ppm): <u>NA</u>			Total Well Depth (ft):					
		Approximate Pump Intake Depth (ft):											
		Remarks: <u>Begin purge @ 1030 - j-plug makes well too high to properly cover - Slight pink hue</u>											
Date (mm/dd/yy)	Time (24 hr)	Water Level (FTOC)	Volume Removed (Gals)	Pumping Rate (Lpm)	Temp. (C)	Specific Conduct. (mS/cm)	DO (mg/L)	pH	ORP (mv)	Turb. (NTU)	Pump Refill/Discharge (Seconds)	PSI (pump)	Note
<u>6/5/13</u>	<u>1035</u>	<u>13.65</u>	<u>initial</u>		<u>12.58</u>	<u>0.687</u>	<u>2.52</u>	<u>7.08</u>	<u>42.0</u>	<u>48.4</u>			<u>20%</u>
	<u>1040</u>	<u>13.65</u>	<u>0.20</u>		<u>12.71</u>	<u>0.678</u>	<u>2.30</u>	<u>6.85</u>	<u>48.4</u>	<u>34.2</u>			<u>23.2</u>
	<u>1045</u>	<u>13.65</u>	<u>0.40</u>		<u>12.38</u>	<u>0.671</u>	<u>1.10</u>	<u>6.87</u>	<u>44.9</u>	<u>29.3</u>			<u>12.5</u>
<u>1049</u>	<u>1050</u>	<u>Pump & batt. not strong enough to purge & thus stopped working - pull tubing up doesn't help - switch out pump & Batt - 1108 Begin purging again</u>											<u>10.2</u>
	<u>1110</u>	<u>13.67</u>	<u>0.40</u>		<u>12.70</u>	<u>0.678</u>	<u>5.05</u>	<u>7.34</u>	<u>80.0</u>	<u>130.4</u>			<u>46.2</u>
	<u>1115</u>	<u>13.68</u>	<u>0.70</u>		<u>11.69</u>	<u>0.651</u>	<u>1.04</u>	<u>7.98</u>	<u>22.4</u>	<u>27.2</u>			<u>9.4</u>
	<u>1120</u>	<u>13.68</u>	<u>0.90</u>		<u>11.70</u>	<u>0.649</u>	<u>0.79</u>	<u>7.00</u>	<u>15.2</u>	<u>15.5</u>			<u>7.2</u>
	<u>1125</u>	<u>13.68</u>	<u>1.20</u>		<u>11.57</u>	<u>0.646</u>	<u>0.52</u>	<u>7.04</u>	<u>3.3</u>	<u>11.87</u>			<u>4.8</u>
	<u>1130</u>	<u>13.68</u>	<u>1.40</u>		<u>11.61</u>	<u>0.646</u>	<u>0.47</u>	<u>7.07</u>	<u>-2.5</u>	<u>9.58</u>			<u>4.3</u>
	<u>1135</u>	<u>13.68</u>	<u>1.60</u>		<u>11.57</u>	<u>0.646</u>	<u>0.41</u>	<u>7.09</u>	<u>-8.2</u>	<u>9.57</u>			<u>3.8</u>
Pumping Rate: ≤ 0.5 L/min Drawdown: ≤ 0.33 ft Measurements: 3-5 min Stabilization: ± 0.5 C, $\pm 3\%$ conductivity, $\pm 10\%$ DO, ± 0.1 pH, ± 10 mv ORP, $\pm 10\%$ turb (≤ 10 NTU ideal) All for 3 consecutive readings													
Sample ID #(s)/Time(s)				# of conds./Vol./Type			Preservative			Parameter(s)			
<u>MW14-060513@ 1142</u>				<u>1142 0.6/2 0.42</u>			<u>HCl</u>			<u>3.9</u>			
				(3) 40 mL VOAs			HCl			VOCs by 8260B			

[illegible]

MONITORING WELL SAMPLE COLLECTION FORM

HANCOCK ANGB - ERP SITE 15

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LOCATION	Site: <u>Hancock</u>	LocID: <u>MW-17</u>	Date: <u>6/4/13</u>
	Project Name: Hancock ANGB ERP Site 15	Project #: 60214697	Recorded By: <u>ELW</u> Checked By:
EQUIPMENT	H2O Quality Meter Type/ID #: <u>YSI 600 XL</u>	Sampling Equipment: <u>Peri Pump</u>	PID Type/ID #: <u>NA</u>
	Water Level Indicator Type/ID #: <u>Geotech Interface Probe</u>	Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse	
WELL INFO	Casing I.D. (in): <u>2"</u>	Ambient PID (ppm):	Initial Depth to Water (ft): <u>11.77</u>
	Ground Condition of Well: <u>good</u>	Well Mouth PID (ppm):	Total Well Depth (ft): <u>17.07</u>
	Approximate Pump Intake Depth (ft):		
	Remarks: <u>Begin purge @ 1344</u>		

Date (mm/dd/yy)	Time (24 hr)	Water Level (FTOC)	Volume Removed (Gals)	Pumping Rate (Lpm)	Temp. (C)	Specific Conduct. (mS/cm)	DO (mg/L)	pH	ORP (mv)	Turb. (NTU)	Pump Refill/Discharge (Seconds)	PSI (pump)	Note	DO%
6/4/13	1345	12.39	initial		12.85	1.146	2.03	6.97	2.9	4.86				16.1
	1350	13.00	0.30		11.87	1.106	0.161	6.45	17.5	3.75				5.7
	1355	13.19	0.55		11.84	1.108	0.55	6.40	7.2	4.19				5.2
	1400	13.39	0.80		11.84	1.117	0.38	6.44	-35.6	4.20				3.3
	1405	13.58	1.05		12.13	1.133	0.31	6.51	-66.1	4.76				2.9
	1410	13.70	1.30		12.15	1.142	0.30	6.56	-88.2	4.97				2.8
	1415	13.81	1.55		12.26	1.146	0.30	6.58	-105.9	5.13				2.8
	1420	13.90	1.80		12.19	1.144	0.28	6.55	-114.7	5.68				2.6
	1425	13.97	2.05		12.37	1.148	0.27	6.58	-131.8	5.36				2.5
	1430	14.03	2.25		12.58	1.155	0.26	6.59	-126.7	5.79				2.5
	1435	14.05	2.45		12.30	1.145	0.26	6.59	-134.0	4.69				2.4

Pumping Rate: ≤ 0.5 L/min Drawdown: < 0.33 ft Measurements: 3-5 min Stabilization: ± 0.5 C, $\pm 3\%$ conductivity, $\pm 10\%$ DO, ± 0.1 pH, ± 10 mv ORP, $\pm 10\%$ turb (≤ 10 NTU ideal) All for 3 consecutive readings

Sample ID #(s)/Time(s)	# of conts./Vol./Type	Preservative	Parameter(s)
MW 17-060413 @ 1437	(3) 40 mL VOAs	HCl	VOCs by 8260B
4 ms/MSD			
2-060413-MS & 413-MSD			
AECUM			

HANCOCK ANGB ERP SITE 15			
LOCATION	Site:	LocID: MW-14	Date: 6/14/13
	Project Name: Hancock ANGB ERP Site 15	Project #: 60214697	Recorded By: TS Checked By:
EQUIPMENT	H2O Quality Meter Type/ID #: YSI / Lamotte	Sampling Equipment: Peristaltic	PID Type/ID #: —
	Water Level Indicator Type/ID #: Solinst	Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse	
WELL INFO	Casing I.D. (in): 2	Ambient PID (ppm): —	Initial Depth to Water (ft): 8.26
	Ground Condition of Well: Good	Well Mouth PID (ppm): —	Total Well Depth (ft): 13.83
	Approximate Pump Intake Depth (ft): 13 ft		
	Remarks:		

Date (mm/dd/yy)	Time (24 hr)	Water Level (FTOC)	Volume Removed (Gals)	Pumping Rate (Lpm)	Temp. (C)	Specific Conduct. (mS/cm)	DO (mg/L)	pH	ORP (mv)	Turb. (NTU)	Pump Refill/Discharge (Seconds)	PSI (pump)	Note
6/4/13	1345	8.26	0	0.1	10.99	1.454	1.94	7.09	13.2	78.1	—	—	—
	1350	8.27	<.25	0.1	11.65	1.455	0.70	6.85	-7.6	43.8	—	—	—
	1355	8.26	.25	0.1	12.10	1.441	.67	6.62	-8.0	37.8	—	—	—
	1400	8.26	<.5	0.1	12.30	1.407	.71	6.54	-34.7	22.4	—	—	—
	1405	8.26	.5	0.1	12.08	1.379	.62	6.50	-52.3	19.2	—	—	—
	1410	8.26	<.75	0.1	12.00	1.329	.62	6.44	-60.7	17.2	—	—	—
	1415	8.26	.75	0.1	12.10	1.312	.66	6.42	-65.9	14.04	—	—	—
	1420	8.26	1	0.1	12.19	1.294	.66	6.39	-65.8	10.02	—	—	—
	1425	8.26	<1.25	0.1	12.61	1.280	.65	6.46	-63.2	8.02	—	—	—
	1430	8.26	<1.25	0.05	12.79	1.282	.65	6.50	-67.0	6.20	—	—	—
1435	8.26	1.25	0.05	12.80	1.279	.65	6.40	-66.2	5.69	—	—	—	

Pumping Rate: <=0.5 L/min **Drawdown:** <0.33 ft **Measurements:** 3-5 min **Stabilization:** +/- 0.5 C, +/- 3% conductivity, +/- 10% DO, +/- 0.1 pH, +/-10 mv ORP, +/- 10% turb (<= 10 NTU ideal) **All for 3 consecutive readings**

[illegible]

HANCOCK AND B - ERP SITE 15

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HANCOCK ANGB - ERP SITE 15													Page 1 of 2
LOCATION		Site:		LocID:		Date:		Project Name: Hancock ANGB ERP Site 15		Project #: 60214697		Recorded By: TS Checked By:	
EQUIPMENT		H2O Quality Meter Type/ID #: YSI lamotte		Sampling Equipment: Peristaltic		PID Type/ID #: -		Water Level Indicator Type/ID #: Geotech		Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse			
WELL INFO		Casing I.D. (in): 2		Ambient PID (ppm): -		Initial Depth to Water (ft): 21.07		Ground Condition of Well: Poor		Well Mouth PID (ppm): -		Total Well Depth (ft): 10.63	
		Approximate Pump Intake Depth (ft): 20'		Remarks: Sheen + odor									
Date (mm/dd/yy)	Time (24 hr)	Water Level (FTOC)	Volume Removed (Gals)	Pumping Rate (Lpm)	Temp. (C)	Specific Conduct. (mS/cm)	DO (mg/L)	pH	ORP (mv)	Turb. (NTU)	Pump Refill/Discharge (Seconds)	PSI (pump)	Note
6/5/13	1130	10.63	0	0.15	11.87	1.075	.67	7.05	-156.5	Limit	-	-	6.3
	1135	10.65	0.25	0.15	11.90	1.068	1.86	6.94	-162.1	Limit	-	-	13.6
	1140	10.65	<.5	0.15	12.03	1.071	.91	6.89	-164.6	Limit	-	-	7.9
	1145	10.66	0.5	0.15	11.79	1.094	.90	6.72	-164.1	19.1	-	-	8.4
	1150	10.66	<.75	0.15	11.71	1.122	.72	6.71	-161.3	16.7	-	-	6.6
	1155	10.65	<1.0	0.15	11.77	1.144	.80	6.75	-165.6	12.7	-	-	7.5
	1200	10.65	1.0	0.15	11.68	1.147	.60	6.73	-167.3	13.93	-	-	5.3
	1205	10.65	1.25	0.15	11.68	1.146	0.43	6.78	-168.4	9.62	-	-	3.8
	1210	10.65	1.5	0.20	11.47	1.132	0.47	6.79	-166.6	23.9	-	-	4.3
	1215	10.65	1.75	0.20	11.65	1.134	0.42	6.79	-168.7	19.7	-	-	3.9
	1220	10.65	2.00	0.20	12.00	1.148	0.40	6.84	-172.4	12.2	-	-	3.8
Pumping Rate: <=0.5 L/min Drawdown: <0.33 ft Measurements: 3-5 min Stabilization: +/- 0.5 C, +/- 3% conductivity, +/- 10% DO, +/- 0.1 pH, +/- 10 mv ORP, +/- 10% turb (<= 10 NTU ideal) All for 3 consecutive readings													
Sample ID #(s)/Time(s)					# of conts./Vol./Type		Preservative		Parameter(s)				
MW1A-060513 collected @ 1230					(3) 40 mL VOAs		HCl		VOCs by 8260B				

Pumping Rate: ≤ 0.5 L/min **Drawdown:** < 0.33 ft **Measurements:** 3-5 min **Stabilization:** ± 0.5 C, ± 3% conductivity, ± 10% DO, ± 0.1 pH, ± 10 mv ORP, ± 10% turb (≤ 10 NTU ideal) **All for 3 consecutive readings**

HANCOCK AREA - ERP SITE 15

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HANCOCK ANGB - ERP SITE 15				Page 1 of 2
LOCATION	Site: <u>Hancock</u>	LocID: <u>MW-103</u>	Date: <u>6/5/13</u>	
	Project Name: Hancock ANGB ERP Site 15	Project #: 60214697	Recorded By: <u>GLW</u>	Checked By:
EQUIPMENT	H2O Quality Meter Type/ID #: <u>YSI 600 XL</u>	Sampling Equipment: <u>Persi Pump</u>	PID Type/ID #: <u>NA</u>	
	Water Level Indicator Type/ID #: <u>Geotech Interface</u>	Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse		
WELL INFO	Casing I.D. (in): <u>2"</u>	Ambient PID (ppm): <u>NA</u>	Initial Depth to Water (ft): <u>9.85</u>	
	Ground Condition of Well: <u>good</u>	Well Mouth PID (ppm): <u>NA</u>	Total Well Depth (ft): <u>133</u>	
	Approximate Pump Intake Depth (ft): <u>132.5</u>			
	Remarks: <u>Begin purge @ 0835; floc in well-heavy-have to keep clearing out YSI</u>			

Date (mm/dd/yy)	Time (24 hr)	Water Level (FTOC)	Volume Removed (Gals)	Pumping Rate (Lpm)	Temp. (C)	Specific Conduct. (mS/cm)	DO (mg/L)	pH	ORP (mv)	Turb. (NTU)	Pump Refill/Discharge (Seconds)	PSI (pump)	Note
6/5/13	0839	9.85	initial		11.49	1.045	3.86	7.63	-27.1	904			Do%
	0844	9.85	0.20		11.89	1.104	4.03	7.28	-48.2	150			30.5
	0849	9.85	0.40		11.43	1.038	0.26	7.03	-46.0	108.7			35.3
	0854	9.85	0.60		11.24	1.012	0.19	7.01	-46.1	78.0			2.3
	0859	missed	reading										1.8
	0904	9.85	0.90		11.52	0.951	1.00	7.04	-43.1	14.6			8.6
	0909	9.85	1.2		11.82	0.947	0.30	7.05	-46.2	9.84			2.7
	0914	9.85	1.5		11.59	0.897	0.31	7.06	-45.8	59.5			2.9
	0919	9.85	1.75		11.53	0.886	0.23	7.03	-39.3	40.6			2.1
	0924	9.85	2.0		11.31	0.881	0.16	7.04	-41.7	140			1.5
	0929	9.85	2.25		11.45	0.883	0.17	7.07	-43.6	24.7			46

Pumping Rate: ≤ 0.5 L/min **Drawdown:** < 0.33 ft **Measurements:** 3-5 min **Stabilization:** ± 0.5 C, ± 3% conductivity, ± 10% DO, ± 0.1 pH, ± 10 mv ORP, ± 10% turb (≤ 10 NTU ideal) **All for 3 consecutive readings**

[illegible]

Pumping Rate: <=0.5 L/min **Drawdown:** <0.33 ft **Measurements:** 3-5 min **Stabilization:** +/- 0.5 C, +/- 3% conductivity, +/- 10% DO, +/- 0.1 pH, +/-10 mv ORP, +/- 10% turb (<= 10 NTU ideal) **All for 3 consecutive readings**

HANCOCK ANGB ERP SITE 15			
LOCATION	Site:	LocID: MW-105	Date: 6/5/13
	Project Name: Hancock ANGB ERP Site 15	Project #: 60214697	Recorded By: TS Checked By:
EQUIPMENT	H2O Quality Meter Type/ID #: YSE / Lamotte	Sampling Equipment: Peristaltic	PID Type/ID #: —
	Water Level Indicator Type/ID #: Solinst	Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse	
WELL INFO	Casing I.D. (in): 2	Ambient PID (ppm): —	Initial Depth to Water (ft): 884
	Ground Condition of Well: good	Well Mouth PID (ppm): —	Total Well Depth (ft): 3375
	Approximate Pump Intake Depth (ft): 33'		
	Remarks:		

[illegible]

Pumping Rate: ≤0.5 L/min **Drawdown:** <0.33 ft **Measurements:** 3-5 min **Stabilization:** ±0.5 C, ±3% conductivity, ±10% DO, ±0.1 pH, ±10 mv ORP, ±10% turb (<10 NTU ideal) **All for 3 consecutive readings**

[illegible]

[illegible]

Pumping Rate: <=0.5 L/min **Drawdown:** <0.33 ft **Measurements:** 3-5 min **Stabilization:** +/- 0.5 C, +/- 3% conductivity, +/- 10% DO, +/- 0.1 pH, +/-10 mv ORP, +/- 10% turb (<= 10 NTU ideal) **All for 3 consecutive readings**

MONITORING WELL SAMPLE COLLECTION FORM

HANCOCK ANGB - ERP SITE 15

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LOCATION	Site:	LocID: <u>RW-01</u>	Date: <u>6/5/13</u>
	Project Name: Hancock ANGB ERP Site 15	Project #: 60214697	Recorded By: _____ Checked By: _____
EQUIPMENT	H2O Quality Meter Type/ID #: <u>YSI / Lamotte</u>	Sampling Equipment: <u>per. static</u>	PID Type/ID #: <u>—</u>
	Water Level Indicator Type/ID #: <u>Salinist</u>	Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse	
WELL INFO	Casing I.D. (in): <u>8</u>	Ambient PID (ppm): <u>—</u>	Initial Depth to Water (ft): <u>11.28</u>
	Ground Condition of Well: <u>good</u>	Well Mouth PID (ppm): <u>—</u>	Total Well Depth (ft): <u>19.45</u>
	Approximate Pump Intake Depth (ft): <u>19</u>		
	Remarks:		

Date (mm/dd/yy)	Time (24 hr)	Water Level (FTOC)	Volume Removed (Gals)	Pumping Rate (Lpm)	Temp. (C)	Specific Conduct. (mS/cm)	DO (mg/L)	pH	ORP (mv)	Turb. (NTU)	Pump Refill/Discharge (Seconds)	PSI (pump)	DO % Note
<u>6/5/13</u>	<u>1030</u>	<u>11.28</u>	<u>0</u>	<u>0.10</u>	<u>12.66</u>	<u>713</u>	<u>1.72</u>	<u>7.40</u>	<u>-131.2</u>	<u>13.4</u>	<u>—</u>	<u>—</u>	<u>15.8</u>
	<u>1035</u>	<u>11.28</u>	<u><.25</u>	<u>0.10</u>	<u>13.40</u>	<u>718</u>	<u>1.06</u>	<u>7.21</u>	<u>-129.6</u>	<u>11.22</u>	<u>—</u>	<u>—</u>	<u>10.2</u>
	<u>1040</u>	<u>11.28</u>	<u><.25</u>	<u>0.10</u>	<u>13.60</u>	<u>722</u>	<u>.87</u>	<u>7.10</u>	<u>-119.0</u>	<u>9.05</u>	<u>—</u>	<u>—</u>	<u>8.2</u>
	<u>1045</u>	<u>11.30</u>	<u>.25</u>	<u>0.10</u>	<u>13.34</u>	<u>714</u>	<u>.56</u>	<u>7.02</u>	<u>-127.0</u>	<u>8.80</u>	<u>—</u>	<u>—</u>	<u>5.5</u>
	<u>1050</u>	<u>11.30</u>	<u><.56</u>	<u>0.10</u>	<u>13.54</u>	<u>717</u>	<u>.49</u>	<u>6.93</u>	<u>-126.6</u>	<u>8.41</u>	<u>—</u>	<u>—</u>	<u>5.0</u>
	<u>1055</u>	<u>11.31</u>	<u>0.50</u>	<u>0.10</u>	<u>13.52</u>	<u>718</u>	<u>.53</u>	<u>6.94</u>	<u>-127.9</u>	<u>8.17</u>	<u>—</u>	<u>—</u>	<u>5.6</u>
	<u>1100</u>	<u>11.31</u>	<u><.75</u>	<u>0.10</u>	<u>13.26</u>	<u>717</u>	<u>.57</u>	<u>6.92</u>	<u>-131.3</u>	<u>8.31</u>	<u>—</u>	<u>—</u>	<u>5.5</u>

Pumping Rate: <=0.5 L/min Drawdown: <0.33 ft Measurements: 3-5 min Stabilization: +/- 0.5 C, +/- 3% conductivity, +/- 10% DO, +/- 0.1 pH, +/- 10 mv ORP, +/- 10% turb (<= 10 NTU ideal) All for 3 consecutive readings

Sample ID #(s)/Time(s)	# of conts./Vol./Type	Preservative	Parameter(s)
<u>BW01-660513 collected @ 1100</u>	(3) 40 mL VOAs	HCl	VOCs by 8260B

MONITORING WELL SAMPLE COLLECTION FORM

HANCOCK AIR BASE - ERP SITE 15

Page 1 of 1

LOCATION	Site: <u>Hancock ANGB</u>	LocID: <u>MW-11</u>	Date: <u>08/29/13</u>
	Project Name: Hancock ANGB ERP Site 15	Project #: 60214697	Recorded By: <u>RSM</u> Checked By:

EQUIPMENT	H2O Quality Meter Type/ID #: <u>YS/556/Hach 2100</u>	Sampling Equipment: <u>peri pump</u>	PID Type/ID #: <u>NA</u>
	Water Level Indicator Type/ID #: <u>interface probe</u>	Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse	

WELL INFO	Casing I.D. (in): <u>2"</u>	Ambient PID (ppm): <u>NA</u>	Initial Depth to Water (ft): <u>11.34</u>
	Ground Condition of Well: <u>Good</u>	Well Mouth PID (ppm): <u>NA</u>	Total Well Depth (ft):
	Approximate Pump Intake Depth (ft):		
	Remarks:		

Date (mm/dd/yy)	Time (24 hr)	Water Level (FTOC)	Volume Removed (Gals)	Pumping Rate (Lpm)	Temp. (C)	Specific Conduct. (mS/cm)	DO (mg/L)	pH	ORP (mv)	Turb. (NTU)	Pump Refill/Discharge (Seconds)	PSI (pump)	Note
<u>08/29/13</u>	<u>15:52</u>	<u>11.40</u>	<u>-</u>	<u>.45</u>	<u>13.92</u>	<u>1.259</u>	<u>1.60</u>	<u>7.39</u>	<u>-112.9</u>	<u>369</u>	<u>-</u>	<u>-</u>	<u>Murky</u>
	<u>15:57</u>	<u>11.40</u>	<u>.7</u>	<u>.26</u>	<u>14.56</u>	<u>1.263</u>	<u>.23</u>	<u>7.10</u>	<u>-119.7</u>	<u>78</u>	<u>-</u>	<u>-</u>	<u>Murky</u>
	<u>16:02</u>	<u>11.40</u>	<u>1.2</u>	<u>.24</u>	<u>14.90</u>	<u>1.276</u>	<u>.17</u>	<u>7.02</u>	<u>-115.0</u>	<u>40</u>	<u>-</u>	<u>-</u>	
	<u>16:07</u>	<u>11.40</u>	<u>1.5</u>	<u>.21</u>	<u>15.19</u>	<u>1.273</u>	<u>.18</u>	<u>6.99</u>	<u>-107.6</u>	<u>35</u>	<u>-</u>	<u>-</u>	
	<u>16:12</u>	<u>11.40</u>	<u>1.75</u>	<u>.28</u>	<u>15.22</u>	<u>1.261</u>	<u>.18</u>	<u>6.96</u>	<u>-107.6</u>	<u>28</u>	<u>-</u>	<u>-</u>	
	<u>16:17</u>	<u>11.40</u>	<u>2</u>	<u>.20</u>	<u>15.77</u>	<u>1.261</u>	<u>.17</u>	<u>6.94</u>	<u>-103.6</u>	<u>23</u>	<u>-</u>	<u>-</u>	
	<u>16:22</u>	<u>11.40</u>	<u>2.25</u>	<u>.20</u>	<u>16.09</u>	<u>1.270</u>	<u>.18</u>	<u>6.94</u>	<u>-100.0</u>	<u>21</u>	<u>-</u>	<u>-</u>	
	<u>16:27</u>												

Pumping Rate: <= 0.5 L/min Drawdown: < 0.33 ft Measurements: 3-5 min Stabilization: +/- 0.5 C, +/- 3% conductivity, +/- 10% DO, +/- 0.1 pH, +/- 10 mv ORP, +/- 10% turb (<= 10 NTU ideal) All for 3 consecutive readings

Sample ID #(s)/Time(s)	# of conts./Vol./Type	Preservative	Parameter(s)
<u>Begin Page @ 15:50</u> <u>Samples Collected @ 16:26</u>	(3) 40 mL VOAs	HCl	VOCs by 8260B <u>BEX only</u>

MONITORING WELL SAMPLE COLLECTION FORM

HANCOCK ANGB - ERP SITE 15

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LOCATION	Site: <u>Hancock</u>	LocID: <u>MW-14</u>	Date: <u>08/29/13</u>
	Project Name: Hancock ANGB ERP Site 15	Project #: 60214697	Recorded By: <u>RSM</u> Checked By:
EQUIPMENT	H2O Quality Meter Type/ID #: <u>YSI 556 / Hach 2100p</u>	Sampling Equipment: <u>peri pump</u>	PID Type/ID #: <u>N/A</u>
	Water Level Indicator Type/ID #:	Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse	
WELL INFO	Casing I.D. (in): <u>2"</u>	Ambient PID (ppm): <u>NA</u>	Initial Depth to Water (ft): <u>14.33</u>
	Ground Condition of Well: <u>Good</u>	Well Mouth PID (ppm):	Total Well Depth (ft):
	Approximate Pump Intake Depth (ft):		
	Remarks:		

Date (mm/dd/yy)	Time (24 hr)	Water Level (FTOC)	Volume Removed (Gals)	Pumping Rate (Lpm)	Temp. (C)	Specific Conduct. (mS/cm)	DO (mg/L)	pH	ORP (mv)	Turb. (NTU)	Pump Refill/Discharge (Seconds)	PSI (pump)	Note
08/29/13	11:49	14.42	—	.15	17.76	.960	1.13	6.82	-14.2	19	—	—	Slight pink hue
	11:54	14.42	.5	.20	15.40	.940	.47	6.85	-15.3	9	—	—	
	11:59	14.41	.72	.12	15.94	.938	.24	6.91	-32.3	4	—	—	
	12:04	14.41	1.0	.12	15.72	.937	.25	6.92	-34.9	2	—	—	
	12:09	14.41	1.25	.12	15.86	.940	.19	6.91	-34.0	2	—	—	
	12:14	14.41	1.5	.12	15.85	.940	.18	6.91	-36.2	1	—	—	
	12:19	14.41	1.75	.12	15.86	.955	.19	6.93	-43.6	1	—	—	

Pumping Rate: <=0.5 L/min Drawdown: <0.33 ft Measurements: 3-5 min Stabilization: +/- 0.5 C, +/- 3% conductivity, +/- 10% DO, +/- 0.1 pH, +/- 10 mv ORP, +/- 10% turb (<= 10 NTU ideal) All for 3 consecutive readings

Sample ID #(s)/Time(s)	# of conts./Vol./Type	Preservative	Parameter(s)
Started pumping @ 11:47 Collected samples @ 12:22	(3) 40 mL VOAs	HCl	VOCs by 8260B <u>PEX only</u>

MONITORING WELL SAMPLE COLLECTION FORM

HANCOCK AREA B - ERP SITE 15

Page 1 of 1

LOCATION	Site: <u>Hancock</u>	LocID: <u>MW-15</u>	Date: <u>8/29/13</u>
	Project Name: Hancock ANGB ERP Site 15	Project #: 60214697	Recorded By: <u>GLW</u> Checked By:
EQUIPMENT	H2O Quality Meter Type/ID #: <u>YSI 556 HACH 2100</u>	Sampling Equipment: <u>Peri. pump</u>	PID Type/ID #: <u>NA</u>
	Water Level Indicator Type/ID #: <u>interface probe</u>	Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse	
WELL INFO	Casing I.D. (in): <u>2"</u>	Ambient PID (ppm): <u>NA</u>	Initial Depth to Water (ft): <u>13.85</u>
	Ground Condition of Well: <u>good</u>	Well Mouth PID (ppm): <u>NA</u>	Total Well Depth (ft):
	Approximate Pump Intake Depth (ft):		
	Remarks: <u>sl. odor</u>		

Date (mm/dd/yy)	Time (24 hr)	Water Level (FTOC)	Volume Removed (Gals)	Pumping Rate (Lpm)	Temp. (C)	Specific Conduct. (mS/cm)	DO (mg/L)	pH	ORP (mv)	Turb. (NTU)	Pump Refill/Discharge (Seconds)	PSI (pump)	Note
08/29/13	1550	13.90	initial	1.25	15.05	0.955	4.41	6.99	-121.1	16.0	-	-	clean
	1555	13.91	0.25	1.75	14.56	0.943	0.98	7.01	-119.8	2.4	-	-	
	1600	13.91	0.50	1.75	14.31	0.936	0.55	7.03	-117.9	1.9	-	-	
	1605	13.91	0.75	1.75	14.39	0.933	0.39	7.05	-105.4	2.0	-	-	
	1610	13.91	1.0	1.75	14.56	0.927	0.35	7.07	-102.8	2.6	-	-	
	1615	13.91	1.25	1.75	14.52	0.919	0.31	7.06	-101.6	1.5	-	-	

Pumping Rate: <=0.5 L/min Drawdown: <0.33 ft Measurements: 3-5 min Stabilization: +/- 0.5 C, +/- 3% conductivity, +/- 10% DO, +/- 0.1 pH, +/- 10 mv ORP, +/- 10% turb (<= 10 NTU ideal) All for 3 consecutive readings

Sample ID #(s)/Time(s)	# of conts./Vol./Type	Preservative	Parameter(s)
Begin purge @ 1547 Collect MW15-082913 @ 1618	(3) 40 mL VOAs	HCl	VOCs by 8260B <u>BEX only</u>

MONITORING WELL SAMPLE COLLECTION FORM

HANCOCK ANGB - ERP SITE 15

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LOCATION	Site: <u>Hancock ANGB</u>	LocID: <u>mw-19</u>	Date: <u>8/29/13</u>
	Project Name: Hancock ANGB ERP Site 15	Project #: 60214697	Recorded By: <u>RSU</u> Checked By:

EQUIPMENT	H2O Quality Meter Type/ID #: <u>YSI 556 / Hack 2100p</u>	Sampling Equipment: <u>Peri. pump</u>	PID Type/ID #: <u>N/A</u>
	Water Level Indicator Type/ID #: <u>interface probe</u>	Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse	

WELL INFO	Casing I.D. (in): <u>2"</u>	Ambient PID (ppm): <u>NA</u>	Initial Depth to Water (ft): <u>8.91'</u>
	Ground Condition of Well: <u>mud</u>	Well Mouth PID (ppm): <u>NA</u>	Total Well Depth (ft):
	Approximate Pump Intake Depth (ft):		
	Remarks:		

Date (mm/dd/yy)	Time (24 hr)	Water Level (FTOC)	Volume Removed (Gals)	Pumping Rate (Lpm)	Temp. (C)	Specific Conduct. (mS/cm)	DO (mg/L)	pH	ORP (mv)	Turb. (NTU)	Pump Refill/Discharge (Seconds)	PSI (pump)	Note
08/29/13	16:45	8.95	~	.5	15.62	1.383	5.67	7.08	-84.9	46	-	-	clear
	16:50	8.95	.5	.15	15.01	1.347	.24	6.83	-93.4	19	-	-	
	16:55	8.95	.75	.18	15.86	1.384	.24	6.80	-97.1	13	-	-	
	17:00	8.95	1.0	.18	15.69	1.419	.21	6.76	-93.6	7	-	-	
	17:05	8.95	1.25	.10	15.49	1.439	.22	6.70	-90.9	5	-	-	
	17:10	8.95	1.5	.10	15.44	1.452	.20	6.68	-80.0	5	-	-	

Pumping Rate: <=0.5 L/min Drawdown: <0.33 ft Measurements: 3-5 min Stabilization: +/- 0.5 C, +/- 3% conductivity, +/- 10% DO, +/- 0.1 pH, +/- 10 mv ORP, +/- 10% turb (<= 10 NTU ideal) All for 3 consecutive readings

Sample ID #(s)/Time(s)	# of conts./Vol./Type	Preservative	Parameter(s)
<u>Begin purge @ 16:42</u> <u>1645-Collect AMB-0829B</u> <u>as ambient Blank</u> <u>Samples taken @ 17:12</u>	(3) 40 mL VOAs	HCl	VOCs by 8260B <u>BEX only</u>

MONITORING WELL SAMPLE COLLECTION FORM HANCOCK ANGB - ERP SITE 15

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LOCATION	Site: <u>Hancock</u>	LocID: <u>MW101</u>	Date: <u>8/29/13</u>
	Project Name: Hancock ANGB ERP Site 15	Project #: 60214697	Recorded By: <u>RM</u> Checked By:
EQUIPMENT	H2O Quality Meter Type/ID #: <u>SI 556, HACH 2100</u>	Sampling Equipment: <u>Perist. pump</u>	PID Type/ID #: <u>NA</u>
	Water Level Indicator Type/ID #: <u>interface probe</u>	Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse	
WELL INFO	Casing I.D. (in): <u>2" good</u>	Ambient PID (ppm): <u>NA</u>	Initial Depth to Water (ft): <u>11.39</u>
	Ground Condition of Well:	Well Mouth PID (ppm): <u>NA</u>	Total Well Depth (ft):
	Approximate Pump Intake Depth (ft):		
	Remarks:		

Date (mm/dd/yy)	Time (24 hr)	Water Level (FTOC)	Volume Removed (Gals)	Pumping Rate (Lpm)	Temp. (C)	Specific Conduct. (mS/cm)	DO (mg/L)	pH	ORP (mv)	Turb. (NTU)	Pump Refill/Discharge (Seconds)	PSI (pump)	Note
8/29/13	12:47	11.43	-	.30	16.32	1.236	1.86	7.43	-155.7	12	-	-	Submerged
	12:52	11.46	0.5	.30	15.37	1.224	.21	7.40	-157.6	5	-	-	
	12:57	11.45	.75	.30	15.62	1.264	.11	7.06	-153.1	4	-	-	
	13:02	11.47	1.5	.3	15.24	1.279	.08	6.94	-135.1	4	-	-	
	13:07	11.46	2	.3	14.96	1.278	.10	6.92	-125.5	3	-	-	
	13:12	11.46	2.5	.3	15.26	1.287	.13	6.92	-129.9	3	-	-	

Pumping Rate: <=0.5 L/min Drawdown: <0.33 ft Measurements: 3-5 min Stabilization: +/- 0.5 C, +/- 3% conductivity, +/- 10% DO, +/- 0.1 pH, +/- 10 mv ORP, +/- 10% turb (<= 10 NTU ideal) All for 3 consecutive readings

Sample ID #(s)/Time(s)	# of conts./Vol./Type	Preservative	Parameter(s)
Begin purge @ 12:45 Collect MW101-082913 @ 13:13 & Dup -082913	(3) 40 mL VOAs	HCl	VOCs by 8260B, <u>BEX only</u>

MONITORING WELL SAMPLE COLLECTION FORM

HANCOCK ANGB - ERP SITE 15

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LOCATION	Site: <u>Hancock</u>	LocID: <u>MW-103</u>	Date: <u>8/29/13</u>
	Project Name: Hancock ANGB ERP Site 15	Project #: 60214697	Recorded By: <u>GLW</u> Checked By:
EQUIPMENT	H2O Quality Meter Type/ID #: <u>YSI 556 & HACH 2100</u>	Sampling Equipment: <u>Perist pump</u>	PID Type/ID #: <u>NA</u>
	Water Level Indicator Type/ID #: <u>interface probe</u>	Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse	
WELL INFO	Casing I.D. (in): <u>2"</u>	Ambient PID (ppm): <u>NA</u>	Initial Depth to Water (ft): <u>10.41</u>
	Ground Condition of Well: <u>good</u>	Well Mouth PID (ppm): <u>NA</u>	Total Well Depth (ft):
	Approximate Pump Intake Depth (ft):		
	Remarks: <u>orange floc in well</u>		

Date (mm/dd/yy)	Time (24 hr)	Water Level (FTOC)	Volume Removed (Gals)	Pumping Rate (Lpm)	Temp. (C)	Specific Conduct. (mS/cm)	DO (mg/L)	pH	ORP (mv)	Turb. (NTU)	Pump Refill/Discharge (Seconds)	PSI (pump)	Note
08/29/13	1445	10.43	initial	200	13.32	1.183	4.69	7.10	-93.0	99.9*	-	-	
	1450	10.43	0.50	200	13.19	1.172	1.06	6.99	-89.1	99.9*	emptied	-	cup of floc &
	1455	10.43	0.90	200	13.30	1.209	1.88	7.00	-71.2	49.5	let it refill @ 1452	-	
	1500	10.43	1.25	200	12.93	1.168	0.71	6.98	-65.3	13.3	-	-	Floc stopped
	1505	10.43	1.70	200	12.72	1.141	0.46	7.00	-72.7	99.9*	-	-	Floc back
	1510	10.43	2.05	200	12.69	1.118	0.32	6.99	-80.4	99.9*	-	-	FLOC
	1515	10.43	2.50	200	12.68	1.118	0.33	6.99	-83.8	51.5	-	-	no Floc

Pumping Rate: <= 0.5 L/min Drawdown: < 0.33 ft Measurements: 3-5 min Stabilization: +/- 0.5 C, +/- 3% conductivity, +/- 10% DO, +/- 0.1 pH, +/- 10 mv ORP, +/- 10% turb (<= 10 NTU ideal) All for 3 consecutive readings

Sample ID #(s)/Time(s)	# of conts./Vol./Type	Preservative	Parameter(s)
Begin purge @ 1443 Collect MW103-082913 @ 1518	(3) 40 mL VOAs	HCl	VOCs by 8260B, BEX only

HANCOCK AREA B - ERP SITE 15

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HANCOCK ANGB-ERP SITE 15				Page 1 of 1
LOCATION	Site: <u>Hancock ANGB</u>	LocID: <u>MW-105</u>	Date: <u>8/29/13</u>	
	Project Name: Hancock ANGB ERP Site 15	Project #: 60214697	Recorded By: <u>RSM</u>	Checked By:
EQUIPMENT	H2O Quality Meter Type/ID #: <u>YSI 556 HACH 2100</u>	Sampling Equipment: <u>per pump</u>	PID Type/ID #: <u>NA</u>	
	Water Level Indicator Type/ID #: <u>interface probe</u>	Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse		
WELL INFO	Casing I.D. (in): <u>2"</u>	Ambient PID (ppm): <u>NA</u>	Initial Depth to Water (ft): <u>9.33'</u>	
	Ground Condition of Well: <u>Good</u>	Well Mouth PID (ppm): <u>NA</u>	Total Well Depth (ft):	
	Approximate Pump Intake Depth (ft):			
	Remarks:			

[illegible]

Pumping Rate: ≤ 0.5 L/min Drawdown: < 0.33 ft Measurements: 3-5 min Stabilization: ± 0.5 C, $\pm 3\%$ conductivity, $\pm 10\%$ DO, ± 0.1 pH, ± 10 mv ORP $\pm 10\%$ turb (≤ 10 NTU ideal) All for 3 consecutive readings

Sample ID #(s)/Time(s)	# of conts./Vol./Type	Preservative	Parameter(s)
Began purge @ 14:37	(3) 40 mL VOAs	HCl	VOCs by 8260B BEX only
1527-Collect MW10S-082913			

MONITORING WELL SAMPLE COLLECTION FORM

HANCOCK ANGB - ERP SITE 15

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LOCATION	Site: <u>Hancock</u>	LocID: <u>MW-106</u>	Date: <u>8/29/13</u>
	Project Name: Hancock ANGB ERP Site 15	Project #: 60214697	Recorded By: <u>GLW</u> Checked By:

EQUIPMENT	H2O Quality Meter Type/ID #: <u>YSI 556 HACH 200</u>	Sampling Equipment: <u>Peri Pump</u>	PID Type/ID #: <u>NA</u>
	Water Level Indicator Type/ID #: <u>interface probe</u>	Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse	

WELL INFO	Casing I.D. (in): <u>2"</u>	Ambient PID (ppm): <u>NA</u>	Initial Depth to Water (ft): <u>2.65</u>
	Ground Condition of Well: <u>Good - in weeds</u>	Well Mouth PID (ppm): <u>NA</u>	Total Well Depth (ft):
	Approximate Pump Intake Depth (ft):		
	Remarks: <u>No odor</u>		

Date (mm/dd/yy)	Time (24 hr)	Water Level (FTOC)	Volume Removed (Gals)	Pumping Rate (Lpm)	Temp. (C)	Specific Conduct. (mS/cm)	DO (mg/L)	pH	ORP (mv)	Turb. (NTU)	Pump Refill/Discharge (Seconds)	PSI (pump)	Note
<u>8/29/13</u>	<u>1033</u>	<u>2.66</u>	<u>initial</u>	<u>200</u>	<u>14.93</u>	<u>0.780</u>	<u>1.48</u>	<u>7.08</u>	<u>-56.9</u>	<u>24.6</u>	<u>-</u>	<u>-</u>	<u>clear</u>
	<u>1038</u>	<u>2.65</u>	<u>0.50</u>	<u>200</u>	<u>14.53</u>	<u>0.782</u>	<u>0.61</u>	<u>7.12</u>	<u>-44.3</u>	<u>6.2</u>	<u>-</u>	<u>-</u>	
	<u>1043</u>	<u>2.66</u>	<u>1.0</u>	<u>225</u>	<u>14.29</u>	<u>0.801</u>	<u>0.64</u>	<u>7.14</u>	<u>-41.5</u>	<u>3.0</u>	<u>-</u>	<u>-</u>	
	<u>1048</u>	<u>2.66</u>	<u>1.40</u>	<u>225</u>	<u>14.35</u>	<u>0.811</u>	<u>0.61</u>	<u>7.15</u>	<u>-35.9</u>	<u>1.3</u>	<u>-</u>	<u>-</u>	

Pumping Rate: <=0.5 L/min Drawdown: <0.33 ft Measurements: 3-5 min Stabilization: +/- 0.5 C, +/- 3% conductivity, +/- 10% DO, +/- 0.1 pH, +/- 10 mv ORP, +/- 10% turb (<= 10 NTU ideal) All for 3 consecutive readings

Sample ID #(s)/Time(s)	# of conts./Vol./Type	Preservative	Parameter(s)
<u>Begin purge @ 1032</u> <u>Collect MW106-082913</u> <u>@ 1053</u>	(3) 40 mL VOAs	HCl	VOCs by 8260B <u>BEX only</u>

MONITORING WELL SAMPLE COLLECTION FORM HANCOCK ANGB - ERP SITE 15

Page 1 of 1

LOCATION	Site: <u>Hancock ANGB</u>	LocID: <u>MW-112</u>	Date: <u>8/29/13</u>
	Project Name: Hancock ANGB ERP Site 15	Project #: 60214697	Recorded By: <u>RM</u> Checked By:

EQUIPMENT	H2O Quality Meter Type/ID #: <u>YS1556/021453</u>	Sampling Equipment: <u>YS1556 / Hach 2100?</u>	PID Type/ID #: <u>N/A</u>
	Water Level Indicator Type/ID #: <u>Guttek</u>	Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse	

WELL INFO	Casing I.D. (in): <u>2"</u>	Ambient PID (ppm): <u>NA</u>	Initial Depth to Water (ft): <u>2.74</u>
	Ground Condition of Well: <u>Good</u>	Well Mouth PID (ppm):	Total Well Depth (ft):
	Approximate Pump Intake Depth (ft):		
	Remarks:		

Date (mm/dd/yy)	Time (24 hr)	Water Level (FTOC)	Volume Removed (Gals)	Pumping Rate (Lpm)	Temp. (C)	Specific Conduct. (mS/cm)	DO (mg/L)	pH	ORP (mv)	Turb. (NTU)	Pump Refill/Discharge (Seconds)	PSI (pump)	Note
<u>8/29/13</u>	<u>0950</u>	<u>2.74</u>	<u>-</u>	<u>.2</u>	<u>17.30</u>	<u>1.232</u>	<u>1.72</u>	<u>5.99</u>	<u>-82.8</u>	<u>7</u>	<u>-</u>	<u>-</u>	
	<u>0956</u>	<u>2.81</u>	<u>0.60</u>	<u>.2</u>	<u>16.52</u>	<u>1.191</u>	<u>.34</u>	<u>6.44</u>	<u>-81.3</u>	<u>14</u>	<u>-</u>	<u>-</u>	<u>clear</u>
	<u>10:01</u>	<u>2.80</u>	<u>0.90</u>	<u>.175</u>	<u>16.49</u>	<u>1.180</u>	<u>.20</u>	<u>6.66</u>	<u>-90.4</u>	<u>7</u>	<u>-</u>	<u>-</u>	<u>clear</u>
	<u>10:06</u>	<u>2.81</u>	<u>1.3</u>	<u>.15</u>	<u>16.82</u>	<u>1.183</u>	<u>.30</u>	<u>6.77</u>	<u>-70.3</u>	<u>4</u>	<u>-</u>	<u>-</u>	<u>clear</u>
	<u>10:11</u>	<u>2.80</u>	<u>1.5</u>	<u>.15</u>	<u>16.66</u>	<u>1.166</u>	<u>.22</u>	<u>6.83</u>	<u>-80.6</u>	<u>5</u>	<u>-</u>	<u>-</u>	<u>"</u>
	<u>10:16</u>	<u>2.79</u>	<u>1.75</u>	<u>.15</u>	<u>16.64</u>	<u>1.158</u>	<u>.17</u>	<u>6.86</u>	<u>-90.3</u>	<u>2</u>	<u>-</u>	<u>-</u>	<u>"</u>
	<u>10:21</u>	<u>2.79</u>	<u>2.10</u>	<u>.15</u>	<u>16.76</u>	<u>1.156</u>	<u>.19</u>	<u>6.90</u>	<u>-83.9</u>	<u>2</u>	<u>-</u>	<u>-</u>	
	<u>10:26</u>	<u>2.80</u>	<u>2.35</u>	<u>.15</u>	<u>16.57</u>	<u>1.150</u>	<u>.19</u>	<u>6.91</u>	<u>-89.4</u>	<u>2</u>	<u>-</u>	<u>-</u>	<u>"</u>

Pumping Rate: <=0.5 L/min Drawdown: <0.33 ft Measurements: 3-5 min Stabilization: +/- 0.5 C, +/- 3% conductivity, +/- 10% DO, +/- 0.1 pH, +/- 10 mv ORP, +/- 10% turb (<= 10 NTU ideal) All for 3 consecutive readings

Sample ID #(s)/Time(s)	# of conts./Vol./Type	Preservative	Parameter(s)
<u>start purge @ 09:49</u> <u>Sample MW-112 @ 10:28</u>	(3) 40 mL VOAs	HCl	VOCs by 8260B <u>BEX only</u>

MONITORING WELL SAMPLE COLLECTION FORM HANCOCK ANGB - ERP SITE 15

Page 1 of 1

LOCATION	Site: <u>Hancock</u>	LocID: <u>RW-1</u>	Date: <u>8/29/13</u>
	Project Name: Hancock ANGB ERP Site 15	Project #: 60214697	Recorded By: <u>GLW</u> Checked By:
EQUIPMENT	H2O Quality Meter Type/ID #: <u>YSI 556, HACH 2100</u>	Sampling Equipment: <u>Peri Pump</u>	PID Type/ID #: <u>NA</u>
	Water Level Indicator Type/ID #: <u>interface probe</u>	Equipment Decon.: Liquinox and Potable Wash/Potable Rinse/Distilled Rinse	
WELL INFO	Casing I.D. (in): <u>8"</u>	Ambient PID (ppm): <u>NA</u>	Initial Depth to Water (ft): <u>11.83</u>
	Ground Condition of Well: <u>good</u>	Well Mouth PID (ppm): <u>NA</u>	Total Well Depth (ft):
	Approximate Pump Intake Depth (ft):		
	Remarks: <u>sl. odor - unable to purge any slower</u>		

Date (mm/dd/yy)	Time (24 hr)	Water Level (FTOC)	Volume Removed (Gals)	Pumping Rate (Lpm)	Temp. (C)	Specific Conduct. (mS/cm)	DO (mg/L)	pH	ORP (mv)	Turb. (NTU)	Pump Refill/Discharge (Seconds)	PSI (pump)	Note
08/29/13	1150	11.91	initial	225	16.42	0.999	3.94	7.16	-108.1	16.5	-	-	B1. f/ecs
	1155	11.96	0.25	200	16.95	1.011	1.20	7.09	-107.8	6.3	-	-	
	1200	11.99	0.50	200	16.80	1.008	0.68	7.13	-111.3	5.7	-	-	
	1205	12.01	0.75	175	16.61	1.001	0.46	7.12	-110.1	4.7	-	-	
	1210	12.01	1.00	175	16.71	1.007	0.38	7.16	-108.5	5.9	-	-	
	1215	12.01	1.20	175	16.51	1.003	0.36	7.14	-113.1	7.1	-	-	
	1220	12.01	1.50	175	16.77	1.008	0.33	7.16	-107.9	5.4	-	-	

Pumping Rate: <=0.5 L/min Drawdown: <0.33 ft Measurements: 3-5 min Stabilization: +/- 0.5 C, +/- 3% conductivity, +/- 10% DO, +/- 0.1 pH, +/- 10 mv ORP, +/- 10% turb. (<= 10 NTU ideal) All for 3 consecutive readings

Sample ID #(s)/Time(s)	# of conts./Vol./Type	Preservative	Parameter(s)
Begin purge @ 1147 Collect RW01-082913 @ 1223 & ms/msd	(3) 40 mL VOAs	HCl	VOCs by 8260B <u>BEX only</u>

Appendix C

Laboratory Data Packages

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Report Date:
26-Jun-13 08:36



- ☒ Final Report
☐ Re-Issued Report
☐ Revised Report

Laboratory Report

AECOM Technical Services, Inc.
40 British American Boulevard
Latham, NY 12110

Work Order: M0903
Project : Hancock ANGB
Project #: HANCOCK ANGB, 60214697

Attn: John Santacroce

<u>Laboratory ID</u>	<u>Client Sample ID</u>	<u>Matrix</u>	<u>Date Sampled</u>	<u>Date Received</u>
M0903-01	MW112-060413	Aqueous	04-Jun-13 11:32	07-Jun-13 10:50
M0903-02	MW106-060413	Aqueous	04-Jun-13 12:15	07-Jun-13 10:50
M0903-03	DUP-060413	Aqueous	04-Jun-13 00:00	07-Jun-13 10:50
M0903-04	TB-060413	Aqueous	04-Jun-13 00:00	07-Jun-13 10:50
M0903-05	MW19-060413	Aqueous	04-Jun-13 14:35	07-Jun-13 10:50
M0903-06	MW17-060413	Aqueous	04-Jun-13 14:37	07-Jun-13 10:50
M0903-07	MW15-060413	Aqueous	04-Jun-13 15:40	07-Jun-13 10:50
M0903-08	MW11-060413	Aqueous	04-Jun-13 16:04	07-Jun-13 10:50
M0903-09	MW105-060513	Aqueous	05-Jun-13 09:20	07-Jun-13 10:50
M0903-10	MW103-060513	Aqueous	05-Jun-13 09:41	07-Jun-13 10:50
M0903-11	RW01-060513	Aqueous	05-Jun-13 11:00	07-Jun-13 10:50
M0903-12	MW14-060513	Aqueous	05-Jun-13 11:42	07-Jun-13 10:50
M0903-13	MW101-060513	Aqueous	05-Jun-13 12:30	07-Jun-13 10:50

I attest that the information contained within the report has been reviewed for accuracy and checked against the quality control requirements for each method. The results relate only to the sample(s) as received. This report may not be reproduced, except in full, without written approval from Spectrum Analytical.

All applicable NELAP or USEPA CLP requirements have been met.

Spectrum Analytical (Rhode Island) is accredited under the National Environmental Laboratory Approval Program (NELAP) and DoD Environmental Laboratory Accreditation Program (ELAP), holds Organic and Inorganic contracts under the USEPA CLP Program and is certified under several states. The current list of our laboratory approvals and certifications is available on the Certifications page on our web site at www.spectrum-analytical.com.

Please contact the Laboratory or Technical Director at 401-732-3400 with any questions regarding the data contained in the laboratory report.

Department of Defense	N/A
Connecticut	PH-0153
Delaware	N/A
Florida	E87664
Maine	2007037
Massachusetts	M-RI907
New Hampshire	2631
New Jersey	RI001
New York	11522
North Carolina	581
Rhode Island	LAI00301
USDA	P330-08-00023
USEPA - ISM	EP-W-09-039
USEPA - SOM	EP-W-11-033



Authorized by:

Yihai Ding
Laboratory Director

Spectrum Analytical Inc. - North Kingstown RI -- Rhode Island Division

WorkOrder: M0903

Client ID: EARTH_NY

Project: Hancock ANGB

WO Name: Hancock ANGB

Location: HANCOCK_ANGB, HANCOCK ANGB, 60214697

Comments: Send copy of invoice to John Santacroce at Latham office.

Case:

SDG:

HC Due: 06/26/13

Fax Due:

☐

Fax Report:

PO: 60214697, 11S-14728-DC30

Report Level: LEVEL 4A

Special Program: **DoD**

EDD: ERP/MS_5

Lab Samp ID	Client Sample ID	Collection Date	Date Recv'd	Matrix	Test Code	Samp / Lab Test Comments	HF	HT	MS	SEL	Storage
M0903-01A	MW112-060413	06/04/2013 11:32	06/07/2013	Aqueous	SW8260_W	/ BEX list				Y	VOA
M0903-02A	MW106-060413	06/04/2013 12:15	06/07/2013	Aqueous	SW8260_W	/ BEX list				Y	VOA
M0903-03A	DUP-060413	06/04/2013 00:00	06/07/2013	Aqueous	SW8260_W	/ BEX list				Y	VOA
M0903-04A	TB-060413	06/04/2013 00:00	06/07/2013	Aqueous	SW8260_W	/ BEX list				Y	VOA
M0903-05A	MW19-060413	06/04/2013 14:35	06/07/2013	Aqueous	SW8260_W	/ BEX list				Y	VOA
M0903-06A	MW17-060413	06/04/2013 14:37	06/07/2013	Aqueous	SW8260_W	/ BEX list			Y	Y	VOA
M0903-07A	MW15-060413	06/04/2013 15:40	06/07/2013	Aqueous	SW8260_W	/ BEX list				Y	VOA
M0903-08A	MW11-060413	06/04/2013 16:04	06/07/2013	Aqueous	SW8260_W	/ BEX list				Y	VOA
M0903-09A	MW105-060513	06/05/2013 09:20	06/07/2013	Aqueous	SW8260_W	/ BEX list				Y	VOA
M0903-10A	MW103-060513	06/05/2013 09:41	06/07/2013	Aqueous	SW8260_W	/ BEX list				Y	VOA
M0903-11A	RW01-060513	06/05/2013 11:00	06/07/2013	Aqueous	SW8260_W	/ BEX list				Y	VOA
M0903-12A	MW14-060513	06/05/2013 11:42	06/07/2013	Aqueous	SW8260_W	/ BEX list				Y	VOA
M0903-13A	MW101-060513	06/05/2013 12:30	06/07/2013	Aqueous	SW8260_W	/ BEX list				Y	VOA

HF = Fraction logged in but all tests have been placed on hold

HT = Test logged in but has been placed on hold

Sample Transmittal Documentation

CHAIN OF CUSTODY RECORD

☒ 11 Almgren Drive
Agawam, MA 01001
(413) 789-9018

☐ 8405 Benjamin Road, Ste A
Tampa, FL 33634
(813) 888-9507

□ 646 Camp Avenue
N Kingstown, RI 02852
(401) 732-3400

Special Handling:

TAT- Ind icate Date Needed: Stand

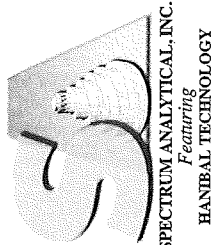
- All TATs subject to laboratory approval. Min. 24-hour notification needed for rushes.
- Samples disposed of after 60 days unless otherwise instructed.

Report To: <u>AECOM</u> <u>40 British American Blvd.</u> <u>Latham, N.Y. 12110</u>	Invoice To: <u>AECOM</u> _____ _____	Project No.: <u>60214697</u>
Telephone #: <u>518-951-2265</u> Project Mgr. <u>John Santacrose</u>	P.O. No.: _____ RQN: _____	Site Name: <u>Hancock ANG B</u>
		Location: <u>Synagogue</u> State: <u>NY</u>
		Sampler(s): <u>GLW, TS</u>

1=Na ₂ S ₂ O ₃	2=HCl	3=H ₂ SO ₄	4=HNO ₃	5=NaOH	6=Ascorbic Acid	7=CH ₃ OH
8=NaHSO ₄	9=Deionized Water	10=H ₃ PO ₄	11=	12=		
DW=Drinking Water GW=Groundwater WW=Wastewater O=Oil SW=Surface Water SO=Soil SL=Sludge A=Air X1= X2= X3=			Containers:		Class Glass Vials	
Analyses:						
2 (Xonly)						
List preservative code below:						QA/QC Reporting Notes:
						QA/QC Reporting Level
						<input type="checkbox"/> Level I <input type="checkbox"/> Level II <input type="checkbox"/> Level III <input type="checkbox"/> Level IV

[illegible]

Relinquished by:	Received by:	Date:	Time:	Temp °C	<input checked="" type="checkbox"/> EDD Format <u>ERPIMS</u> <input checked="" type="checkbox"/> E-mail to <u>John.Santacrose@AECOM.com</u>	Condition upon Receipt: Custody Seals: <input type="checkbox"/> Present <input type="checkbox"/> Broken <input type="checkbox"/> Ambient <input checked="" type="checkbox"/> Ice <input type="checkbox"/> Refrigerated <input type="checkbox"/> DI VOA Frozen <input type="checkbox"/> Soil Jar Frozen
		6/5/13	1355	14.2		
		6/5/13	1750			
		6/5/13	2100	1.2		
		6/7/13	0815			



M0903

CHAIN OF CUSTODY RECORD

Page 2 of 2

☒ 11 Almgren Drive
Agawam, MA 01001
(413) 789-9018

☐ 646 Camp Avenue
N Kingstown, RI 02852
(401) 732-3400

Special Handling:

TAT- Ind icate Date Needed: Std
· All TATs subject to laboratory approval.
· Min. 24-hour notification needed for rushes.
· Samples disposed of after 60 days unless otherwise instructed.

Report To: AECOM40 British American Blvd
Latham, NY 12110Telephone #: 518-951-2265Project Mgr. John Santacrose

1=Na₂S₂O₃ 2=HCl 3=H₂SO₄ 4=HNO₃ 5=NaOH 6=Ascorbic Acid 7=CH₃OH
8= NaHSO₄ 9= Deionized Water 10=H₃PO₄ 11= 12=

DW=Drinking Water GW=Groundwater WW=Wastewater
O=Oil SW= Surface Water SO=Soil SL=Sludge A=Air
X1= X2= X3=

Invoice To: AECOM P.O. No.: RQN: Project No.: 160214697Site Name: Hancock ANGBLocation: Syracuse State: NYSampler(s): GLW, TS

List preservative code below:

2

QA/QC Reporting Notes:

QA/QC Reporting Level

☐ Level I ☐ Level II
☐ Level III ☐ Level IV
☐ Other Std

Matrix

Time:

Date:

Sample Id:

Lab Id:

Type

of VOA Vials
of Amber Glass
of Clear Glass
of Plastic

Relinquished by:

Received by:

Date:

Time:

Temp °C

☒ EDD Format ERPIMS☒ E-mail to John.Santacrose@aecom.com

Condition upon receipt: ☐ Ambient ☐ Iced ☐ Refrigerated ☐ DI VOA Frozen ☐ Present ☐ Intact ☐ Broken ☐ Soil Jar Frozen

www.spectrum-analytical.com

Revised Feb 2013

Page 4 of 127

Received By: VB		Page 01 of 00	
Reviewed By: AED		Log-in Date 06/07/2013	
Work Order: M0903		Client Name: AECOM Technical Services, Inc.	
Project Name/Event: Hancock ANGB / Hancock ANGB			
Remarks: (1/2) Please see associated sample/extract transfer logbook pages submitted with this data package.			
		Preservation (pH)	
Lab Sample ID		HNO3	H2SO4
		HCl	NaOH
		H3PO4	VOA Matrix
1. Custody Seal(s) Present / Absent			
Intact / Broken			
2. Custody Seal Nos. N/A			
3. Traffic Reports/ Chain of Custody Records (TR/COCs) or Packing Lists Present / Absent			
4. Airbill AirBill / Sticker			
Present / Absent			
5. Airbill No. Courier N/A			
6. Sample Tags Present / Absent			
Sample Tag Numbers Listed /			
Not Listed on Chain-of-Custody			
7. Sample Condition Intact / Broken / Leaking			
8. Cooler Temperature Indicator Bottle Present / Absent			
9. Cooler Temperature 4 °C			
10. Does information on TR/COCs and sample tags agree? Yes / No			
11. Date Received at Laboratory 06/07/2013			
12. Time Received 10:50			
Sample Transfer			
Fraction (1) TVOA/VOA	Fraction (2) SVOA/PEST/ARO		
Area #	Area #		
By	By		
On	On		
IR Temp Gun ID: MT-1		VOA Matrix Key:	
Coolant Condition: ICE		US = Unpreserved Soil A = Air	
Preservative Name/Lot No:		UA = Unpreserved Aqueous H = HCl	
		M = MeOH E = Encore	
		N = NaHSO4 F = Freeze	
		See Sample Condition Notification/Corrective Action Form Yes / No	
		Rad OK Yes / No	



SPECTRUM ANALYTICAL, INC.
Featuring
HANIBAL TECHNOLOGY

*** Volatiles ***

REPORT NARRATIVE

Spectrum Analytical, Inc. Featuring Hanibal Technology, RI Division.

Client : AECOM Technical Services, Inc.

Project: Hancock ANGB

Laboratory Workorder / SDG #: M0903

SW846 8260C, VOC by GC-MS

I. SAMPLE RECEIPT

No exceptions or unusual conditions were encountered unless a Sample Condition Notification Form, or other record of communication is included with the Sample Receipt Documentation.

II. HOLDING TIMES

A. Sample Preparation:

All samples were prepared within the method-specified holding times.

B. Sample Analysis:

All samples were analyzed within the method-specified holding times.

III. METHODS

Samples were analyzed following procedures in laboratory test code:
SW846 8260C

IV. PREPARATION

Aqueous Samples were prepared following procedures in laboratory test code: SW5030B

V. INSTRUMENTATION

The following instrumentation was used

Instrument Code: V5
Instrument Type: GCMS-VOA

Description: HP6890 / HP6890
Manufacturer: Hewlett-Packard
Model: 6890 / 6890

VI. ANALYSIS

A. Calibration:

Calibrations met the method/SOP acceptance criteria.

B. Blanks:

All method blanks were within the acceptance criteria.

C. Surrogates:

Surrogate standard percent recoveries were within the QC limits.

D. Spikes:

1. Laboratory Control Spikes (LCS):

Percent recoveries for lab control samples were within the QC limits.

Replicate RPDs were within the advisory QC limits.

2. Matrix Spike / Matrix Spike Duplicate (MS/MSD):

Matrix spikes were performed on samples: MW17-060413 (M0903-06AMS) and MW17-060413 (M0903-06AMSD).

Percent recoveries were within the QC limits.

Replicate RPDs were within the advisory QC limits.

E. Internal Standards:

Internal standard peak areas were within the QC limits.

F. Dilutions:

No sample in this SDG required analysis at dilution.

G. Samples:

No other unusual occurrences were noted during sample analysis.

H. Manual Integration

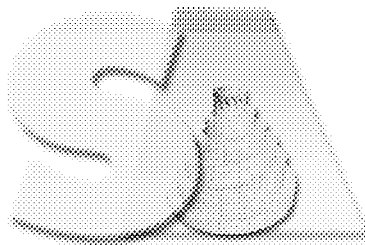
No manual integrations were performed on any sample or standard.

I certify that this data package is in compliance with the terms and conditions agreed to by the client and Spectrum, both technically and for completeness, except for the conditions noted above. Release of the data contained in this hardcopy data package has been authorized by the Laboratory Manager or designated person, as verified by the following signature.

A handwritten signature in black ink, appearing to be 'J. H. P.', is written over a horizontal line.

Signed: _____

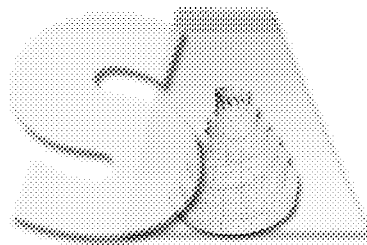
Date: _____ 6/25/2013 _____



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HANIBAL TECHNOLOGY

Data Flag/Qualifiers:

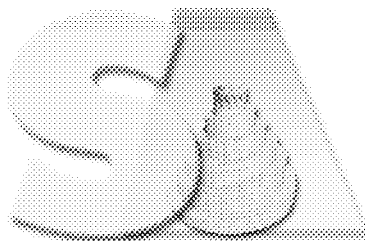
- U Not Detected. This compound was analyzed-for but not detected. For most analyses the reporting limit (lowest standard concentration) is the value listed. For Department of Defense programs, this is the Limit of Detection (LOD).
- J This flag indicates an estimated value due to either
- the compound was detected below the reporting limit, or
 - estimated concentration for Tentatively Identified Compound
- B This flag indicates the compound was also detected in the associated Method Blank. The B flag has an alternative meaning for Inorganics analyses reported using CLP ILM-type metals forms, indicating a “trace” concentration below the reporting limit and equal to or above the detection limit.
- D For Organics analysis, this flag indicates the compound concentration was obtained from a secondary dilution analysis
- E This flag indicates the compound concentration exceeded the Calibration Range. The E flag has an alternative meaning for Inorganics analyses reported using CLP metals forms, indicating an estimated concentration due to the presence of interferences, as determined by the serial dilution analysis.
- P This flag is used for pesticides/PCB/herbicide compound when there is a greater than 40% difference for detected concentration between the two GC columns used for primary and confirmation analyses. This difference typically indicates an interference, causing one value to be unusually high. The **lower** of the two values is generally reported on the Form 1, and both values reported on the Form 10.
- A Used to flag semivolatile organic Tentatively Identified Compound library search results for compounds identified as aldol condensation byproducts.
- N Used to flag results for volatile and semivolatile Organics analysis Tentatively Identified Compounds where an analyte has passed the identification criteria, and is considered to be positively identified. For Inorganics analysis the N flag indicates the matrix spike recovery falls outside of the control limit.
- * For Inorganics analysis the * flag indicates Relative Percent Difference for duplicate analyses is outside of the control limit.



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HANIBAL TECHNOLOGY

Sample ID Suffixes

- DL** Diluted analysis. The sample was diluted and reanalyzed. The DL may be followed by a digit if more than one diluted reanalysis is provided. The DL suffix is not attached to an analysis initially performed at dilution, only to reanalyses performed at dilution
- RE** Reanalysis. Appended to the client sample ID to indicate a reextraction and reanalysis or a reanalysis of the original sample extract.
- RA** Reanalysis. Appended to the laboratory sample ID indicates a reanalysis of the original sample extract.
- RX** Reextraction. Appended to the laboratory sample ID indicates a reextraction of the sample.
- MS** Matrix Spike.
- MSD** Matrix Spike Duplicate
- DUP** Duplicate analysis
- SD** Serial Dilution
- PS** Post-digestion or Post-distillation spike. For metals or inorganic analyses



SPECTRUM ANALYTICAL, INC.

Featuring

HANIBAL TECHNOLOGY

*** Sample Data ***

175 Metro Center Boulevard • Warwick, RI 02886-1755 • 401-732-3400 • FAX 401-732-3499
www.spectrum-analytical.com

1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

MW112-060413

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: M0903-01A
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V504002.D
Level: (TRACE/LOW/MED) LOW Date Received: 06/07/2013
% Moisture: not dec. Date Analyzed: 06/07/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	0.50	U	0.33	0.50	1.0
100-41-4	Ethylbenzene	0.50	U	0.35	0.50	1.0
179601-23-1	m,p-Xylene	1.0	U	0.77	1.0	1.0
95-47-6	o-Xylene	0.50	U	0.36	0.50	1.0
1330-20-7	Xylene (Total)	1.0	U	0.36	1.0	1.0

1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

MW106-060413

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: M0903-02A
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V504003.D
Level: (TRACE/LOW/MED) LOW Date Received: 06/07/2013
% Moisture: not dec. Date Analyzed: 06/07/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	0.50	U	0.33	0.50	1.0
100-41-4	Ethylbenzene	0.50	U	0.35	0.50	1.0
179601-23-1	m,p-Xylene	1.0	U	0.77	1.0	1.0
95-47-6	o-Xylene	0.50	U	0.36	0.50	1.0
1330-20-7	Xylene (Total)	1.0	U	0.36	1.0	1.0

1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.
DUP-060413

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: M0903-03A
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V504004.D
Level: (TRACE/LOW/MED) LOW Date Received: 06/07/2013
% Moisture: not dec. Date Analyzed: 06/07/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	0.50	U	0.33	0.50	1.0
100-41-4	Ethylbenzene	0.50	U	0.35	0.50	1.0
179601-23-1	m,p-Xylene	1.0	U	0.77	1.0	1.0
95-47-6	o-Xylene	0.50	U	0.36	0.50	1.0
1330-20-7	Xylene (Total)	1.0	U	0.36	1.0	1.0

1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

TB-060413

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: M0903-04A
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V504005.D
Level: (TRACE/LOW/MED) LOW Date Received: 06/07/2013
% Moisture: not dec. Date Analyzed: 06/07/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	0.50	U	0.33	0.50	1.0
100-41-4	Ethylbenzene	0.50	U	0.35	0.50	1.0
179601-23-1	m,p-Xylene	1.0	U	0.77	1.0	1.0
95-47-6	o-Xylene	0.50	U	0.36	0.50	1.0
1330-20-7	Xylene (Total)	1.0	U	0.36	1.0	1.0

1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

MW19-060413

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: M0903-05A
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V504006.D
Level: (TRACE/LOW/MED) LOW Date Received: 06/07/2013
% Moisture: not dec. Date Analyzed: 06/07/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	0.50	U	0.33	0.50	1.0
100-41-4	Ethylbenzene	0.50	U	0.35	0.50	1.0
179601-23-1	m,p-Xylene	1.0	U	0.77	1.0	1.0
95-47-6	o-Xylene	0.50	U	0.36	0.50	1.0
1330-20-7	Xylene (Total)	1.0	U	0.36	1.0	1.0

1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

MW17-060413

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: M0903-06A
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V504007.D
Level: (TRACE/LOW/MED) LOW Date Received: 06/07/2013
% Moisture: not dec. Date Analyzed: 06/07/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	0.50	U	0.33	0.50	1.0
100-41-4	Ethylbenzene	0.50	U	0.35	0.50	1.0
179601-23-1	m,p-Xylene	1.0	U	0.77	1.0	1.0
95-47-6	o-Xylene	0.50	U	0.36	0.50	1.0
1330-20-7	Xylene (Total)	1.0	U	0.36	1.0	1.0

1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

MW17-060413MS

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: M0903-06AMS
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V503995.D
Level: (TRACE/LOW/MED) LOW Date Received: 06/07/2013
% Moisture: not dec. Date Analyzed: 06/07/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	56		0.33	0.50	1.0
100-41-4	Ethylbenzene	46		0.35	0.50	1.0
179601-23-1	m,p-Xylene	92		0.77	1.0	1.0
95-47-6	o-Xylene	48		0.36	0.50	1.0
1330-20-7	Xylene (Total)	140		0.36	1.0	1.0

1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.
MW17-060413MSD

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: M0903-06AMSD
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V503996.D
Level: (TRACE/LOW/MED) LOW Date Received: 06/07/2013
% Moisture: not dec. Date Analyzed: 06/07/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	56		0.33	0.50	1.0
100-41-4	Ethylbenzene	46		0.35	0.50	1.0
179601-23-1	m,p-Xylene	93		0.77	1.0	1.0
95-47-6	o-Xylene	47		0.36	0.50	1.0
1330-20-7	Xylene (Total)	140		0.36	1.0	1.0

1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

MW15-060413

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: M0903-07A
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V504008.D
Level: (TRACE/LOW/MED) LOW Date Received: 06/07/2013
% Moisture: not dec. Date Analyzed: 06/07/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	0.50	U	0.33	0.50	1.0
100-41-4	Ethylbenzene	0.50	U	0.35	0.50	1.0
179601-23-1	m,p-Xylene	1.0	U	0.77	1.0	1.0
95-47-6	o-Xylene	0.50	U	0.36	0.50	1.0
1330-20-7	Xylene (Total)	1.0	U	0.36	1.0	1.0

1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

MW11-060413

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: M0903-08A
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V504009.D
Level: (TRACE/LOW/MED) LOW Date Received: 06/07/2013
% Moisture: not dec. Date Analyzed: 06/07/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	0.50	U	0.33	0.50	1.0
100-41-4	Ethylbenzene	0.50	U	0.35	0.50	1.0
179601-23-1	m,p-Xylene	1.0	U	0.77	1.0	1.0
95-47-6	o-Xylene	0.50	U	0.36	0.50	1.0
1330-20-7	Xylene (Total)	1.0	U	0.36	1.0	1.0

1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

MW105-060513

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: M0903-09A
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V504010.D
Level: (TRACE/LOW/MED) LOW Date Received: 06/07/2013
% Moisture: not dec. Date Analyzed: 06/07/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	0.50	U	0.33	0.50	1.0
100-41-4	Ethylbenzene	0.50	U	0.35	0.50	1.0
179601-23-1	m,p-Xylene	1.0	U	0.77	1.0	1.0
95-47-6	o-Xylene	0.50	U	0.36	0.50	1.0
1330-20-7	Xylene (Total)	1.0	U	0.36	1.0	1.0

1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

MW103-060513

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: M0903-10A
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V504011.D
Level: (TRACE/LOW/MED) LOW Date Received: 06/07/2013
% Moisture: not dec. Date Analyzed: 06/07/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	0.50	U	0.33	0.50	1.0
100-41-4	Ethylbenzene	0.50	U	0.35	0.50	1.0
179601-23-1	m,p-Xylene	1.0	U	0.77	1.0	1.0
95-47-6	o-Xylene	0.50	U	0.36	0.50	1.0
1330-20-7	Xylene (Total)	1.0	U	0.36	1.0	1.0

1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

RW01-060513

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: M0903-11A
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V504012.D
Level: (TRACE/LOW/MED) LOW Date Received: 06/07/2013
% Moisture: not dec. Date Analyzed: 06/07/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	0.50	U	0.33	0.50	1.0
100-41-4	Ethylbenzene	0.50	U	0.35	0.50	1.0
179601-23-1	m,p-Xylene	1.0	U	0.77	1.0	1.0
95-47-6	o-Xylene	0.50	U	0.36	0.50	1.0
1330-20-7	Xylene (Total)	1.0	U	0.36	1.0	1.0

1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

MW14-060513

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: M0903-12A
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V504013.D
Level: (TRACE/LOW/MED) LOW Date Received: 06/07/2013
% Moisture: not dec. Date Analyzed: 06/07/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	0.50	U	0.33	0.50	1.0
100-41-4	Ethylbenzene	0.50	U	0.35	0.50	1.0
179601-23-1	m,p-Xylene	1.0	U	0.77	1.0	1.0
95-47-6	o-Xylene	0.50	U	0.36	0.50	1.0
1330-20-7	Xylene (Total)	1.0	U	0.36	1.0	1.0

1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

MW101-060513

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: M0903-13A
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V504014.D
Level: (TRACE/LOW/MED) LOW Date Received: 06/07/2013
% Moisture: not dec. Date Analyzed: 06/08/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

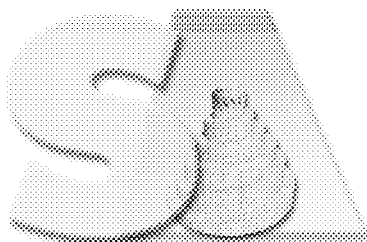
CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	0.50	U	0.33	0.50	1.0
100-41-4	Ethylbenzene	5.7		0.35	0.50	1.0
179601-23-1	m,p-Xylene	7.2		0.77	1.0	1.0
95-47-6	o-Xylene	0.50	U	0.36	0.50	1.0
1330-20-7	Xylene (Total)	7.2		0.36	1.0	1.0

1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.
LCS-72123

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: LCS-72123
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V503994.D
Level: (TRACE/LOW/MED) LOW Date Received: _____
% Moisture: not dec. Date Analyzed: 06/07/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	52		0.33	0.50	5.0
100-41-4	Ethylbenzene	42		0.35	0.50	5.0
179601-23-1	m,p-Xylene	86		0.77	1.0	5.0
95-47-6	o-Xylene	44		0.36	0.50	5.0
1330-20-7	Xylene (Total)	130		0.36	1.0	5.0



SPECTRUM ANALYTICAL, INC.
Featuring
HANIBAL TECHNOLOGY

* QC Summary *

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1A - FORM I VOA-1
VOLATILE ORGANICS ANALYSIS DATA SHEET

EPA SAMPLE NO.

MB-72123

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Matrix: (SOIL/SED/WATER) WATER Lab Sample ID: MB-72123
Sample wt/vol: 5.00 (g/mL) ML Lab File ID: V503993.D
Level: (TRACE/LOW/MED) LOW Date Received: _____
% Moisture: not dec. Date Analyzed: 06/07/2013
GC Column: DB-624 ID: 0.25 (mm) Dilution Factor: 1.0
Soil Extract Volume: _____ (uL) Soil Aliquot Volume: _____ (uL)
Purge Volume: 5.0 (mL)

CAS NO.	COMPOUND	CONCENTRATION: UG/L	Q	DL	LOD	LOQ
71-43-2	Benzene	0.50	U	0.33	0.50	1.0
100-41-4	Ethylbenzene	0.50	U	0.35	0.50	1.0
179601-23-1	m,p-Xylene	1.0	U	0.77	1.0	1.0
95-47-6	o-Xylene	0.50	U	0.36	0.50	1.0
1330-20-7	Xylene (Total)	1.0	U	0.36	1.0	1.0

4A - FORM IV VOA
VOLATILE METHOD BLANK SUMMARY

EPA SAMPLE NO.

MB-72123

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____

Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903

Lab File ID: V503993.D Lab Sample ID: MB-72123

Instrument ID: V5

Matrix: (SOIL/SED/WATER) WATER Date Analyzed: 06/07/2013

Level: (TRACE or LOW/MED) LOW Time Analyzed: 15:14

GC Column: DB-624 ID: 0.25 (mm) Heated Purge: (Y/N) N

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	TIME ANALYZED
01	LCS-72123	LCS-72123	V503994.D	15:39
02	MW17-060413M S	M0903-06AMS	V503995.D	16:05
03	MW17-060413M SD	M0903-06AMSD	V503996.D	16:30
04	MW112-060413	M0903-01A	V504002.D	19:04
05	MW106-060413	M0903-02A	V504003.D	19:30
06	DUP-060413	M0903-03A	V504004.D	19:56
07	TB-060413	M0903-04A	V504005.D	20:22
08	MW19-060413	M0903-05A	V504006.D	20:48
09	MW17-060413	M0903-06A	V504007.D	21:13
10	MW15-060413	M0903-07A	V504008.D	21:39
11	MW11-060413	M0903-08A	V504009.D	22:05
12	MW105-060513	M0903-09A	V504010.D	22:30
13	MW103-060513	M0903-10A	V504011.D	22:56
14	RW01-060513	M0903-11A	V504012.D	23:21
15	MW14-060513	M0903-12A	V504013.D	23:47
16	MW101-060513	M0903-13A	V504014.D	0:12

COMMENTS:

2B - FORM II VOA-2
WATER VOLATILE DEUTERATED MONITORING COMPOUND RECOVERY

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
 Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
 Level: (TRACE or LOW) LOW

	EPA SAMPLE NO.	VDMC1 (DBFM) #	VDMC2 (DCE) #	VDMC3 (TOL) #	VDMC4 (BFB) #				TOT OUT
01	MB-72123	106	101	93	104				0
02	LCS-72123	104	99	93	108				0
03	MW17-060413M S	107	94	93	106				0
04	MW17-060413M SD	106	99	92	106				0
05	MW112-060413	107	102	94	106				0
06	MW106-060413	103	99	94	105				0
07	DUP-060413	105	96	94	105				0
08	TB-060413	104	96	92	102				0
09	MW19-060413	105	98	94	105				0
10	MW17-060413	104	100	93	105				0
11	MW15-060413	104	98	91	104				0
12	MW11-060413	107	95	93	103				0
13	MW105-060513	106	98	94	105				0
14	MW103-060513	106	99	94	103				0
15	RW01-060513	107	97	94	106				0
16	MW14-060513	106	99	93	102				0
17	MW101-060513	107	99	95	105				0

QC LIMITS
 (85-115)
 (70-120)
 (85-120)
 (75-120)

VDMC1 (DBFM) Dibromofluoromethane
 VDMC2 (DCE) = 1,2-Dichloroethane-d4
 VDMC3 (TOL) = Toluene-d8
 VDMC4 (BFB) = Bromofluorobenzene
 # Column to be used to flag recovery values
 * Values outside of contract required QC limits

som13.06.03.A

5A - FORM V VOA
VOLATILE ORGANIC INSTRUMENT
PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

EPA SAMPLE NO.

BFB5V

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Lab File ID: V503800.D BFB Injection Date: 05/30/2013
Instrument ID: V5 BFB Injection Time: 12:22
GC Column: DB-624 ID: 0.25 (mm)

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	22.4
75	30.0 - 60.0% of mass 95	44.0
95	Base peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	6.8
173	Less than 2.0% of mass 174	0.0 (0.0)1
174	Greater than 50.0% of mass 95	74.6
175	5.0 - 9.0% of mass 174	4.0 (5.3)1
176	95.0 - 101.0% of mass 174	72.2 (96.9)1
177	5.0 - 9.0% of mass 176	4.8 (6.7)2

1 - Value is % mass 174

2 - Value is % mass 176

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	VSTD0015V	VSTD0015V	V503801.D	05/30/2013	12:48
02	VSTD0055V	VSTD0055V	V503802.D	05/30/2013	13:14
03	VSTD0205V	VSTD0205V	V503803.D	05/30/2013	13:39
04	VSTD0505V	VSTD0505V	V503804.D	05/30/2013	14:05
05	VSTD1005V	VSTD1005V	V503805.D	05/30/2013	14:32
06	VSTD2005V	VSTD2005V	V503806.D	05/30/2013	14:58
07	VICV0505V	VICV0505V	V503807.D	05/30/2013	15:46

5A - FORM V VOA
VOLATILE ORGANIC INSTRUMENT
PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

EPA SAMPLE NO.

BFB5

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Lab File ID: V503990.D BFB Injection Date: 06/07/2013
Instrument ID: V5 BFB Injection Time: 13:58
GC Column: DB-624 ID: 0.25 (mm)

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	24.8
75	30.0 - 60.0% of mass 95	44.8
95	Base peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	6.3
173	Less than 2.0% of mass 174	0.0 (0.0)1
174	Greater than 50.0% of mass 95	72.2
175	5.0 - 9.0% of mass 174	5.8 (8.0)1
176	95.0 - 101.0% of mass 174	69.8 (96.7)1
177	5.0 - 9.0% of mass 176	4.4 (6.3)2

1 - Value is % mass 174

2 - Value is % mass 176

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	VSTD050D5	VSTD050D5	V503991.D	06/07/2013	14:23
02	MB-72123	MB-72123	V503993.D	06/07/2013	15:14
03	LCS-72123	LCS-72123	V503994.D	06/07/2013	15:39
04	MW17-060413MS	M0903-06AMS	V503995.D	06/07/2013	16:05
05	MW17-060413MSD	M0903-06AMSD	V503996.D	06/07/2013	16:30
06	MW112-060413	M0903-01A	V504002.D	06/07/2013	19:04
07	MW106-060413	M0903-02A	V504003.D	06/07/2013	19:30
08	DUP-060413	M0903-03A	V504004.D	06/07/2013	19:56
09	TB-060413	M0903-04A	V504005.D	06/07/2013	20:22
10	MW19-060413	M0903-05A	V504006.D	06/07/2013	20:48
11	MW17-060413	M0903-06A	V504007.D	06/07/2013	21:13
12	MW15-060413	M0903-07A	V504008.D	06/07/2013	21:39
13	MW11-060413	M0903-08A	V504009.D	06/07/2013	22:05
14	MW105-060513	M0903-09A	V504010.D	06/07/2013	22:30
15	MW103-060513	M0903-10A	V504011.D	06/07/2013	22:56
16	RW01-060513	M0903-11A	V504012.D	06/07/2013	23:21
17	MW14-060513	M0903-12A	V504013.D	06/07/2013	23:47
18	MW101-060513	M0903-13A	V504014.D	06/08/2013	0:12

8A - FORM VIII VOA
VOLATILE INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
 Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
 GC Column: DB-624 ID: 0.25 (mm) Init. Calib. Date(s): 05/30/2013 05/30/2013
 EPA Sample No.(VSTD#####): VSTD050D5 Date Analyzed: 06/07/2013
 Lab File ID (Standard): V503991.D Time Analyzed: 14:23
 Instrument ID: V5 Heated Purge: (Y/N) N

	IS1 (S1)		IS2 (S2)		IS3 (S3)	
	AREA #	RT #	AREA #	RT #	AREA #	RT #
12 HOUR STD	793954	5.604	665170	9.077	311190	12.237
UPPER LIMIT	1587908	6.104	1330340	9.577	622380	12.737
LOWER LIMIT	396977	5.104	332585	8.577	155595	11.737
EPA SAMPLE NO.						
01 MB-72123	779985	5.605	649391	9.078	281655	12.238
02 LCS-72123	788569	5.599	658540	9.072	306088	12.244
03 MW17-060413M S	793004	5.611	665941	9.084	312009	12.244
04 MW17-060413M SD	790614	5.600	667620	9.073	304729	12.245
05 MW112-060413	796720	5.605	652480	9.078	284647	12.238
06 MW106-060413	801398	5.602	658057	9.076	280142	12.235
07 DUP-060413	776334	5.606	637969	9.079	274792	12.239
08 TB-060413	778639	5.600	642363	9.073	274583	12.245
09 MW19-060413	778963	5.604	630329	9.078	278498	12.237
10 MW17-060413	762182	5.600	623305	9.073	264311	12.244
11 MW15-060413	751663	5.599	625719	9.073	264861	12.244
12 MW11-060413	737897	5.605	616471	9.079	261716	12.238
13 MW105-060513	746710	5.605	620570	9.078	269318	12.238
14 MW103-060513	735924	5.602	614291	9.075	257071	12.235

IS1 () = Fluorobenzene

IS2 () = Chlorobenzene-d5

IS3 () = 1,4-Dichlorobenzene-d4

AREA UPPER LIMIT = 200% (Low-Medium Volatiles) and 140% (Trace Volatiles) of
internal standard area

AREA LOWER LIMIT = 50% (Low-Medium Volatiles) and 60% (Trace Volatiles) of
internal standard area

RT UPPER LIMIT = +0.50 (Low-Medium Volatiles) and +0.33 (Trace Volatiles)
minutes of internal standard RT

RT LOWER LIMIT = -0.50 (Low-Medium Volatiles) and -0.33 (Trace Volatiles)
minutes of internal standard RT

Column used to flag values outside contract required QC limits with an asterisk.

8A - FORM VIII VOA
VOLATILE INTERNAL STANDARD AREA AND RETENTION TIME SUMMARY

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____

Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903

GC Column: DB-624 ID: 0.25 (mm) Init. Calib. Date(s): 05/30/2013 05/30/2013

EPA Sample No.(VSTD#####): VSTD050D5 Date Analyzed: 06/07/2013

Lab File ID (Standard): V503991.D Time Analyzed: 14:23

Instrument ID: V5 Heated Purge: (Y/N) N

		IS1 (S1)		IS2 (S2)		IS3 (S3)	
		AREA	# RT #	AREA	# RT #	AREA	# RT #
	12 HOUR STD	793954	5.604	665170	9.077	311190	12.237
	UPPER LIMIT	1587908	6.104	1330340	9.577	622380	12.737
	LOWER LIMIT	396977	5.104	332585	8.577	155595	11.737
	EPA SAMPLE NO.						
15	RW01-060513	737920	5.605	604449	9.079	260408	12.238
16	MW14-060513	728580	5.600	604231	9.073	255556	12.244
17	MW101-060513	732159	5.605	612002	9.079	272002	12.238

IS1 () = Fluorobenzene

IS2 () = Chlorobenzene-d5

IS3 () = 1,4-Dichlorobenzene-d4

AREA UPPER LIMIT = 200% (Low-Medium Volatiles) and 140% (Trace Volatiles) of internal standard area

AREA LOWER LIMIT = 50% (Low-Medium Volatiles) and 60% (Trace Volatiles) of internal standard area

RT UPPER LIMIT = +0.50 (Low-Medium Volatiles) and +0.33 (Trace Volatiles) minutes of internal standard RT

RT LOWER LIMIT = -0.50 (Low-Medium Volatiles) and -0.33 (Trace Volatiles) minutes of internal standard RT

Column used to flag values outside contract required QC limits with an asterisk.

3A - FORM III VOA-1
WATER VOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
 Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
 Matrix Spike - EPA Sample No.: MW17-060413 Level: (TRACE or LOW) LOW

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE CONCENTRATION (ug/L)	MS CONCENTRATION (ug/L)	MS %REC	#	QC. LIMITS REC.
Benzene	50.0000	0.0000	55.5099	111		80-120
Ethylbenzene	50.0000	0.0000	46.3680	93		75-125
m,p-Xylene	100.0000	0.0000	92.3896	92		75-130
o-Xylene	50.0000	0.0000	47.6515	95		80-120
Xylene (Total)	150.0000	0.0000	140.0411	93		81-121

COMPOUND	SPIKE ADDED (ug/L)	MSD CONCENTRATION (ug/L)	MSD %REC	#	%RPD	#	QC LIMITS	
							RPD	REC.
Benzene	50.0000	55.7700	112		0		0-40	80-120
Ethylbenzene	50.0000	46.4063	93		0		0-40	75-125
m,p-Xylene	100.0000	92.9766	93		1		0-40	75-130
o-Xylene	50.0000	46.5698	93		2		0-40	80-120
Xylene (Total)	150.0000	139.5464	93		0		0-40	81-121

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

RPD: 0 out of 5 outside limits

Spike Recovery: 0 out of 10 outside limits

COMMENTS: _____

3 - FORM III
WATER LABORATORY CONTROL
SAMPLE RECOVERY

EPA SAMPLE NO.

LCS-72123

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Lab Sample ID: LCS-72123 LCS Lot No.: _____
Date Extracted: 06/07/2013 Date Analyzed (1): 06/07/2013

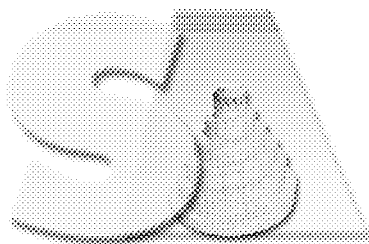
COMPOUND	SPIKE ADDED	SAMPLE CONCENTRATION	LCS CONCENTRATION	LCS %REC	#	QC. LIMITS REC.
Benzene	50.0000	0.0000	52.3818	105		80 - 120
Ethylbenzene	50.0000	0.0000	42.4445	85		75 - 125
m,p-Xylene	100.0000	0.0000	86.4939	86		75 - 130
o-Xylene	50.0000	0.0000	43.8419	88		80 - 120
Xylene (Total)	150.0000	0.0000	130.3358	87		81 - 121

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

Spike Recovery: 0 out of 5 outside limits

COMMENTS:



SPECTRUM ANALYTICAL, INC.
Featuring
HANIBAL TECHNOLOGY

*** Standards Data ***

175 Metro Center Boulevard · Warwick, RI 02886-1755 · 401-732-3400 · FAX 401-732-3499
www.spectrum-analytical.com

Lab Name: Spectrum Analytical, Inc.

Contract:

Lab Code: MITKEM

Case No.: M0903

SAS No.:

SDG No.: SM0903

Instrument ID: V5

Calibration Date(s): 05/30/2013

05/30/2013

Heated Purge: (Y/N) N

Calibration Times: 12:48

14:58

Purge Volume: 5

(mL)

GC Column: DB-624

ID: 0.25

(mm)

Length: 30

(mm)

LAB FILE ID: RRF005 = <u>V5O3802.D</u> RRF020 = <u>V5O3803.D</u> RRF050 = <u>V5O3804.D</u> RRF100 = <u>V5O3805.D</u> RRF200 = <u>V5O3806.D</u>												
RRF001 = <u>V5O3801.D</u>												
COMPOUND	RRF005	RRF020	RRF050	RRF100	RRF200	RRF001						
Benzene	1.623	1.632	1.544	1.990	1.888	1.689					1.728	10.0
Ethylbenzene	0.688	0.698	0.672	0.887	0.871	0.760					0.763	12.5
m,p-Xylene	0.872	0.891	0.815	1.087	1.039	0.895					0.933	11.3
o-Xylene	0.860	0.858	0.839	1.075	1.055	0.857					0.924	11.9
Xylene (Total)	0.868	0.880	0.823	1.083	1.045	0.883					0.930	11.4

Lab Code: MITKEM

Case No.: M0903

SAS No.:

SDG No.: SM0903

Instrument ID: V5

Calibration Date(s): 05/30/2013

05/30/2013

Heated Purge: (Y/N) N

Calibration Times: 12:48

14:58

Purge Volume: 5

(mL)

GC Column: DB-624

ID: 0.25

(mm)

Length: 30

(mm)

LAB FILE ID: RRF005 = <u>V5O3802.D</u> RRF020 = <u>V5O3803.D</u> RRF050 = <u>V5O3804.D</u> RRF100 = <u>V5O3805.D</u> RRF200 = <u>V5O3806.D</u>									
RRF001 = <u>V5O3801.D</u>									
COMPOUND	RRF005	RRF020	RRF050	RRF100	RRF200	RRF001		RRF	% RSD
Dibromofluoromethane	0.315	0.311	0.316	0.302	0.307	0.320		0.312	2.1
1,2-Dichloroethane-d4	0.072	0.068	0.067	0.069	0.069	0.072		0.070	2.9
Toluene-d8	1.295	1.293	1.258	1.330	1.275	1.296		1.291	1.9
Bromofluorobenzene	0.520	0.521	0.525	0.515	0.507	0.521		0.518	1.2

7A - FORM VII VOA-1
VOLATILE CONTINUING CALIBRATION DATA

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____

Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903

Instrument ID: V5 Calibration Date: 05/30/2013 Time: 15:46

Lab File ID: V503807.D Init. Calib. Date(s): 05/30/2013 05/30/2013

EPA Sample No.(VSTD####) VICV0505V Init. Calib. Time(s): 12:48 14:58

Heated Purge: (Y/N) N GC Column: DB-624 ID: 0.25 (mm) Length: 30 (m)

Purge Volume: 5.0 (mL)

COMPOUND	RRF	RRF050	MIN RRF	%D	MAX %D
Benzene	1.728	1.504	0.500	-13.0	20.0
Ethylbenzene	0.763	0.658	0.100	-13.7	20.0
m,p-Xylene	0.933	0.808	0.100	-13.5	20.0
o-Xylene	0.924	0.806	0.300	-12.8	20.0
Xylene (Total)	0.930	0.807	0.000	-13.2	20.0

7B - FORM VII VOA-2
VOLATILE CONTINUING CALIBRATION DATA

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____

Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903

Instrument ID: V5 Calibration Date: 05/30/2013 Time: 15:46

Lab File ID: V503807.D Init. Calib. Date(s): 05/30/2013 05/30/2013

EPA Sample No.(VSTD####) VICV0505V Init. Calib. Time(s): 12:48 14:58

Heated Purge: (Y/N) N GC Column: DB-624 ID: 0.25 (mm) Length: 30 (m)

Purge Volume: 5.0 (mL)

COMPOUND	RRF	RRF050	MIN RRF	%D	MAX %D
Dibromofluoromethane	0.312	0.310	0.100	-0.6	20.0
1,2-Dichloroethane-d4	0.070	0.067	0.100	-3.3	20.0
Toluene-d8	1.291	1.272	0.100	-1.5	20.0
Bromofluorobenzene	0.518	0.508	0.100	-1.9	20.0

7A - FORM VII VOA-1
VOLATILE CONTINUING CALIBRATION DATA

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____

Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903

Instrument ID: V5 Calibration Date: 06/07/2013 Time: 14:23

Lab File ID: V503991.D Init. Calib. Date(s): 05/30/2013 05/30/2013

EPA Sample No.(VSTD####) VSTD050D5 Init. Calib. Time(s): 12:48 14:58

Heated Purge: (Y/N) N GC Column: DB-624 ID: 0.25 (mm) Length: 30 (m)

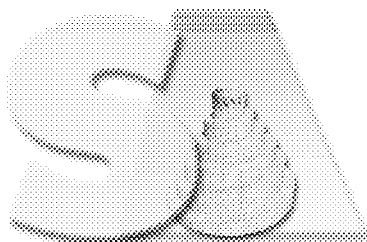
Purge Volume: 5.0 (mL)

COMPOUND	RRF	RRF050	MIN RRF	%D	MAX %D
Benzene	1.728	1.885	0.500	9.1	20.0
Ethylbenzene	0.763	0.694	0.100	-9.0	20.0
m,p-Xylene	0.933	0.849	0.100	-9.1	20.0
o-Xylene	0.924	0.863	0.300	-6.6	20.0
Xylene (Total)	0.930	0.853	0.000	-8.3	20.0

7B - FORM VII VOA-2
VOLATILE CONTINUING CALIBRATION DATA

Lab Name: SPECTRUM ANALYTICAL, INC. Contract: _____
Lab Code: MITKEM Case No.: M0903 Mod. Ref No.: _____ SDG No.: SM0903
Instrument ID: V5 Calibration Date: 06/07/2013 Time: 14:23
Lab File ID: V503991.D Init. Calib. Date(s): 05/30/2013 05/30/2013
EPA Sample No.(VSTD####) VSTD050D5 Init. Calib. Time(s): 12:48 14:58
Heated Purge: (Y/N) N GC Column: DB-624 ID: 0.25 (mm) Length: 30 (m)
Purge Volume: 5.0 (mL)

COMPOUND	RRF	RRF050	MIN RRF	%D	MAX %D
Dibromofluoromethane	0.312	0.325	0.100	4.3	20.0
1,2-Dichloroethane-d4	0.070	0.065	0.100	-6.0	20.0
Toluene-d8	1.291	1.200	0.100	-7.1	20.0
Bromofluorobenzene	0.518	0.553	0.100	6.6	20.0



SPECTRUM ANALYTICAL, INC.
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*** Raw Data ***

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www.spectrum-analytical.com

Spectrum Analytical, Inc. RI Division

Data file : \\avogadro\organics\V5.I\130530.B\V503800.D
 Lab Smp Id: BFB5V Client Smp ID: BFB5V
 Inj Date : 30-MAY-2013 12:22
 Operator : WL SRC: WL Inst ID: V5.i
 Smp Info : 2UL,BFB5V,BFB5V
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130530.B\bfb8260.m
 Meth Date : 30-May-2013 14:16 wluo Quant Type: ISTD
 Cal Date : 21-MAR-2011 17:17 Cal File: V5M6831.D
 Als bottle: 2 QC Sample: BFB
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: all.sub
 Target Version: 4.14 Sample Matrix: WATER
 Processing Host: TARGET103

Concentration Formula: Amt * DF * Uf * Vf * VI * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vf	1.000	Volumetric correction factor
VI	1.000	Injection Volume
Cpnd Variable		Local Compound Variable

CONCENTRATIONS									
				ON-COL		FINAL			
RT	EXP RT	REL RT	MASS	RESPONSE	(ug/L)	(ug/L)	TARGET	RANGE	RATIO
=====	=====	=====	=====	=====	=====	=====	=====	=====	=====
1 bfb				CAS #: 460-00-4					
10.635	11.000	(0.000)	95	104568			0.00-	100.00	100.00
10.635	11.000	(0.000)	50	23408			15.00-	40.00	22.39
10.635	11.000	(0.000)	75	46024			30.00-	60.00	44.01
10.635	11.000	(0.000)	96	7126			5.00-	9.00	6.81
10.635	11.000	(0.000)	173	0	0.0	0.0	0.00-	2.00	0.00
10.635	11.000	(0.000)	174	77960			50.00-	100.00	74.55
10.635	11.000	(0.000)	175	4144			5.00-	9.00	5.32
10.635	11.000	(0.000)	176	75544			95.00-	101.00	96.90
10.635	11.000	(0.000)	177	5030			5.00-	9.00	6.66

Date : 30-MAY-2013 12:22

Client ID: BFB5V

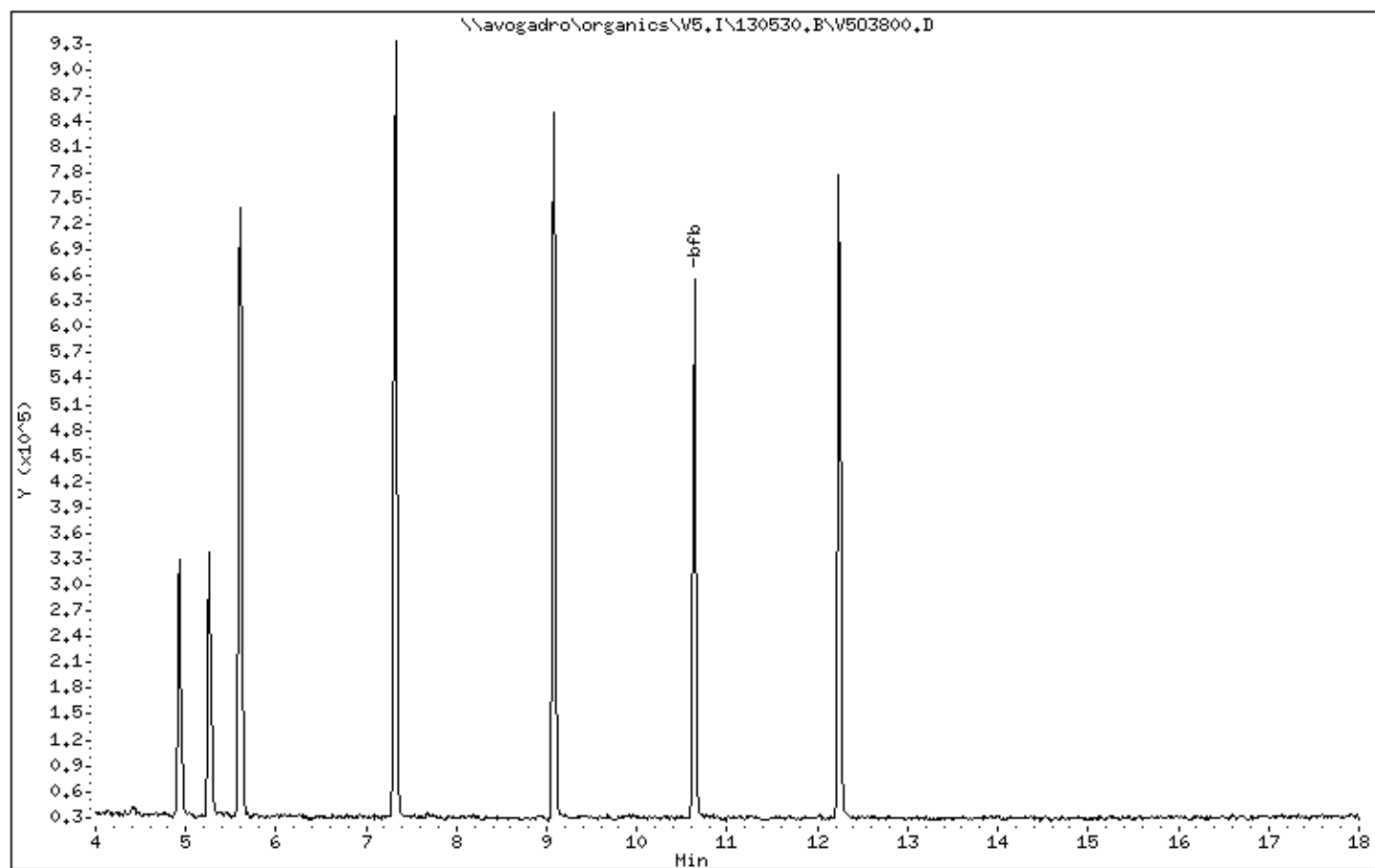
Instrument: V5.i

Sample Info: 2UL,BFB5V,BFB5V

Operator: WL SRC: WL

Column phase: DB-624

Column diameter: 0.25



Date : 30-MAY-2013 12:22

Client ID: BFB5V

Instrument: V5.i

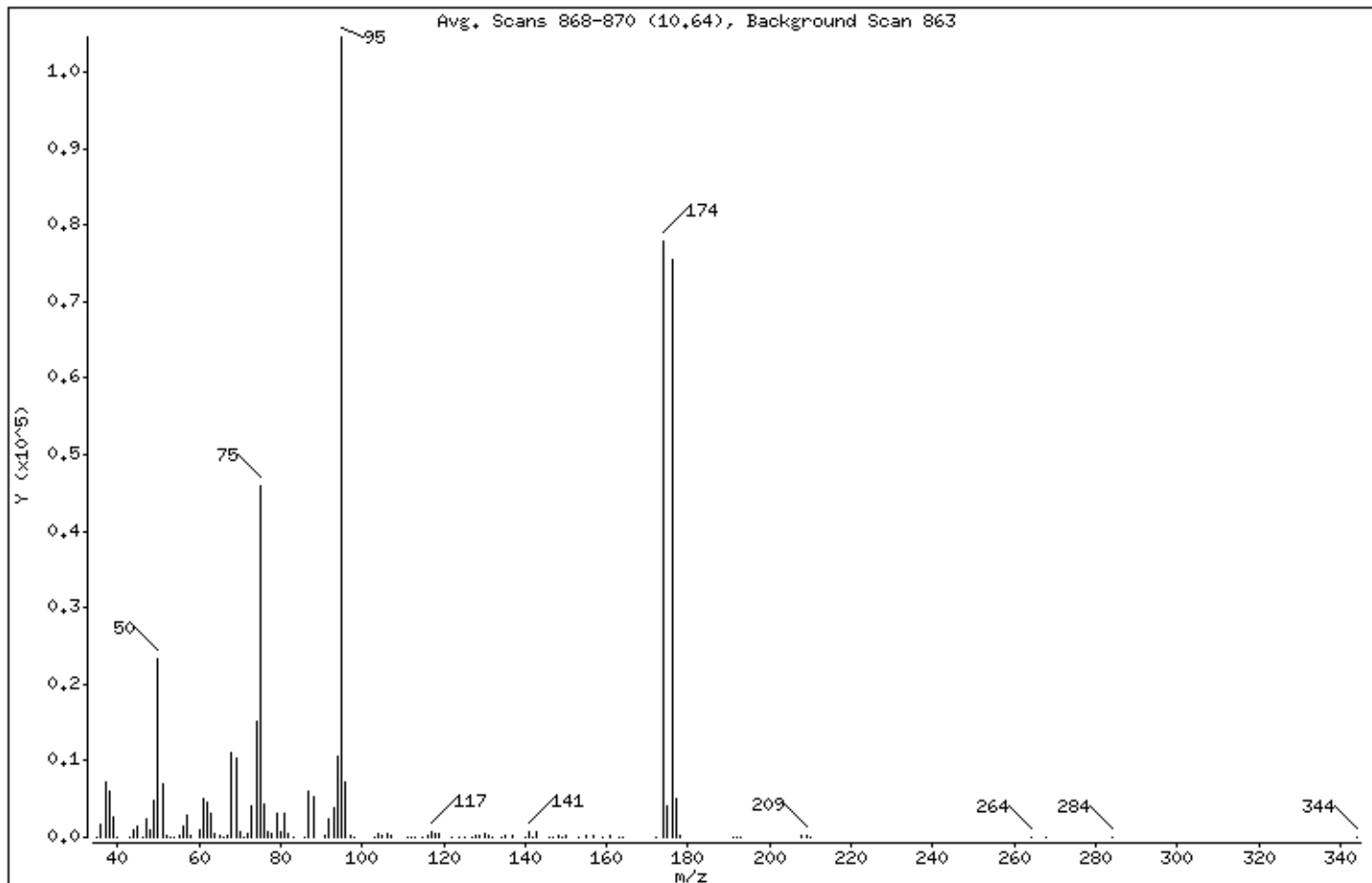
Sample Info: 2UL,BFB5V,BFB5V

Operator: WL SRC: WL

Column phase: DB-624

Column diameter: 0.25

1 bfb



m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
95	Base Peak, 100% relative abundance	100.00
50	15.00 - 40.00% of mass 95	22.39
75	30.00 - 60.00% of mass 95	44.01
96	5.00 - 9.00% of mass 95	6.81
173	Less than 2.00% of mass 174	0.00 (0.00)
174	50.00 - 100.00% of mass 95	74.55
175	5.00 - 9.00% of mass 174	3.96 (5.32)
176	95.00 - 101.00% of mass 174	72.24 (96.90)
177	5.00 - 9.00% of mass 176	4.81 (6.66)

Date : 30-MAY-2013 12:22

Client ID: BFB5V

Instrument: V5.i

Sample Info: 2UL,BFB5V,BFB5V

Operator: WL SRC: WL

Column phase: DB-624

Column diameter: 0.25

Data File: V503800.D

Spectrum: Avg. Scans 868-870 (10.64), Background Scan 863

Location of Maximum: 95.00

Number of points: 114

m/z	Y	m/z	Y	m/z	Y	m/z	Y

35.00	37	67.00	290	104.00	425	147.00	85
36.00	1715	68.00	11022	105.00	274	148.00	212
37.00	7302	69.00	10416	106.00	464	149.00	73
38.00	6081	70.00	822	107.00	154	150.00	121
39.00	2568	71.00	57	111.00	34	153.00	45

40.00	50	72.00	515	112.00	55	155.00	173
43.00	80	73.00	3988	113.00	52	157.00	186
44.00	876	74.00	15198	115.00	115	159.00	97
45.00	1428	75.00	46024	116.00	324	161.00	164
46.00	111	76.00	4221	117.00	701	163.00	1

47.00	2428	77.00	784	118.00	415	164.00	39
48.00	1030	78.00	540	119.00	377	172.00	118
49.00	4814	79.00	3202	122.00	82	174.00	77960
50.00	23408	80.00	793	124.00	47	175.00	4144
51.00	6980	81.00	3204	125.00	37	176.00	75544

52.00	305	82.00	570	127.00	35	177.00	5030
53.00	104	83.00	41	128.00	303	178.00	225
54.00	13	86.00	55	129.00	259	191.00	110
55.00	237	87.00	6004	130.00	362	192.00	93
56.00	1489	88.00	5398	131.00	250	193.00	16

57.00	2959	91.00	244	132.00	41	208.00	160
58.00	303	92.00	2499	134.00	49	209.00	262
60.00	981	93.00	3950	135.00	238	210.00	37
61.00	4988	94.00	10523	137.00	150	264.00	81
62.00	4562	95.00	104568	140.00	50	268.00	37

63.00	3199	96.00	7126	141.00	723	284.00	41
64.00	463	97.00	136	142.00	109	344.00	34
65.00	158	98.00	38	143.00	704		
66.00	98	103.00	34	146.00	93		

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130530.B\V503801.D
 Lab Smp Id: VSTD0015V Client Smp ID: VSTD0015V
 Inj Date : 30-MAY-2013 12:48
 Operator : WL SRC: WL Inst ID: V5.i
 Smp Info : 5ML,VSTD0015V,VSTD0015V
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130530.B\v5_8260W.m
 Meth Date : 03-Jun-2013 13:35 wluo Quant Type: ISTD
 Cal Date : 30-MAY-2013 12:48 Cal File: V503801.D
 Als bottle: 3 Calibration Sample, Level: 6
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: FULL.sub
 Target Version: 4.14
 Processing Host: TARGET103

Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG						AMOUNTS	
		MASS	RT	EXP RT	REL RT	RESPONSE	CAL-AMT (ug/L)	ON-COL (ug/L)
1 Dichlorodifluoromethane	85		1.603	1.583 (0.286)		9020	1.00000	1
3 Chloromethane	50		1.708	1.734 (0.305)		15957	1.00000	1
4 Vinyl Chloride	62		1.836	1.850 (0.328)		12648	1.00000	1
5 Bromomethane	94		2.138	2.141 (0.382)		7914	1.00000	1
6 Chloroethane	64		2.242	2.233 (0.400)		7606	1.00000	1
7 Trichlorofluoromethane	101		2.440	2.454 (0.436)		9497	1.00000	1
8 Ethanol	46		2.637	2.617 (0.471)		5822	500.000	110
10 Acrolein	56		2.823	2.814 (0.504)		9008	1.00000	5
13 Acetone	58		2.951	2.942 (0.527)		2618	1.00000	2
15 Carbon Disulfide	76		3.102	3.105 (0.554)		37439	1.00000	1
16 Acetonitrile	41		3.218	3.221 (0.575)		31924	1.00000	10
17 Allyl Chloride	39		3.218	3.221 (0.575)		12573	1.00000	1(T)
18 Methyl Acetate	43		3.241	3.232 (0.579)		15460	1.00000	1
19 Methylene Chloride	84		3.334	3.325 (0.595)		10114	1.00000	1
20 tert-Butanol	59		3.439	3.430 (0.614)		2297	1.00000	3
22 trans-1,2-Dichloroethene	96		3.578	3.569 (0.639)		8124	1.00000	1
23 Methyl tert-butyl ether	73		3.578	3.581 (0.639)		19675	1.00000	1
24 1,1-Dichloroethane	63		3.950	3.941 (0.705)		14895	1.00000	1
25 Vinyl acetate	43		4.008	3.999 (0.716)		30094	1.00000	1
26 Diisopropyl Ether	45		4.020	4.011 (0.718)		36721	1.00000	1
27 2-Chloro-1,3-Butadiene	53		4.043	4.034 (0.722)		11382	1.00000	1
28 Ethyl tert-butyl ether	59		4.345	4.348 (0.776)		24800	1.00000	1

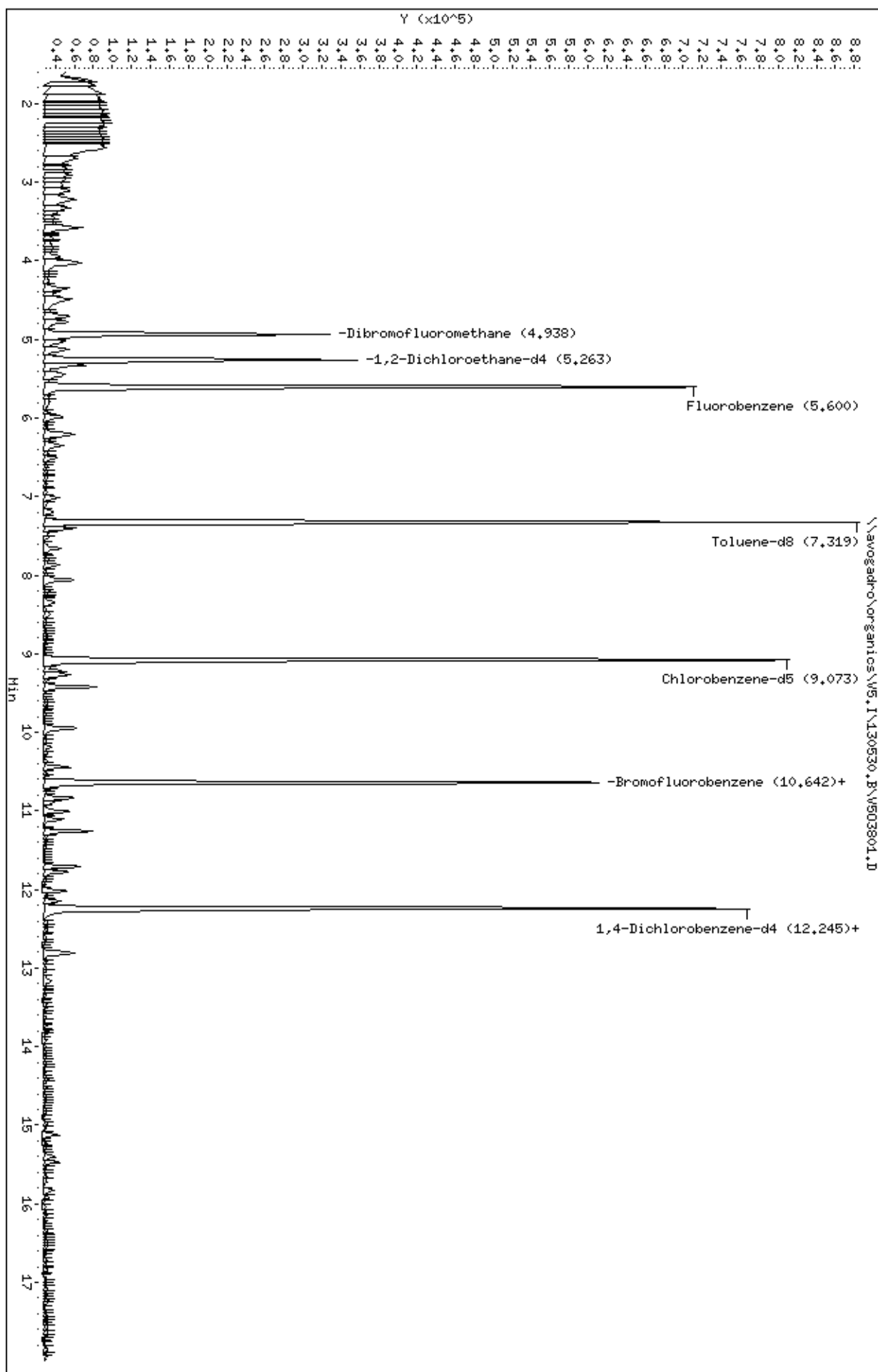
Compounds	QUANT SIG						AMOUNTS	
	MASS	RT	EXP RT	REL RT	RESPONSE		CAL-AMT (ug/L)	ON-COL (ug/L)
=====	=====	=====	=====	=====	=====		=====	=====
29 cis-1,2-Dichloroethene	96	4.484	4.475 (0.801)		8548		1.00000	1
30 2,2-Dichloropropane	77	4.484	4.487 (0.801)		7956		1.00000	1
31 2-Butanone	72	4.496	4.487 (0.803)		1170		1.00000	1(T)
35 Tetrahydrofuran	72	4.775	4.754 (0.853)		2087		1.00000	2
36 Chloroform	83	4.786	4.777 (0.855)		12579		1.00000	1
\$ 37 Dibromofluoromethane	113	4.937	4.928 (0.882)		241902		1.00000	52
38 1,1,1-Trichloroethane	97	4.972	4.975 (0.888)		9767		1.00000	1
40 1,1-Dichloropropene	110	5.135	5.126 (0.917)		3494		1.00000	1
41 Carbon Tetrachloride	117	5.135	5.138 (0.917)		6909		1.00000	1
\$ 43 1,2-Dichloroethane-d4	102	5.263	5.265 (0.940)		54612		1.00000	52
* 47 Fluorobenzene	96	5.599	5.602 (1.000)		756380		50.0000	
48 Trichloroethene	130	5.983	5.986 (1.068)		6662		1.00000	1
49 Methylcyclohexane	83	6.203	6.195 (1.108)		8337		1.00000	1
50 1,2-Dichloropropane	63	6.215	6.218 (1.110)		8694		1.00000	1
52 Methyl Methacrylate	69	6.343	6.346 (1.133)		9823		1.00000	2
53 1,4-Dioxane	88	6.389	6.357 (1.141)		1416		1.00000	20(T)
M 55 1,2-Dichloroethene (Total)	96				16672		1.00000	2
56 2-Chloroethyl vinyl ether	63	6.761	6.846 (1.207)		142		1.00000	1(T)
\$ 59 Toluene-d8	98	7.319	7.322 (0.806)		689005		1.00000	50
63 1,1,2-Trichloroethane	97	7.876	7.867 (1.407)		7516		1.00000	1
65 1,3-Dichloropropane	76	8.074	8.077 (0.889)		10612		1.00000	1
* 69 Chlorobenzene-d5	117	9.084	9.076 (1.000)		531545		50.0000	
70 1-Chlorohexane	91	9.108	9.099 (1.003)		9649		1.00000	1
72 1,1,1,2-Tetrachloroethane	131	9.224	9.227 (1.015)		5585		1.00000	1
79 trans-1,4-Dichloro-2-butene	75	10.629	10.528 (1.170)		122996		1.00000	68(T)
\$ 80 Bromofluorobenzene	95	10.629	10.632 (1.170)		276913		1.00000	50
* 93 1,4-Dichlorobenzene-d4	152	12.244	12.247 (1.000)		236920		50.0000	
97 n-Butylbenzene	91	12.813	12.816 (1.046)		17552		1.00000	1
99 1,2-Dibromo-3-chloropropane	75	13.917	13.908 (1.137)		1746		1.00000	1

QC Flag Legend

T - Target compound detected outside RT window.

Data File: \\avogadro\organics\W5.I\130530.B\W503801.D
Date : 30-May-2013 12:48
Client ID: WSTD0015V
Sample Info: 5HL,WSTD0015V,WSTD0015V
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: ML
Column diameter: 0.25



Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130530.B\V503802.D
 Lab Smp Id: VSTD0055V Client Smp ID: VSTD0055V
 Inj Date : 30-MAY-2013 13:14
 Operator : WL SRC: WL Inst ID: V5.i
 Smp Info : 5ML,VSTD0055V,VSTD0055V
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130530.B\v5_8260W.m
 Meth Date : 03-Jun-2013 13:35 wluo Quant Type: ISTD
 Cal Date : 30-MAY-2013 13:14 Cal File: V503802.D
 Als bottle: 4 Calibration Sample, Level: 1
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: FULL.sub
 Target Version: 4.14
 Processing Host: TARGET103

Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG						AMOUNTS	
		MASS	RT	EXP RT	REL RT	RESPONSE	CAL-AMT (ug/L)	ON-COL (ug/L)
1 Dichlorodifluoromethane	85		1.579	1.583 (0.282)		32922	5.00000	4
3 Chloromethane	50		1.707	1.734 (0.305)		61804	5.00000	4
4 Vinyl Chloride	62		1.835	1.850 (0.328)		54772	5.00000	4
5 Bromomethane	94		2.125	2.141 (0.380)		39290	5.00000	5
6 Chloroethane	64		2.230	2.233 (0.398)		31362	5.00000	5
7 Trichlorofluoromethane	101		2.439	2.454 (0.436)		42325	5.00000	4
8 Ethanol	46		2.613	2.617 (0.467)		21583	500.000	410
9 Ether	59		2.694	2.698 (0.481)		32536	5.00000	5
10 Acrolein	56		2.811	2.814 (0.502)		39871	25.0000	23
11 1,1-Dichloroethene	96		2.904	2.907 (0.519)		34105	5.00000	5
12 1,1,2-Trichloro-1,2,2-Trifluo	101		2.915	2.919 (0.521)		29489	5.00000	5
13 Acetone	58		2.938	2.942 (0.525)		5797	5.00000	5
14 Iodomethane	142		3.043	3.047 (0.544)		52041	5.00000	4
15 Carbon Disulfide	76		3.101	3.105 (0.554)		139701	5.00000	4
16 Acetonitrile	41		3.217	3.221 (0.575)		143368	50.0000	47
17 Allyl Chloride	39		3.217	3.221 (0.575)		57752	5.00000	5
18 Methyl Acetate	43		3.229	3.232 (0.577)		49139	5.00000	4
19 Methylene Chloride	84		3.322	3.325 (0.593)		38562	5.00000	4
20 tert-Butanol	59		3.438	3.430 (0.614)		7830	10.0000	9
21 Acrylonitrile	53		3.542	3.534 (0.633)		14717	5.00000	4
22 trans-1,2-Dichloroethene	96		3.566	3.569 (0.637)		32920	5.00000	4
23 Methyl tert-butyl ether	73		3.577	3.581 (0.639)		91622	5.00000	5

Compounds	QUANT SIG						AMOUNTS	
	MASS	RT	EXP RT	REL RT	RESPONSE		CAL-AMT (ug/L)	ON-COL (ug/L)
=====	=====	=====	=====	=====	=====		=====	=====
24 1,1-Dichloroethane	63	3.949	3.941 (0.705)		69572		5.00000	5
25 Vinyl acetate	43	3.996	3.999 (0.714)		142460		5.00000	5
26 Diisopropyl Ether	45	4.007	4.011 (0.716)		169147		5.00000	5
27 2-Chloro-1,3-Butadiene	53	4.030	4.034 (0.720)		50535		5.00000	4
28 Ethyl tert-butyl ether	59	4.344	4.348 (0.776)		114754		5.00000	5
29 cis-1,2-Dichloroethene	96	4.483	4.475 (0.801)		35653		5.00000	5
30 2,2-Dichloropropane	77	4.483	4.487 (0.801)		34359		5.00000	5
31 2-Butanone	72	4.495	4.487 (0.803)		3400		5.00000	4
32 Propionitrile	54	4.541	4.534 (0.811)		51810		50.0000	43
33 Methacrylonitrile	41	4.681	4.685 (0.836)		53566		10.0000	9
34 Bromochloromethane	128	4.704	4.708 (0.840)		17123		5.00000	5
35 Tetrahydrofuran	72	4.762	4.754 (0.851)		8803		10.0000	9
36 Chloroform	83	4.774	4.777 (0.853)		58574		5.00000	5
\$ 37 Dibromofluoromethane	113	4.925	4.928 (0.880)		238884		50.0000	51
38 1,1,1-Trichloroethane	97	4.971	4.975 (0.888)		37539		5.00000	4
39 Cyclohexane	56	5.029	5.033 (0.898)		52972		5.00000	4
40 1,1-Dichloropropene	110	5.122	5.126 (0.915)		14352		5.00000	4
41 Carbon Tetrachloride	117	5.134	5.138 (0.917)		30157		5.00000	4
42 Isobutyl Alcohol	43	5.227	5.219 (0.934)		40986		100.000	96
\$ 43 1,2-Dichloroethane-d4	102	5.262	5.265 (0.940)		54458		50.0000	52
44 Benzene	78	5.331	5.335 (0.952)		122934		5.00000	5
45 1,2-Dichloroethane	62	5.331	5.335 (0.952)		45746		5.00000	5
46 tert-Amyl methyl ether	73	5.436	5.440 (0.971)		87435		5.00000	5
* 47 Fluorobenzene	96	5.599	5.602 (1.000)		757597		50.0000	
48 Trichloroethene	130	5.982	5.986 (1.068)		29721		5.00000	4
49 Methylcyclohexane	83	6.203	6.195 (1.108)		35952		5.00000	4
50 1,2-Dichloropropane	63	6.214	6.218 (1.110)		39228		5.00000	5
51 Dibromomethane	93	6.342	6.346 (1.133)		24011		5.00000	5
52 Methyl Methacrylate	69	6.342	6.346 (1.133)		25680		5.00000	4
53 1,4-Dioxane	88	6.377	6.357 (1.139)		7471		100.000	110
54 Bromodichloromethane	83	6.505	6.508 (1.162)		41864		5.00000	5
M 55 1,2-Dichloroethene (Total)	96				68573		10.0000	9
56 2-Chloroethyl vinyl ether	63	6.865	6.846 (1.226)		629		5.00000	5(T)
57 cis-1,3-Dichloropropene	75	7.016	7.008 (1.253)		51361		5.00000	4
58 4-Methyl-2-pentanone	43	7.190	7.182 (1.284)		49351		5.00000	4
\$ 59 Toluene-d8	98	7.318	7.322 (0.806)		694258		50.0000	50
60 Toluene	91	7.399	7.403 (1.322)		115751		5.00000	5
61 trans-1,3-Dichloropropene	75	7.655	7.647 (1.367)		43663		5.00000	4
62 Ethyl Methacrylate	69	7.771	7.775 (1.388)		27987		5.00000	4
63 1,1,2-Trichloroethane	97	7.875	7.867 (1.407)		25958		5.00000	4
64 Tetrachloroethene	164	8.061	8.053 (0.887)		23517		5.00000	4
65 1,3-Dichloropropane	76	8.073	8.077 (0.889)		48916		5.00000	5
66 2-Hexanone	43	8.189	8.181 (0.902)		31594		5.00000	4
67 Dibromochloromethane	129	8.340	8.344 (0.918)		26886		5.00000	4
68 1,2-Dibromoethane	107	8.491	8.483 (0.935)		30560		5.00000	4
* 69 Chlorobenzene-d5	117	9.084	9.076 (1.000)		536169		50.0000	
70 1-Chlorohexane	91	9.095	9.099 (1.001)		39412		5.00000	4
71 Chlorobenzene	112	9.118	9.110 (1.004)		72726		5.00000	5
72 1,1,1,2-Tetrachloroethane	131	9.223	9.227 (1.015)		25244		5.00000	4
73 Ethylbenzene	106	9.269	9.273 (1.020)		36902		5.00000	4
74 m,p-Xylene	106	9.420	9.424 (1.037)		93531		10.0000	9
75 o-Xylene	106	9.943	9.935 (1.095)		46137		5.00000	5
76 Styrene	104	9.955	9.958 (1.096)		62060		5.00000	4
77 Bromoform	173	10.187	10.179 (1.121)		16014		5.00000	4

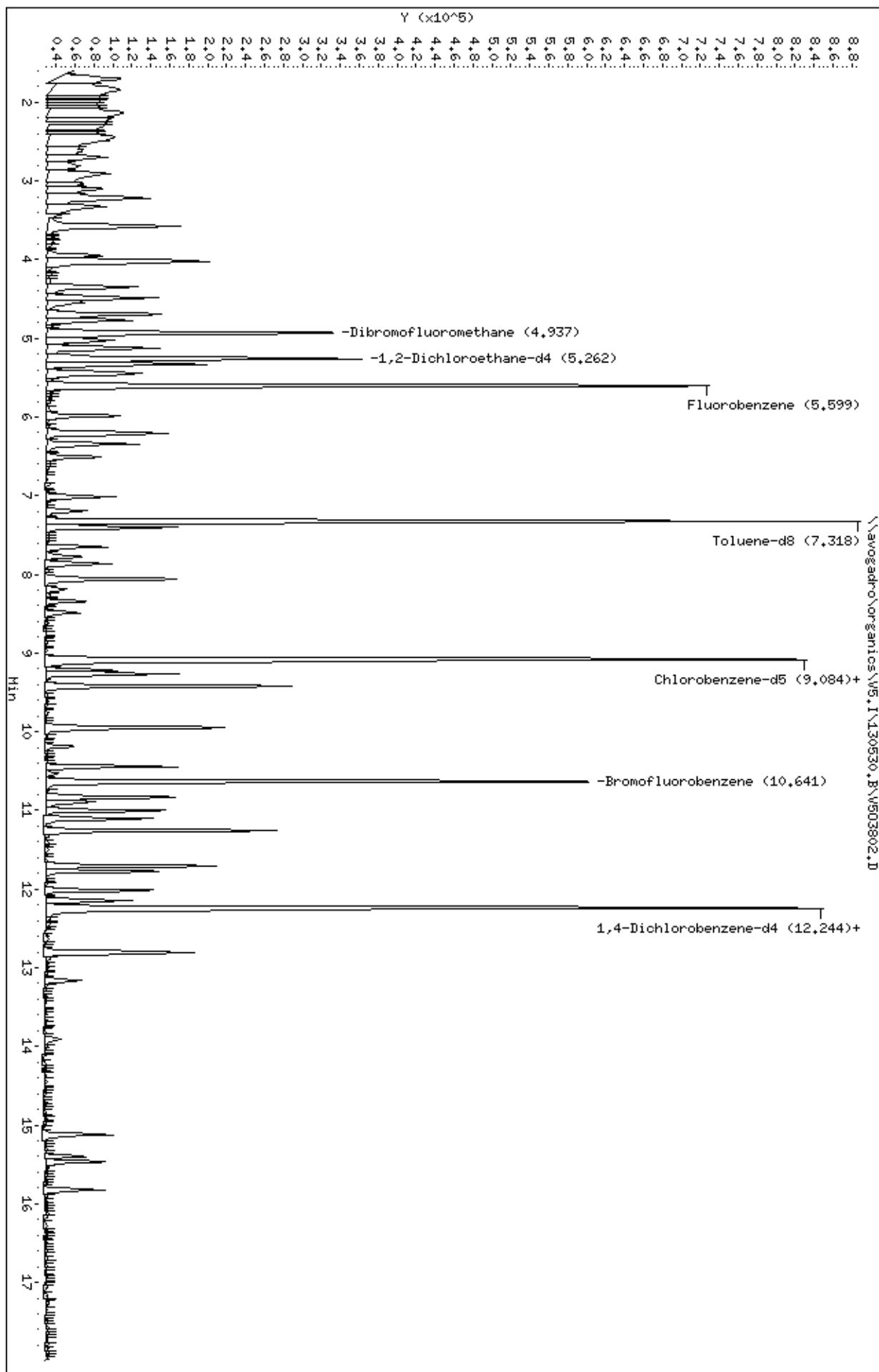
Compounds	QUANT SIG						AMOUNTS	
	MASS	RT	EXP RT	REL RT	RESPONSE		CAL-AMT (ug/L)	ON-COL (ug/L)
=====	=====	=====	=====	=====	=====		=====	=====
78 Isopropylbenzene	105	10.443	10.446 (1.150)		111115		5.00000	4
79 trans-1,4-Dichloro-2-butene	75	10.524	10.528 (1.159)		4381		5.00000	2
\$ 80 Bromofluorobenzene	95	10.629	10.632 (1.170)		278682		50.0000	50
81 Bromobenzene	156	10.838	10.830 (0.885)		29245		5.00000	4
82 1,1,2,2-Tetrachloroethane	83	10.838	10.841 (0.885)		41336		5.00000	5
83 1,2,3-Trichloropropane	75	10.896	10.888 (0.890)		35514		5.00000	4
84 n-Propylbenzene	120	11.000	11.004 (0.898)		26268		5.00000	4
85 2-Chlorotoluene	126	11.105	11.109 (0.907)		26411		5.00000	4
86 4-Chlorotoluene	126	11.256	11.260 (0.919)		27881		5.00000	4
87 1,3,5-Trimethylbenzene	105	11.256	11.260 (0.919)		88431		5.00000	5
88 tert-Butylbenzene	119	11.709	11.713 (0.956)		80932		5.00000	4
89 1,2,4-Trimethylbenzene	105	11.779	11.771 (0.962)		87860		5.00000	4
90 sec-Butylbenzene	105	12.011	12.015 (0.981)		108837		5.00000	4
91 1,3-Dichlorobenzene	146	12.150	12.142 (0.992)		48240		5.00000	4
92 4-Isopropyltoluene	119	12.232	12.235 (0.999)		79677		5.00000	4
* 93 1,4-Dichlorobenzene-d4	152	12.243	12.247 (1.000)		238114		50.0000	
94 1,4-Dichlorobenzene	146	12.278	12.270 (1.003)		52171		5.00000	5
95 1,2-Dichlorobenzene	146	12.789	12.793 (1.045)		47722		5.00000	4
M 96 Xylene (Total)	106				139668		15.0000	14
97 n-Butylbenzene	91	12.812	12.816 (1.046)		80665		5.00000	4
98 Hexachloroethane	117	13.161	13.165 (1.075)		11777		5.00000	4
99 1,2-Dibromo-3-chloropropane	75	13.916	13.908 (1.137)		5967		5.00000	5
100 1,3,5-Trichlorobenzene	182	15.124	15.128 (2.701)		26953		5.00000	4
101 1,2,4-Trichlorobenzene	180	15.124	15.128 (1.235)		28494		5.00000	4
102 Hexachlorobutadiene	225	15.403	15.407 (1.258)		11276		5.00000	4
103 Naphthalene	128	15.461	15.465 (1.263)		64029		5.00000	4
104 1,2,3-Trichlorobenzene	180	15.821	15.825 (1.292)		25188		5.00000	4

QC Flag Legend

T - Target compound detected outside RT window.

Data File: \\avogadro\organics\W5.I\130530.B\W503802.D
Date : 30-May-2013 13:14
Client ID: WSTD0055V
Sample Info: 5HL,WSTD0055V,WSTD0055V
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: ML
Column diameter: 0.25



Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
Data file : \\avogadro\organics\V5.I\130530.B\V503803.D
Lab Smp Id: VSTD0205V Client Smp ID: VSTD0205V
Inj Date : 30-MAY-2013 13:39
Operator : WL SRC: WL Inst ID: V5.i
Smp Info : 5ML,VSTD0205V,VSTD0205V
Misc Info :
Comment :
Method : \\avogadro\organics\V5.I\130530.B\v5_8260W.m
Meth Date : 03-Jun-2013 13:35 wluo Quant Type: ISTD
Cal Date : 30-MAY-2013 13:39 Cal File: V503803.D
Als bottle: 5 Calibration Sample, Level: 2
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: FULL.sub
Target Version: 4.14
Processing Host: TARGET103

Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG	AMOUNTS					
		RT	EXP RT	REL RT	RESPONSE	CAL-AMT (ug/L)	ON-COL (ug/L)
-----	----	----	-----	-----	-----	-----	-----
1 Dichlorodifluoromethane	85	1.586	1.583 (0.283)		132228	20.0000	17
3 Chloromethane	50	1.725	1.734 (0.308)		260586	20.0000	18
4 Vinyl Chloride	62	1.829	1.850 (0.326)		219664	20.0000	18
5 Bromomethane	94	2.143	2.141 (0.382)		145201	20.0000	20
6 Chloroethane	64	2.236	2.233 (0.399)		119377	20.0000	18
7 Trichlorofluoromethane	101	2.457	2.454 (0.438)		173035	20.0000	18
8 Ethanol	46	2.619	2.617 (0.467)		91530	2000.00	1700
9 Ether	59	2.701	2.698 (0.482)		128172	20.0000	19
10 Acrolein	56	2.817	2.814 (0.503)		158857	100.000	91
11 1,1-Dichloroethene	96	2.910	2.907 (0.519)		138203	20.0000	20
12 1,1,2-Trichloro-1,2,2-Trifluo	101	2.921	2.919 (0.521)		111805	20.0000	18
13 Acetone	58	2.945	2.942 (0.525)		19551	20.0000	18
14 Iodomethane	142	3.038	3.047 (0.542)		215791	20.0000	18
15 Carbon Disulfide	76	3.107	3.105 (0.554)		570514	20.0000	18
16 Acetonitrile	41	3.212	3.221 (0.573)		580455	200.000	190
17 Allyl Chloride	39	3.223	3.221 (0.575)		231822	20.0000	19(T)
18 Methyl Acetate	43	3.235	3.232 (0.577)		203923	20.0000	18
19 Methylene Chloride	84	3.328	3.325 (0.594)		150832	20.0000	18
20 tert-Butanol	59	3.433	3.430 (0.612)		33643	40.0000	40
21 Acrylonitrile	53	3.537	3.534 (0.631)		63364	20.0000	20
22 trans-1,2-Dichloroethene	96	3.572	3.569 (0.637)		135791	20.0000	18
23 Methyl tert-butyl ether	73	3.584	3.581 (0.639)		354663	20.0000	18

Compounds	QUANT SIG						AMOUNTS	
	MASS	RT	EXP RT	REL RT	RESPONSE		CAL-AMT (ug/L)	ON-COL (ug/L)
=====	=====	=====	=====	=====	=====		=====	=====
24 1,1-Dichloroethane	63	3.944	3.941	(0.704)	279934		20.0000	19
25 Vinyl acetate	43	4.002	3.999	(0.714)	580431		20.0000	20
26 Diisopropyl Ether	45	4.013	4.011	(0.716)	669454		20.0000	19
27 2-Chloro-1,3-Butadiene	53	4.037	4.034	(0.720)	208127		20.0000	19
28 Ethyl tert-butyl ether	59	4.350	4.348	(0.776)	468067		20.0000	19
29 cis-1,2-Dichloroethene	96	4.478	4.475	(0.799)	141901		20.0000	17
30 2,2-Dichloropropane	77	4.490	4.487	(0.801)	129657		20.0000	18
31 2-Butanone	72	4.501	4.487	(0.803)	17224		20.0000	19
32 Propionitrile	54	4.536	4.534	(0.809)	229763		200.000	190
33 Methacrylonitrile	41	4.687	4.685	(0.836)	241772		40.0000	40
34 Bromochloromethane	128	4.710	4.708	(0.840)	71119		20.0000	19
35 Tetrahydrofuran	72	4.757	4.754	(0.849)	34150		40.0000	36
36 Chloroform	83	4.780	4.777	(0.853)	230796		20.0000	19
\$ 37 Dibromofluoromethane	113	4.931	4.928	(0.880)	235426		50.0000	50
38 1,1,1-Trichloroethane	97	4.966	4.975	(0.886)	154687		20.0000	18
39 Cyclohexane	56	5.036	5.033	(0.898)	217638		20.0000	17
40 1,1-Dichloropropene	110	5.129	5.126	(0.915)	57733		20.0000	18
41 Carbon Tetrachloride	117	5.140	5.138	(0.917)	120850		20.0000	18
42 Isobutyl Alcohol	43	5.222	5.219	(0.932)	160044		400.000	370
\$ 43 1,2-Dichloroethane-d4	102	5.268	5.265	(0.940)	51774		50.0000	49
44 Benzene	78	5.326	5.335	(0.950)	494920		20.0000	19
45 1,2-Dichloroethane	62	5.338	5.335	(0.952)	184683		20.0000	19
46 tert-Amyl methyl ether	73	5.442	5.440	(0.971)	356739		20.0000	19
* 47 Fluorobenzene	96	5.605	5.602	(1.000)	758047		50.0000	
48 Trichloroethene	130	5.988	5.986	(1.068)	124434		20.0000	19
49 Methylcyclohexane	83	6.197	6.195	(1.106)	137636		20.0000	17
50 1,2-Dichloropropane	63	6.221	6.218	(1.110)	152873		20.0000	18
51 Dibromomethane	93	6.337	6.346	(1.131)	97631		20.0000	19
52 Methyl Methacrylate	69	6.337	6.346	(1.131)	109704		20.0000	19
53 1,4-Dioxane	88	6.360	6.357	(1.135)	26524		400.000	380
54 Bromodichloromethane	83	6.511	6.508	(1.162)	165828		20.0000	19
M 55 1,2-Dichloroethene (Total)	96				277692		40.0000	36
56 2-Chloroethyl vinyl ether	63	6.859	6.846	(1.224)	3029		20.0000	22
57 cis-1,3-Dichloropropene	75	7.010	7.008	(1.251)	213355		20.0000	19
58 4-Methyl-2-pentanone	43	7.185	7.182	(1.282)	208699		20.0000	19
\$ 59 Toluene-d8	98	7.324	7.322	(0.807)	691005		50.0000	50
60 Toluene	91	7.405	7.403	(1.321)	469198		20.0000	19
61 trans-1,3-Dichloropropene	75	7.649	7.647	(1.365)	185036		20.0000	19
62 Ethyl Methacrylate	69	7.777	7.775	(1.388)	116837		20.0000	18
63 1,1,2-Trichloroethane	97	7.870	7.867	(1.404)	111247		20.0000	18
64 Tetrachloroethene	164	8.056	8.053	(0.887)	96654		20.0000	19
65 1,3-Dichloropropane	76	8.068	8.077	(0.889)	197814		20.0000	19
66 2-Hexanone	43	8.184	8.181	(0.901)	136250		20.0000	19
67 Dibromochloromethane	129	8.346	8.344	(0.919)	119805		20.0000	20
68 1,2-Dibromoethane	107	8.486	8.483	(0.935)	131082		20.0000	19
* 69 Chlorobenzene-d5	117	9.078	9.076	(1.000)	534273		50.0000	
70 1-Chlorohexane	91	9.101	9.099	(1.003)	160897		20.0000	18
71 Chlorobenzene	112	9.113	9.110	(1.004)	295038		20.0000	19
72 1,1,1,2-Tetrachloroethane	131	9.218	9.227	(1.015)	104892		20.0000	19
73 Ethylbenzene	106	9.264	9.273	(1.020)	149142		20.0000	18
74 m,p-Xylene	106	9.427	9.424	(1.038)	380985		40.0000	38
75 o-Xylene	106	9.938	9.935	(1.095)	183407		20.0000	18
76 Styrene	104	9.961	9.958	(1.097)	277645		20.0000	19
77 Bromoform	173	10.182	10.179	(1.122)	72115		20.0000	20

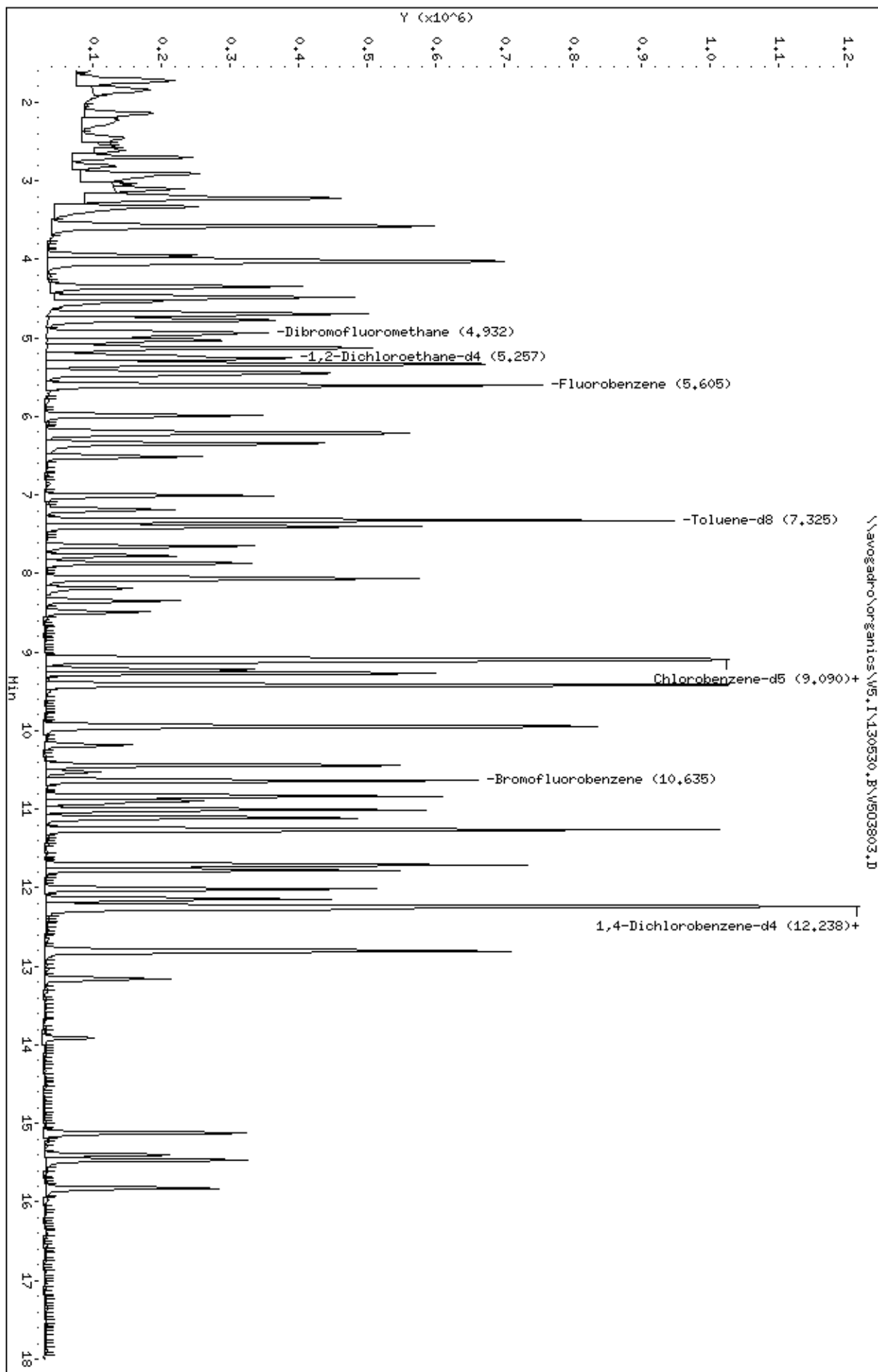
Compounds	QUANT SIG						AMOUNTS	
	MASS	RT	EXP RT	REL RT	RESPONSE		CAL-AMT (ug/L)	ON-COL (ug/L)
=====	=====	=====	=====	=====	=====		=====	=====
78 Isopropylbenzene	105	10.437	10.446	(1.150)	461467		20.0000	19
79 trans-1,4-Dichloro-2-butene	75	10.519	10.528	(1.159)	24774		20.0000	14
\$ 80 Bromofluorobenzene	95	10.635	10.632	(1.171)	278137		50.0000	50
81 Bromobenzene	156	10.832	10.830	(0.885)	121311		20.0000	18
82 1,1,2,2-Tetrachloroethane	83	10.844	10.841	(0.886)	171259		20.0000	19
83 1,2,3-Trichloropropane	75	10.890	10.888	(0.890)	148635		20.0000	18
84 n-Propylbenzene	120	11.007	11.004	(0.899)	112447		20.0000	19
85 2-Chlorotoluene	126	11.111	11.109	(0.908)	111002		20.0000	19
86 4-Chlorotoluene	126	11.262	11.260	(0.920)	114418		20.0000	18
87 1,3,5-Trimethylbenzene	105	11.262	11.260	(0.920)	363083		20.0000	18
88 tert-Butylbenzene	119	11.704	11.713	(0.956)	344106		20.0000	19
89 1,2,4-Trimethylbenzene	105	11.773	11.771	(0.962)	364776		20.0000	18
90 sec-Butylbenzene	105	12.017	12.015	(0.982)	453667		20.0000	19
91 1,3-Dichlorobenzene	146	12.145	12.142	(0.992)	206548		20.0000	19
92 4-Isopropyltoluene	119	12.226	12.235	(0.999)	347457		20.0000	19
* 93 1,4-Dichlorobenzene-d4	152	12.238	12.247	(1.000)	244213		50.0000	
94 1,4-Dichlorobenzene	146	12.273	12.270	(1.003)	213504		20.0000	19
95 1,2-Dichlorobenzene	146	12.796	12.793	(1.046)	199226		20.0000	18
M 96 Xylene (Total)	106				564392		60.0000	57
97 n-Butylbenzene	91	12.819	12.816	(1.047)	338572		20.0000	19
98 Hexachloroethane	117	13.156	13.165	(1.075)	55767		20.0000	18
99 1,2-Dibromo-3-chloropropane	75	13.911	13.908	(1.137)	24555		20.0000	18
100 1,3,5-Trichlorobenzene	182	15.119	15.128	(2.697)	122685		20.0000	19
101 1,2,4-Trichlorobenzene	180	15.119	15.128	(1.235)	126515		20.0000	18
102 Hexachlorobutadiene	225	15.398	15.407	(1.258)	50575		20.0000	18
103 Naphthalene	128	15.467	15.465	(1.264)	296617		20.0000	19
104 1,2,3-Trichlorobenzene	180	15.827	15.825	(1.293)	109579		20.0000	19

QC Flag Legend

T - Target compound detected outside RT window.

Data File: \\avogadro\organics\W5.1\130530.B\W503803.D
Date : 30-May-2013 13:39
Client ID: WSTD0205V
Sample Info: 5HL,WSTD0205V,WSTD0205V
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: ML
Column diameter: 0.25



Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130530.B\V503804.D
 Lab Smp Id: VSTD0505V Client Smp ID: VSTD0505V
 Inj Date : 30-MAY-2013 14:05
 Operator : WL SRC: WL Inst ID: V5.i
 Smp Info : 5ML,VSTD0505V,VSTD0505V
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130530.B\v5_8260W.m
 Meth Date : 03-Jun-2013 13:35 wluo Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:05 Cal File: V503804.D
 Als bottle: 6 Calibration Sample, Level: 3
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: FULL.sub
 Target Version: 4.14
 Processing Host: TARGET103

Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

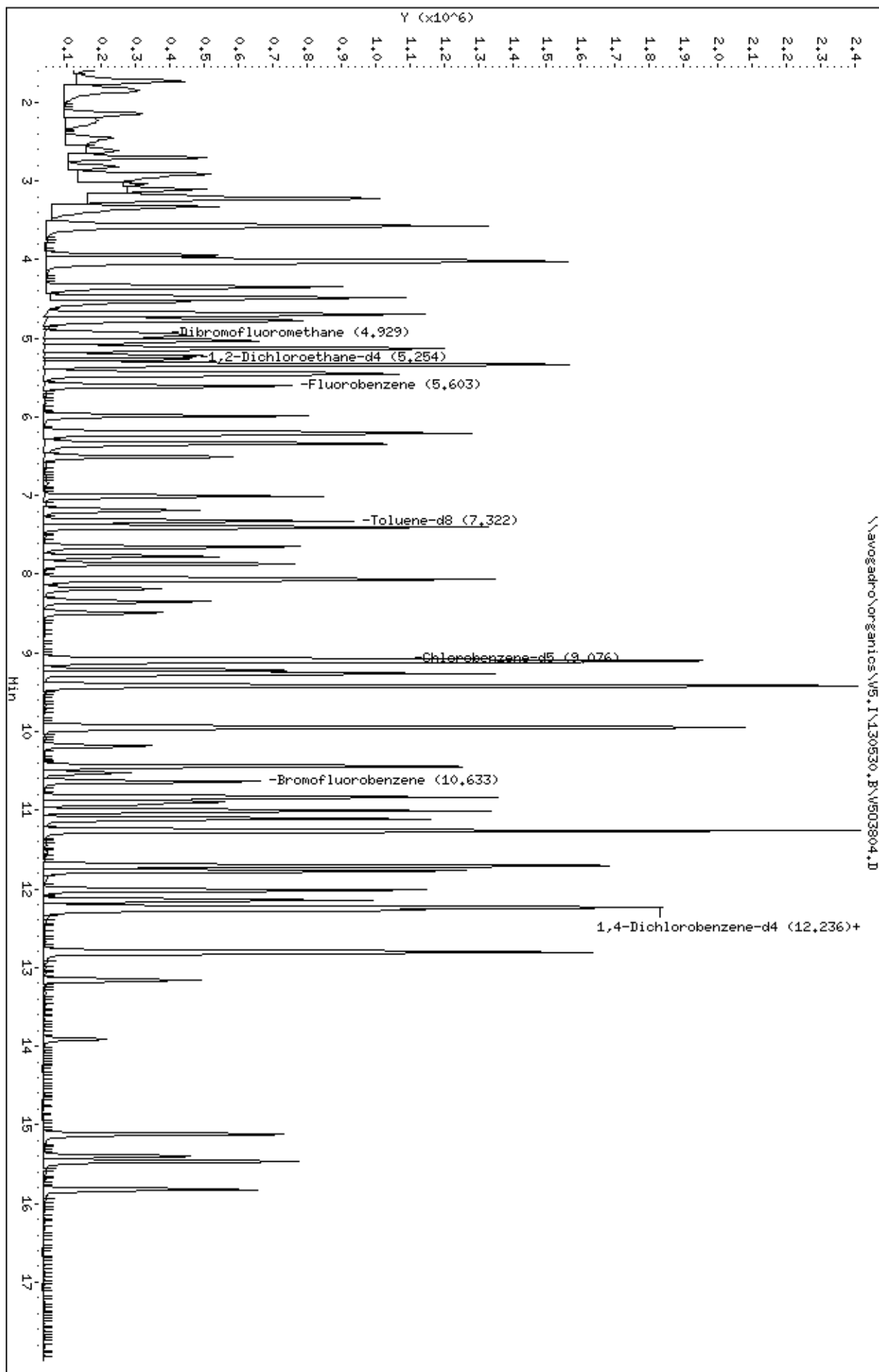
Compounds	QUANT SIG						AMOUNTS	
		MASS	RT	EXP RT	REL RT	RESPONSE	CAL-AMT (ug/L)	ON-COL (ug/L)
1 Dichlorodifluoromethane	85		1.583	1.583 (0.283)		361103	50.0000	46
3 Chloromethane	50		1.734	1.734 (0.310)		656487	50.0000	45
4 Vinyl Chloride	62		1.850	1.850 (0.330)		538132	50.0000	44
5 Bromomethane	94		2.141	2.141 (0.382)		337417	50.0000	45
6 Chloroethane	64		2.233	2.233 (0.399)		284691	50.0000	43
7 Trichlorofluoromethane	101		2.454	2.454 (0.438)		427870	50.0000	45
8 Ethanol	46		2.617	2.617 (0.467)		242832	5000.00	4500
9 Ether	59		2.698	2.698 (0.482)		306194	50.0000	44
10 Acrolein	56		2.814	2.814 (0.502)		360295	250.000	200
11 1,1-Dichloroethene	96		2.907	2.907 (0.519)		309697	50.0000	44
12 1,1,2-Trichloro-1,2,2-Trifluo	101		2.919	2.919 (0.521)		283948	50.0000	46
13 Acetone	58		2.942	2.942 (0.525)		41152	50.0000	37
14 Iodomethane	142		3.047	3.047 (0.544)		514176	50.0000	44
15 Carbon Disulfide	76		3.105	3.105 (0.554)		1347691	50.0000	42
16 Acetonitrile	41		3.221	3.221 (0.575)		1382919	500.000	450
17 Allyl Chloride	39		3.221	3.221 (0.575)		524208	50.0000	44
18 Methyl Acetate	43		3.232	3.232 (0.577)		472297	50.0000	41
19 Methylene Chloride	84		3.325	3.325 (0.594)		355469	50.0000	42
20 tert-Butanol	59		3.430	3.430 (0.612)		75327	100.000	89
21 Acrylonitrile	53		3.534	3.534 (0.631)		156854	50.0000	49
22 trans-1,2-Dichloroethene	96		3.569	3.569 (0.637)		316252	50.0000	43
23 Methyl tert-butyl ether	73		3.581	3.581 (0.639)		841644	50.0000	44

Compounds	QUANT SIG						AMOUNTS	
	MASS	RT	EXP RT	REL RT	RESPONSE		CAL-AMT (ug/L)	ON-COL (ug/L)
=====	=====	=====	=====	=====	=====		=====	=====
24 1,1-Dichloroethane	63	3.941	3.941	(0.704)	657619		50.0000	44
25 Vinyl acetate	43	3.999	3.999	(0.714)	1347846		50.0000	45
26 Diisopropyl Ether	45	4.011	4.011	(0.716)	1536465		50.0000	44
27 2-Chloro-1,3-Butadiene	53	4.034	4.034	(0.720)	495274		50.0000	45
28 Ethyl tert-butyl ether	59	4.348	4.348	(0.776)	1102936		50.0000	44
29 cis-1,2-Dichloroethene	96	4.475	4.475	(0.799)	336778		50.0000	40
30 2,2-Dichloropropane	77	4.487	4.487	(0.801)	298412		50.0000	41
31 2-Butanone	72	4.487	4.487	(0.801)	43916		50.0000	48
32 Propionitrile	54	4.534	4.534	(0.809)	548180		500.000	450
33 Methacrylonitrile	41	4.685	4.685	(0.836)	570699		100.000	93
34 Bromochloromethane	128	4.708	4.708	(0.840)	166568		50.0000	44
35 Tetrahydrofuran	72	4.754	4.754	(0.849)	85853		100.000	91
36 Chloroform	83	4.777	4.777	(0.853)	525965		50.0000	43
\$ 37 Dibromofluoromethane	113	4.928	4.928	(0.880)	240525		50.0000	51
38 1,1,1-Trichloroethane	97	4.975	4.975	(0.888)	362036		50.0000	43
39 Cyclohexane	56	5.033	5.033	(0.898)	588104		50.0000	47
40 1,1-Dichloropropene	110	5.126	5.126	(0.915)	144574		50.0000	45
41 Carbon Tetrachloride	117	5.138	5.138	(0.917)	309111		50.0000	46
42 Isobutyl Alcohol	43	5.219	5.219	(0.932)	398164		1000.00	920
\$ 43 1,2-Dichloroethane-d4	102	5.265	5.265	(0.940)	51048		50.0000	48
44 Benzene	78	5.335	5.335	(0.952)	1176164		50.0000	45
45 1,2-Dichloroethane	62	5.335	5.335	(0.952)	429211		50.0000	45
46 tert-Amyl methyl ether	73	5.440	5.440	(0.971)	848966		50.0000	44
* 47 Fluorobenzene	96	5.602	5.602	(1.000)	761578		50.0000	
48 Trichloroethene	130	5.986	5.986	(1.068)	306119		50.0000	46
49 Methylcyclohexane	83	6.195	6.195	(1.106)	356801		50.0000	44
50 1,2-Dichloropropane	63	6.218	6.218	(1.110)	365450		50.0000	44
51 Dibromomethane	93	6.346	6.346	(1.133)	225657		50.0000	45
52 Methyl Methacrylate	69	6.346	6.346	(1.133)	259610		50.0000	45
53 1,4-Dioxane	88	6.357	6.357	(1.135)	59461		1000.00	850
54 Bromodichloromethane	83	6.508	6.508	(1.162)	399972		50.0000	45
M 55 1,2-Dichloroethene (Total)	96				653030		100.000	83
56 2-Chloroethyl vinyl ether	63	6.845	6.846	(1.222)	8519		50.0000	63
57 cis-1,3-Dichloropropene	75	7.008	7.008	(1.251)	519104		50.0000	45
58 4-Methyl-2-pentanone	43	7.182	7.182	(1.282)	490680		50.0000	45
\$ 59 Toluene-d8	98	7.322	7.322	(0.807)	687685		50.0000	49
60 Toluene	91	7.403	7.403	(1.321)	1094587		50.0000	45
61 trans-1,3-Dichloropropene	75	7.647	7.647	(1.365)	461524		50.0000	47
62 Ethyl Methacrylate	69	7.775	7.775	(1.388)	302233		50.0000	45
63 1,1,2-Trichloroethane	97	7.867	7.867	(1.404)	264275		50.0000	43
64 Tetrachloroethene	164	8.053	8.053	(0.887)	229750		50.0000	44
65 1,3-Dichloropropane	76	8.077	8.077	(0.890)	473633		50.0000	44
66 2-Hexanone	43	8.181	8.181	(0.901)	340483		50.0000	46
67 Dibromochloromethane	129	8.344	8.344	(0.919)	294934		50.0000	47
68 1,2-Dibromoethane	107	8.483	8.483	(0.935)	312059		50.0000	45
* 69 Chlorobenzene-d5	117	9.076	9.076	(1.000)	546450		50.0000	
70 1-Chlorohexane	91	9.099	9.099	(1.003)	387045		50.0000	44
71 Chlorobenzene	112	9.110	9.110	(1.004)	717418		50.0000	45
72 1,1,1,2-Tetrachloroethane	131	9.227	9.227	(1.017)	249915		50.0000	44
73 Ethylbenzene	106	9.273	9.273	(1.022)	367397		50.0000	44
74 m,p-Xylene	106	9.424	9.424	(1.038)	891238		100.000	87
75 o-Xylene	106	9.935	9.935	(1.095)	458316		50.0000	45
76 Styrene	104	9.958	9.958	(1.097)	709616		50.0000	48
77 Bromoform	173	10.179	10.179	(1.122)	181239		50.0000	48

Compounds	QUANT SIG						AMOUNTS	
	MASS	RT	EXP RT	REL RT	RESPONSE		CAL-AMT (ug/L)	ON-COL (ug/L)
=====	=====	=====	=====	=====	=====		=====	=====
78 Isopropylbenzene	105	10.446	10.446	(1.151)	1109253		50.0000	45
79 trans-1,4-Dichloro-2-butene	75	10.528	10.528	(1.160)	84315		50.0000	46
\$ 80 Bromofluorobenzene	95	10.632	10.632	(1.172)	287018		50.0000	51
81 Bromobenzene	156	10.830	10.830	(0.884)	289604		50.0000	44
82 1,1,2,2-Tetrachloroethane	83	10.841	10.841	(0.885)	408181		50.0000	44
83 1,2,3-Trichloropropane	75	10.888	10.888	(0.889)	376605		50.0000	45
84 n-Propylbenzene	120	11.004	11.004	(0.899)	270271		50.0000	45
85 2-Chlorotoluene	126	11.109	11.109	(0.907)	264929		50.0000	44
86 4-Chlorotoluene	126	11.260	11.260	(0.919)	275130		50.0000	44
87 1,3,5-Trimethylbenzene	105	11.260	11.260	(0.919)	882193		50.0000	45
88 tert-Butylbenzene	119	11.713	11.713	(0.956)	823973		50.0000	44
89 1,2,4-Trimethylbenzene	105	11.771	11.771	(0.961)	895788		50.0000	45
90 sec-Butylbenzene	105	12.015	12.015	(0.981)	1098867		50.0000	45
91 1,3-Dichlorobenzene	146	12.142	12.142	(0.991)	496416		50.0000	45
92 4-Isopropyltoluene	119	12.235	12.235	(0.999)	820906		50.0000	44
* 93 1,4-Dichlorobenzene-d4	152	12.247	12.247	(1.000)	245450		50.0000	
94 1,4-Dichlorobenzene	146	12.270	12.270	(1.002)	508048		50.0000	44
95 1,2-Dichlorobenzene	146	12.793	12.793	(1.045)	481995		50.0000	45
M 96 Xylene (Total)	106				1349554		150.000	130
97 n-Butylbenzene	91	12.816	12.816	(1.046)	809118		50.0000	45
98 Hexachloroethane	117	13.165	13.165	(1.075)	146631		50.0000	48
99 1,2-Dibromo-3-chloropropane	75	13.908	13.908	(1.136)	54041		50.0000	40
100 1,3,5-Trichlorobenzene	182	15.128	15.128	(2.700)	300741		50.0000	47
101 1,2,4-Trichlorobenzene	180	15.128	15.128	(1.235)	307813		50.0000	45
102 Hexachlorobutadiene	225	15.407	15.407	(1.258)	118552		50.0000	43
103 Naphthalene	128	15.465	15.465	(1.263)	737789		50.0000	46
104 1,2,3-Trichlorobenzene	180	15.825	15.825	(1.292)	263900		50.0000	45

Data File: \\avogadro\organics\W5.I\130530.B\W503804.D
Date : 30-May-2013 14:05
Client ID: WSTD0505V
Sample Info: 5HL,WSTD0505V,WSTD0505V
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: ML
Column diameter: 0.25



Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130530.B\V503805.D
 Lab Smp Id: VSTD1005V Client Smp ID: VSTD1005V
 Inj Date : 30-MAY-2013 14:32
 Operator : WL SRC: WL Inst ID: V5.i
 Smp Info : 5ML,VSTD1005V,VSTD1005V
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130530.B\v5_8260W.m
 Meth Date : 03-Jun-2013 13:35 wluo Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V503805.D
 Als bottle: 7 Calibration Sample, Level: 4
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: FULL.sub
 Target Version: 4.14
 Processing Host: TARGET103

Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG MASS					AMOUNTS	
		RT	EXP RT	REL RT	RESPONSE	CAL-AMT (ug/L)	ON-COL (ug/L)
1 Dichlorodifluoromethane	85	1.588	1.583 (0.283)		569698	100.000	110
3 Chloromethane	50	1.728	1.734 (0.308)		1057565	100.000	110
4 Vinyl Chloride	62	1.844	1.850 (0.329)		900500	100.000	120
5 Bromomethane	94	2.146	2.141 (0.383)		487758	100.000	100
6 Chloroethane	64	2.227	2.233 (0.397)		452063	100.000	110
7 Trichlorofluoromethane	101	2.448	2.454 (0.437)		705030	100.000	120
8 Ethanol	46	2.622	2.617 (0.468)		439733	10000.0	13000
9 Ether	59	2.704	2.698 (0.482)		511013	100.000	120
10 Acrolein	56	2.808	2.814 (0.501)		675124	500.000	600
11 1,1-Dichloroethene	96	2.901	2.907 (0.517)		533133	100.000	120
12 1,1,2-Trichloro-1,2,2-Trifluo	101	2.913	2.919 (0.519)		489621	100.000	120
13 Acetone	58	2.947	2.942 (0.526)		89179	100.000	130
14 Iodomethane	142	3.040	3.047 (0.542)		857307	100.000	120
15 Carbon Disulfide	76	3.098	3.105 (0.553)		2228024	100.000	110
16 Acetonitrile	41	3.215	3.221 (0.573)		2229701	1000.00	1100
17 Allyl Chloride	39	3.215	3.221 (0.573)		866132	100.000	110(T)
18 Methyl Acetate	43	3.238	3.232 (0.577)		787362	100.000	110
19 Methylene Chloride	84	3.331	3.325 (0.594)		604267	100.000	110
20 tert-Butanol	59	3.435	3.430 (0.613)		117388	200.000	220
21 Acrylonitrile	53	3.540	3.534 (0.631)		251152	100.000	120
22 trans-1,2-Dichloroethene	96	3.575	3.569 (0.638)		527857	100.000	110
23 Methyl tert-butyl ether	73	3.575	3.581 (0.638)		1392294	100.000	110

Compounds	QUANT SIG						AMOUNTS	
	MASS	RT	EXP RT	REL RT	RESPONSE		CAL-AMT (ug/L)	ON-COL (ug/L)
=====	=====	=====	=====	=====	=====		=====	=====
24 1,1-Dichloroethane	63	3.946	3.941 (0.704)		1085248		100.000	110
25 Vinyl acetate	43	4.005	3.999 (0.714)		2098275		100.000	110
26 Diisopropyl Ether	45	4.016	4.011 (0.716)		2438187		100.000	110
27 2-Chloro-1,3-Butadiene	53	4.028	4.034 (0.718)		817442		100.000	120
28 Ethyl tert-butyl ether	59	4.341	4.348 (0.774)		1786827		100.000	110
29 cis-1,2-Dichloroethene	96	4.481	4.475 (0.799)		549410		100.000	110
30 2,2-Dichloropropane	77	4.492	4.487 (0.801)		525760		100.000	110
31 2-Butanone	72	4.492	4.487 (0.801)		71464		100.000	120
32 Propionitrile	54	4.539	4.534 (0.809)		894819		1000.00	1200
33 Methacrylonitrile	41	4.690	4.685 (0.836)		878294		200.000	220
34 Bromochloromethane	128	4.702	4.708 (0.838)		281229		100.000	120
35 Tetrahydrofuran	72	4.760	4.754 (0.849)		135291		200.000	220
36 Chloroform	83	4.771	4.777 (0.851)		862237		100.000	110
\$ 37 Dibromofluoromethane	113	4.934	4.928 (0.880)		146163		50.0000	49
38 1,1,1-Trichloroethane	97	4.969	4.975 (0.886)		589887		100.000	110
39 Cyclohexane	56	5.027	5.033 (0.896)		996312		100.000	120
40 1,1-Dichloropropene	110	5.131	5.126 (0.915)		230093		100.000	110
41 Carbon Tetrachloride	117	5.131	5.138 (0.915)		480915		100.000	110
42 Isobutyl Alcohol	43	5.224	5.219 (0.932)		616843		2000.00	2200
\$ 43 1,2-Dichloroethane-d4	102	5.259	5.265 (0.938)		33595		50.0000	50
44 Benzene	78	5.329	5.335 (0.950)		1926274		100.000	120
45 1,2-Dichloroethane	62	5.340	5.335 (0.952)		685558		100.000	110
46 tert-Amyl methyl ether	73	5.445	5.440 (0.971)		1386990		100.000	110
* 47 Fluorobenzene	96	5.608	5.602 (1.000)		483935		50.0000	
48 Trichloroethene	130	5.979	5.986 (1.066)		485138		100.000	110
49 Methylcyclohexane	83	6.200	6.195 (1.106)		634854		100.000	120
50 1,2-Dichloropropane	63	6.212	6.218 (1.108)		598321		100.000	110
51 Dibromomethane	93	6.339	6.346 (1.130)		365468		100.000	110
52 Methyl Methacrylate	69	6.339	6.346 (1.130)		422044		100.000	120
53 1,4-Dioxane	88	6.363	6.357 (1.135)		99321		2000.00	2200
54 Bromodichloromethane	83	6.502	6.508 (1.159)		634435		100.000	110
M 55 1,2-Dichloroethene (Total)	96				1077267		200.000	220
56 2-Chloroethyl vinyl ether	63	6.851	6.846 (1.222)		7679		100.000	88
57 cis-1,3-Dichloropropene	75	7.013	7.008 (1.251)		830294		100.000	110
58 4-Methyl-2-pentanone	43	7.188	7.182 (1.282)		778832		100.000	110
\$ 59 Toluene-d8	98	7.327	7.322 (0.807)		448197		50.0000	52
60 Toluene	91	7.397	7.403 (1.319)		1744568		100.000	110
61 trans-1,3-Dichloropropene	75	7.652	7.647 (1.365)		713389		100.000	110
62 Ethyl Methacrylate	69	7.768	7.775 (1.385)		507623		100.000	120
63 1,1,2-Trichloroethane	97	7.861	7.867 (1.402)		423910		100.000	110
64 Tetrachloroethene	164	8.059	8.053 (0.887)		369360		100.000	110
65 1,3-Dichloropropane	76	8.070	8.077 (0.889)		745505		100.000	110
66 2-Hexanone	43	8.175	8.181 (0.900)		550602		100.000	120
67 Dibromochloromethane	129	8.349	8.344 (0.919)		455390		100.000	120
68 1,2-Dibromoethane	107	8.489	8.483 (0.935)		486325		100.000	110
* 69 Chlorobenzene-d5	117	9.081	9.076 (1.000)		336924		50.0000	
70 1-Chlorohexane	91	9.104	9.099 (1.003)		644632		100.000	120
71 Chlorobenzene	112	9.116	9.110 (1.004)		1152835		100.000	120
72 1,1,1,2-Tetrachloroethane	131	9.220	9.227 (1.015)		394608		100.000	110
73 Ethylbenzene	106	9.267	9.273 (1.020)		597862		100.000	120
74 m,p-Xylene	106	9.418	9.424 (1.037)		1464909		200.000	230
75 o-Xylene	106	9.941	9.935 (1.095)		724382		100.000	120
76 Styrene	104	9.952	9.958 (1.096)		1115720		100.000	120
77 Bromoform	173	10.185	10.179 (1.122)		270165		100.000	120

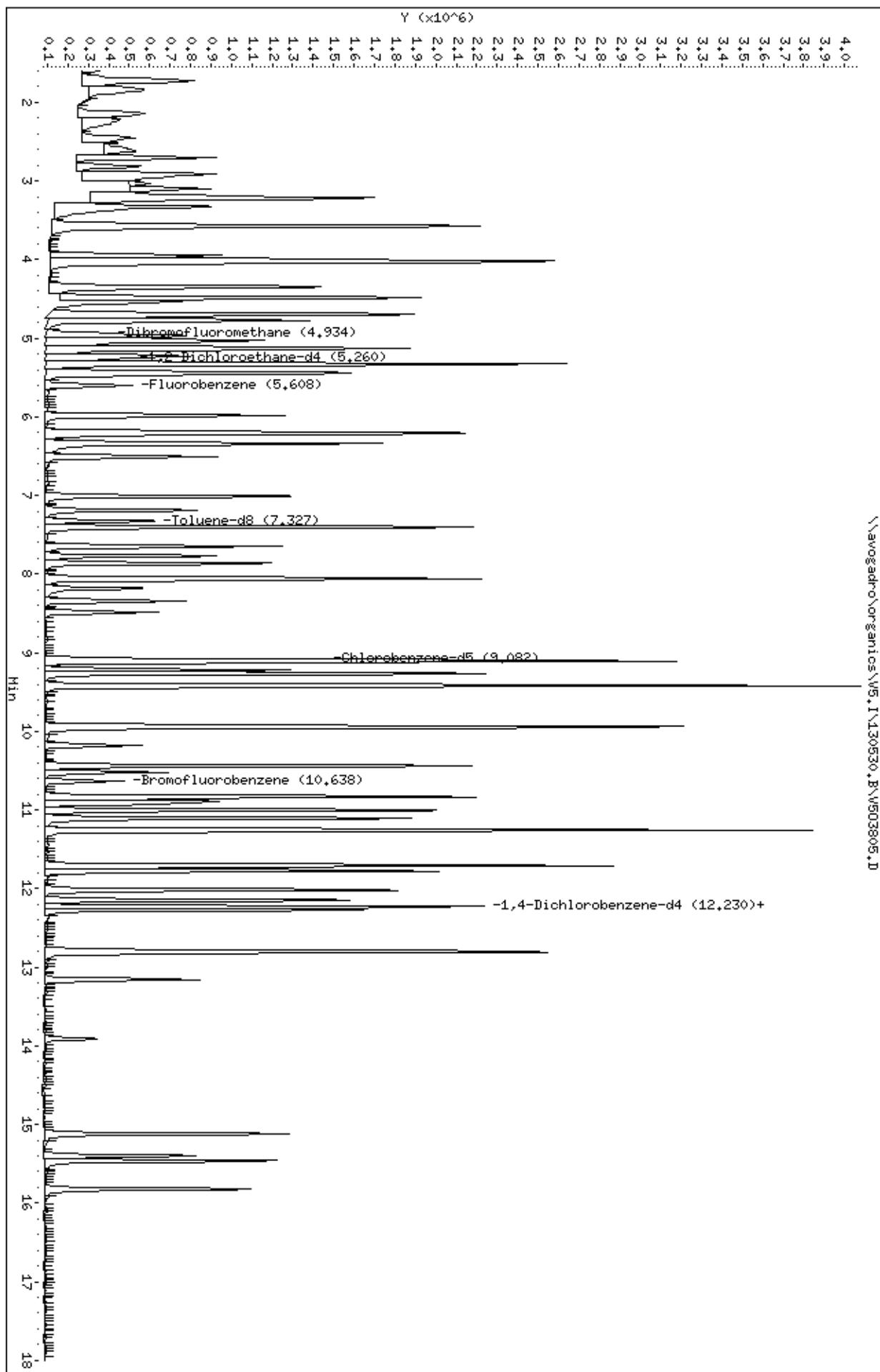
Compounds	QUANT SIG						AMOUNTS	
	MASS	RT	EXP RT	REL RT	RESPONSE		CAL-AMT (ug/L)	ON-COL (ug/L)
=====	=====	=====	=====	=====	=====		=====	=====
78 Isopropylbenzene	105	10.440	10.446	(1.150)	1770436		100.000	120
79 trans-1,4-Dichloro-2-butene	75	10.521	10.528	(1.159)	182650		100.000	160
\$ 80 Bromofluorobenzene	95	10.638	10.632	(1.171)	173633		50.0000	50
81 Bromobenzene	156	10.835	10.830	(0.885)	471234		100.000	120
82 1,1,2,2-Tetrachloroethane	83	10.835	10.841	(0.885)	640140		100.000	120
83 1,2,3-Trichloropropane	75	10.893	10.888	(0.890)	586294		100.000	120
84 n-Propylbenzene	120	11.009	11.004	(0.899)	442646		100.000	120
85 2-Chlorotoluene	126	11.102	11.109	(0.907)	423385		100.000	120
86 4-Chlorotoluene	126	11.253	11.260	(0.919)	441923		100.000	120
87 1,3,5-Trimethylbenzene	105	11.253	11.260	(0.919)	1380807		100.000	120
88 tert-Butylbenzene	119	11.706	11.713	(0.956)	1322381		100.000	120
89 1,2,4-Trimethylbenzene	105	11.776	11.771	(0.962)	1388282		100.000	120
90 sec-Butylbenzene	105	12.020	12.015	(0.982)	1736108		100.000	120
91 1,3-Dichlorobenzene	146	12.148	12.142	(0.992)	774758		100.000	120
92 4-Isopropyltoluene	119	12.229	12.235	(0.999)	1329087		100.000	120
* 93 1,4-Dichlorobenzene-d4	152	12.241	12.247	(1.000)	145240		50.0000	
94 1,4-Dichlorobenzene	146	12.276	12.270	(1.003)	797725		100.000	120
95 1,2-Dichlorobenzene	146	12.798	12.793	(1.046)	760213		100.000	120
M 96 Xylene (Total)	106				2189291		300.000	350
97 n-Butylbenzene	91	12.810	12.816	(1.047)	1254263		100.000	120
98 Hexachloroethane	117	13.158	13.165	(1.075)	223332		100.000	120
99 1,2-Dibromo-3-chloropropane	75	13.902	13.908	(1.136)	80771		100.000	100
100 1,3,5-Trichlorobenzene	182	15.122	15.128	(2.696)	476375		100.000	120
101 1,2,4-Trichlorobenzene	180	15.122	15.128	(1.235)	501678		100.000	120
102 Hexachlorobutadiene	225	15.400	15.407	(1.258)	200770		100.000	120
103 Naphthalene	128	15.459	15.465	(1.263)	1183537		100.000	120
104 1,2,3-Trichlorobenzene	180	15.819	15.825	(1.292)	430643		100.000	120

QC Flag Legend

T - Target compound detected outside RT window.

Data File: \\avogadro\organics\W5,I\130530.B\W503805.D
Date : 30-May-2013 14:32
Client ID: WSTD1005V
Sample Info: 5HL,WSTD1005V,WSTD1005V
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: ML
Column diameter: 0.25



Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
Data file : \\avogadro\organics\V5.I\130530.B\V503806.D
Lab Smp Id: VSTD2005V Client Smp ID: VSTD2005V
Inj Date : 30-MAY-2013 14:58
Operator : WL SRC: WL Inst ID: V5.i
Smp Info : 5ML,VSTD2005V,VSTD2005V
Misc Info :
Comment :
Method : \\avogadro\organics\V5.I\130530.B\v5_8260W.m
Meth Date : 03-Jun-2013 13:35 wluo Quant Type: ISTD
Cal Date : 30-MAY-2013 14:58 Cal File: V503806.D
Als bottle: 8 Calibration Sample, Level: 5
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: FULL.sub
Target Version: 4.14
Processing Host: TARGET103

Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG MASS					AMOUNTS	
		RT	EXP RT	REL RT	RESPONSE	CAL-AMT (ug/L)	ON-COL (ug/L)
1 Dichlorodifluoromethane	85	1.589	1.583 (0.283)		1024552	200.000	210 (A)
3 Chloromethane	50	1.729	1.734 (0.308)		2005056	200.000	220 (A)
4 Vinyl Chloride	62	1.833	1.850 (0.327)		1663708	200.000	220 (A)
5 Bromomethane	94	2.135	2.141 (0.381)		859337	200.000	190
6 Chloroethane	64	2.240	2.233 (0.399)		854717	200.000	210 (A)
7 Trichlorofluoromethane	101	2.449	2.454 (0.437)		1260546	200.000	220 (A)
8 Ethanol	46	2.623	2.617 (0.468)		733495	20000.0	22000 (A)
9 Ether	59	2.704	2.698 (0.482)		959581	200.000	230 (A)
10 Acrolein	56	2.809	2.814 (0.501)		1235422	1000.00	1100 (A)
11 1,1-Dichloroethene	96	2.902	2.907 (0.517)		958296	200.000	220 (A)
12 1,1,2-Trichloro-1,2,2-Trifluo	101	2.913	2.919 (0.520)		845427	200.000	220 (A)
13 Acetone	58	2.937	2.942 (0.524)		139045	200.000	200 (A)
14 Iodomethane	142	3.041	3.047 (0.542)		1643178	200.000	230 (A)
15 Carbon Disulfide	76	3.099	3.105 (0.553)		4104622	200.000	210 (A)
16 Acetonitrile	41	3.216	3.221 (0.573)		4073578	2000.00	2200 (A)
17 Allyl Chloride	39	3.216	3.221 (0.573)		1531346	200.000	210 (A)
18 Methyl Acetate	43	3.239	3.232 (0.578)		1413747	200.000	200 (A)
19 Methylene Chloride	84	3.332	3.325 (0.594)		1106408	200.000	210 (A)
20 tert-Butanol	59	3.436	3.430 (0.613)		227238	400.000	440 (A)
21 Acrylonitrile	53	3.541	3.534 (0.631)		470578	200.000	240 (A)
22 trans-1,2-Dichloroethene	96	3.576	3.569 (0.638)		981532	200.000	220 (A)
23 Methyl tert-butyl ether	73	3.576	3.581 (0.638)		2584617	200.000	220 (A)

Compounds	QUANT SIG						AMOUNTS	
	MASS	RT	EXP RT	REL RT	RESPONSE		CAL-AMT (ug/L)	ON-COL (ug/L)
=====	=====	=====	=====	=====	=====		=====	=====
24 1,1-Dichloroethane	63	3.947	3.941 (0.704)		2002226		200.000	220 (A)
25 Vinyl acetate	43	3.994	3.999 (0.712)		3836327		200.000	210 (A)
26 Diisopropyl Ether	45	4.017	4.011 (0.716)		4454319		200.000	210 (A)
27 2-Chloro-1,3-Butadiene	53	4.029	4.034 (0.718)		1489240		200.000	220 (A)
28 Ethyl tert-butyl ether	59	4.342	4.348 (0.774)		3340826		200.000	220 (A)
29 cis-1,2-Dichloroethene	96	4.482	4.475 (0.799)		1031346		200.000	200 (A)
30 2,2-Dichloropropane	77	4.482	4.487 (0.799)		994715		200.000	220 (A)
31 2-Butanone	72	4.493	4.487 (0.801)		125192		200.000	220 (A)
32 Propionitrile	54	4.540	4.534 (0.809)		1660189		2000.00	2200 (A)
33 Methacrylonitrile	41	4.691	4.685 (0.836)		1615495		400.000	430 (A)
34 Bromochloromethane	128	4.702	4.708 (0.838)		514023		200.000	220 (A)
35 Tetrahydrofuran	72	4.761	4.754 (0.849)		256800		400.000	440 (A)
36 Chloroform	83	4.772	4.777 (0.851)		1621994		200.000	220 (A)
\$ 37 Dibromofluoromethane	113	4.935	4.928 (0.880)		143501		50.0000	50
38 1,1,1-Trichloroethane	97	4.970	4.975 (0.886)		1112457		200.000	210 (A)
39 Cyclohexane	56	5.028	5.033 (0.896)		1765501		200.000	230 (A)
40 1,1-Dichloropropene	110	5.121	5.126 (0.913)		444043		200.000	220 (A)
41 Carbon Tetrachloride	117	5.132	5.138 (0.915)		915248		200.000	220 (A)
42 Isobutyl Alcohol	43	5.225	5.219 (0.932)		1120085		4000.00	4200 (A)
\$ 43 1,2-Dichloroethane-d4	102	5.260	5.265 (0.938)		32126		50.0000	50
44 Benzene	78	5.330	5.335 (0.950)		3529485		200.000	220 (A)
45 1,2-Dichloroethane	62	5.341	5.335 (0.952)		1261820		200.000	220 (A)
46 tert-Amyl methyl ether	73	5.446	5.440 (0.971)		2576898		200.000	220 (A)
* 47 Fluorobenzene	96	5.609	5.602 (1.000)		467336		50.0000	
48 Trichloroethene	130	5.992	5.986 (1.068)		921506		200.000	220 (A)
49 Methylcyclohexane	83	6.201	6.195 (1.106)		1029042		200.000	210 (A)
50 1,2-Dichloropropane	63	6.213	6.218 (1.108)		1119616		200.000	220 (A)
51 Dibromomethane	93	6.340	6.346 (1.130)		691605		200.000	220 (A)
52 Methyl Methacrylate	69	6.340	6.346 (1.130)		772499		200.000	220 (A)
53 1,4-Dioxane	88	6.364	6.357 (1.135)		176492		4000.00	4100 (A)
54 Bromodichloromethane	83	6.503	6.508 (1.159)		1216518		200.000	220 (A)
M 55 1,2-Dichloroethene (Total)	96				2012878		400.000	420
56 2-Chloroethyl vinyl ether	63	6.852	6.846 (1.222)		13000		200.000	160
57 cis-1,3-Dichloropropene	75	7.014	7.008 (1.251)		1573980		200.000	220 (A)
58 4-Methyl-2-pentanone	43	7.188	7.182 (1.282)		1481310		200.000	220 (A)
\$ 59 Toluene-d8	98	7.328	7.322 (0.807)		421483		50.0000	49
60 Toluene	91	7.397	7.403 (1.319)		3193851		200.000	210 (A)
61 trans-1,3-Dichloropropene	75	7.653	7.647 (1.365)		1380333		200.000	230 (A)
62 Ethyl Methacrylate	69	7.769	7.775 (1.385)		944985		200.000	230 (A)
63 1,1,2-Trichloroethane	97	7.874	7.867 (1.404)		820788		200.000	220 (A)
64 Tetrachloroethene	164	8.060	8.053 (0.887)		694707		200.000	220 (A)
65 1,3-Dichloropropane	76	8.071	8.077 (0.889)		1407049		200.000	220 (A)
66 2-Hexanone	43	8.176	8.181 (0.900)		999079		200.000	220 (A)
67 Dibromochloromethane	129	8.350	8.344 (0.919)		885519		200.000	240 (A)
68 1,2-Dibromoethane	107	8.489	8.483 (0.935)		945907		200.000	220 (A)
* 69 Chlorobenzene-d5	117	9.082	9.076 (1.000)		330520		50.0000	
70 1-Chlorohexane	91	9.105	9.099 (1.003)		1190137		200.000	220 (A)
71 Chlorobenzene	112	9.117	9.110 (1.004)		2131899		200.000	220 (A)
72 1,1,1,2-Tetrachloroethane	131	9.221	9.227 (1.015)		762243		200.000	220 (A)
73 Ethylbenzene	106	9.268	9.273 (1.020)		1152054		200.000	230 (A)
74 m,p-Xylene	106	9.419	9.424 (1.037)		2748390		400.000	440 (A)
75 o-Xylene	106	9.942	9.935 (1.095)		1394609		200.000	230 (A)
76 Styrene	104	9.953	9.958 (1.096)		2088598		200.000	230 (A)
77 Bromoform	173	10.185	10.179 (1.122)		538889		200.000	240 (A)

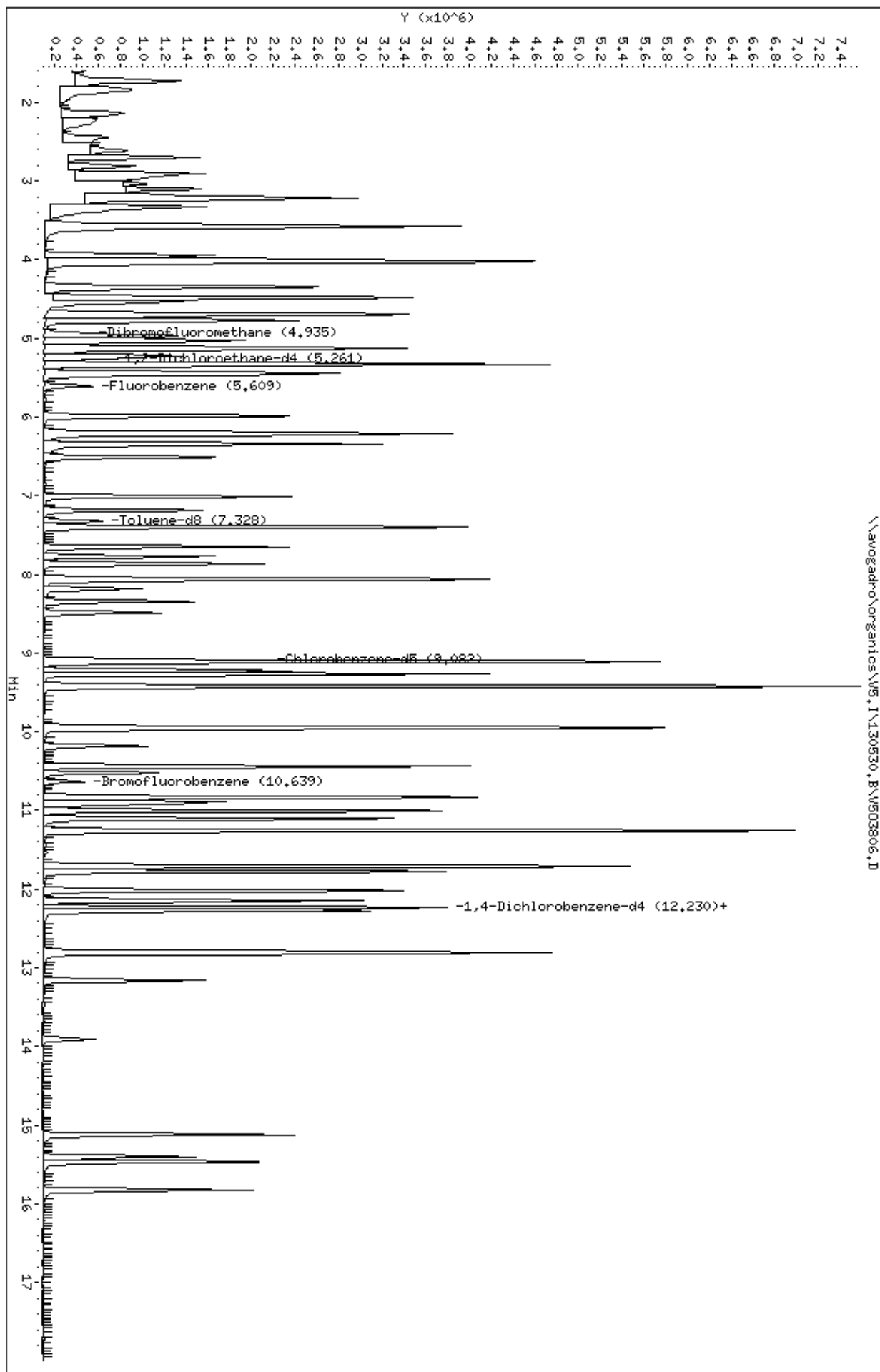
						AMOUNTS	
QUANT SIG						CAL-AMT	ON-COL
Compounds	MASS	RT	EXP RT	REL RT	RESPONSE	(ug/L)	(ug/L)
=====	=====	=====	=====	=====	=====	=====	=====
78 Isopropylbenzene	105	10.441	10.446	(1.150)	3281213	200.000	220(A)
79 trans-1,4-Dichloro-2-butene	75	10.522	10.528	(1.159)	311860	200.000	280(A)
\$ 80 Bromofluorobenzene	95	10.639	10.632	(1.171)	167726	50.0000	49
81 Bromobenzene	156	10.836	10.830	(0.885)	875905	200.000	220(A)
82 1,1,2,2-Tetrachloroethane	83	10.836	10.841	(0.885)	1194250	200.000	220(A)
83 1,2,3-Trichloropropane	75	10.894	10.888	(0.890)	1222015	200.000	240(A)
84 n-Propylbenzene	120	11.010	11.004	(0.899)	843627	200.000	230(A)
85 2-Chlorotoluene	126	11.115	11.109	(0.908)	809708	200.000	230(A)
86 4-Chlorotoluene	126	11.266	11.260	(0.920)	851957	200.000	230(A)
87 1,3,5-Trimethylbenzene	105	11.254	11.260	(0.919)	2561828	200.000	220(A)
88 tert-Butylbenzene	119	11.707	11.713	(0.956)	2456448	200.000	220(A)
89 1,2,4-Trimethylbenzene	105	11.777	11.771	(0.962)	2623493	200.000	220(A)
90 sec-Butylbenzene	105	12.021	12.015	(0.982)	3208280	200.000	220(A)
91 1,3-Dichlorobenzene	146	12.149	12.142	(0.992)	1481522	200.000	220(A)
92 4-Isopropyltoluene	119	12.230	12.235	(0.999)	2457017	200.000	220(A)
* 93 1,4-Dichlorobenzene-d4	152	12.242	12.247	(1.000)	146849	50.0000	
94 1,4-Dichlorobenzene	146	12.276	12.270	(1.003)	1505803	200.000	220(A)
95 1,2-Dichlorobenzene	146	12.799	12.793	(1.046)	1461689	200.000	230(A)
M 96 Xylene (Total)	106				4142999	600.000	670
97 n-Butylbenzene	91	12.811	12.816	(1.046)	2336754	200.000	220(A)
98 Hexachloroethane	117	13.159	13.165	(1.075)	436048	200.000	240(A)
99 1,2-Dibromo-3-chloropropane	75	13.914	13.908	(1.137)	153880	200.000	190
100 1,3,5-Trichlorobenzene	182	15.123	15.128	(2.696)	928046	200.000	240(A)
101 1,2,4-Trichlorobenzene	180	15.123	15.128	(1.235)	944724	200.000	230(A)
102 Hexachlorobutadiene	225	15.401	15.407	(1.258)	368145	200.000	220(A)
103 Naphthalene	128	15.471	15.465	(1.264)	2173009	200.000	230(A)
104 1,2,3-Trichlorobenzene	180	15.820	15.825	(1.292)	820918	200.000	230(A)

QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.

Data File: \\avogadro\organics\W5,I\130530.B\W503806.D
Date : 30-May-2013 14:58
Client ID: WSTD2005V
Sample Info: 5HL,WSTD2005V,WSTD2005V
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: ML
Column diameter: 0.25



Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130530.B\V503807.D
 Lab Smp Id: VICV0505V Client Smp ID: VICV0505V
 Inj Date : 30-MAY-2013 15:46
 Operator : WL SRC: WL Inst ID: V5.i
 Smp Info : 5ML,VICV0505V,VICV0505V
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130530.B\v5_8260W.m
 Meth Date : 03-Jun-2013 13:35 wluo Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V503805.D
 Als bottle: 10 Calibration Sample, Level: 3
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: FULL.sub
 Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG						AMOUNTS	
		MASS	RT	EXP RT	REL RT	RESPONSE	CAL-AMT (ug/L)	ON-COL (ug/L)
1 Dichlorodifluoromethane	85		1.586	1.583 (0.283)		341042	50.0000	41
3 Chloromethane	50		1.737	1.734 (0.310)		671838	50.0000	43
4 Vinyl Chloride	62		1.841	1.850 (0.329)		557906	50.0000	43
5 Bromomethane	94		2.143	2.141 (0.382)		349157	50.0000	44
6 Chloroethane	64		2.248	2.233 (0.401)		295529	50.0000	42
7 Trichlorofluoromethane	101		2.445	2.454 (0.436)		432044	50.0000	43
8 Ethanol	46		2.620	2.617 (0.467)		219462	5000.00	3800
9 Ether	59		2.701	2.698 (0.482)		323078	50.0000	44
10 Acrolein	56		2.817	2.814 (0.503)		385301	250.000	200
11 1,1-Dichloroethene	96		2.910	2.907 (0.519)		335171	50.0000	45
12 1,1,2-Trichloro-1,2,2-Trifluo	101		2.922	2.919 (0.521)		297868	50.0000	45
13 Acetone	58		2.945	2.942 (0.525)		45082	50.0000	38
14 Iodomethane	142		3.038	3.047 (0.542)		553448	50.0000	44
15 Carbon Disulfide	76		3.108	3.105 (0.554)		1412367	50.0000	42
16 Acetonitrile	41		3.224	3.221 (0.575)		1420327	500.000	430
17 Allyl Chloride	39		3.224	3.221 (0.575)		551550	50.0000	43 (T)
18 Methyl Acetate	43		3.235	3.232 (0.577)		486830	50.0000	40
19 Methylene Chloride	84		3.328	3.325 (0.594)		370410	50.0000	41
20 tert-Butanol	59		3.444	3.430 (0.615)		73719	100.000	82
21 Acrylonitrile	53		3.537	3.534 (0.631)		163308	50.0000	48
22 trans-1,2-Dichloroethene	96		3.572	3.569 (0.637)		334185	50.0000	42
23 Methyl tert-butyl ether	73		3.584	3.581 (0.639)		887490	50.0000	43
24 1,1-Dichloroethane	63		3.944	3.941 (0.704)		690524	50.0000	43

Compounds	QUANT	SIG						AMOUNTS	
			MASS	RT	EXP RT	REL RT	RESPONSE	CAL-AMT (ug/L)	ON-COL (ug/L)
=====	=====	=====	=====	=====	=====	=====	=====	=====	=====
25 Vinyl acetate	43		4.002	3.999	(0.714)		1398121	50.0000	44
26 Diisopropyl Ether	45		4.014	4.011	(0.716)		1583931	50.0000	42
27 2-Chloro-1,3-Butadiene	53		4.037	4.034	(0.720)		517269	50.0000	44
28 Ethyl tert-butyl ether	59		4.351	4.348	(0.776)		1147254	50.0000	43
29 cis-1,2-Dichloroethene	96		4.478	4.475	(0.799)		350996	50.0000	42
30 2,2-Dichloropropane	77		4.490	4.487	(0.801)		324602	50.0000	42
31 2-Butanone	72		4.490	4.487	(0.801)		45367	50.0000	46
32 Propionitrile	54		4.536	4.534	(0.809)		567513	500.000	440
33 Methacrylonitrile	41		4.687	4.685	(0.836)		588701	100.000	90
34 Bromochloromethane	128		4.711	4.708	(0.840)		172339	50.0000	43
35 Tetrahydrofuran	72		4.757	4.754	(0.849)		89936	100.000	90
36 Chloroform	83		4.780	4.777	(0.853)		544998	50.0000	42
\$ 37 Dibromofluoromethane	113		4.931	4.928	(0.880)		251556	50.0000	50
38 1,1,1-Trichloroethane	97		4.978	4.975	(0.888)		369436	50.0000	41
39 Cyclohexane	56		5.036	5.033	(0.898)		553651	50.0000	42
40 1,1-Dichloropropene	110		5.129	5.126	(0.915)		152719	50.0000	44
41 Carbon Tetrachloride	117		5.141	5.138	(0.917)		305718	50.0000	43
42 Isobutyl Alcohol	43		5.222	5.219	(0.932)		390723	1000.00	850
\$ 43 1,2-Dichloroethane-d4	102		5.268	5.265	(0.940)		54610	50.0000	49
44 Benzene	78		5.326	5.335	(0.950)		1220432	50.0000	44
45 1,2-Dichloroethane	62		5.338	5.335	(0.952)		433136	50.0000	43
46 tert-Amyl methyl ether	73		5.443	5.440	(0.971)		884280	50.0000	44
* 47 Fluorobenzene	96		5.605	5.602	(1.000)		811616	50.0000	
48 Trichloroethene	130		5.989	5.986	(1.068)		316227	50.0000	44
49 Methylcyclohexane	83		6.198	6.195	(1.106)		386007	50.0000	45
50 1,2-Dichloropropane	63		6.221	6.218	(1.110)		384907	50.0000	43
51 Dibromomethane	93		6.337	6.346	(1.131)		233911	50.0000	44
52 Methyl Methacrylate	69		6.337	6.346	(1.131)		263778	50.0000	43
53 1,4-Dioxane	88		6.360	6.357	(1.135)		57508	1000.00	770
54 Bromodichloromethane	83		6.511	6.508	(1.162)		416670	50.0000	44
M 55 1,2-Dichloroethene (Total)	96						685181	100.000	85
56 2-Chloroethyl vinyl ether	63		6.848	6.846	(1.222)		6634	50.0000	46
57 cis-1,3-Dichloropropene	75		7.011	7.008	(1.251)		535829	50.0000	44
58 4-Methyl-2-pentanone	43		7.185	7.182	(1.282)		515678	50.0000	44
\$ 59 Toluene-d8	98		7.324	7.322	(0.807)		740306	50.0000	49
60 Toluene	91		7.406	7.403	(1.321)		1132118	50.0000	44
61 trans-1,3-Dichloropropene	75		7.650	7.647	(1.365)		472791	50.0000	45
62 Ethyl Methacrylate	69		7.777	7.775	(1.388)		322740	50.0000	46
63 1,1,2-Trichloroethane	97		7.870	7.867	(1.404)		280389	50.0000	43
64 Tetrachloroethene	164		8.056	8.053	(0.887)		239622	50.0000	43
65 1,3-Dichloropropane	76		8.068	8.077	(0.889)		485948	50.0000	43
66 2-Hexanone	43		8.184	8.181	(0.901)		355325	50.0000	44
67 Dibromochloromethane	129		8.347	8.344	(0.919)		305134	50.0000	46
68 1,2-Dibromoethane	107		8.486	8.483	(0.935)		322360	50.0000	44
* 69 Chlorobenzene-d5	117		9.079	9.076	(1.000)		582029	50.0000	
70 1-Chlorohexane	91		9.102	9.099	(1.003)		414846	50.0000	44
71 Chlorobenzene	112		9.113	9.110	(1.004)		743245	50.0000	44
72 1,1,1,2-Tetrachloroethane	131		9.218	9.227	(1.015)		264563	50.0000	44
73 Ethylbenzene	106		9.264	9.273	(1.020)		383124	50.0000	43
74 m,p-Xylene	106		9.427	9.424	(1.038)		940429	100.000	86
75 o-Xylene	106		9.938	9.935	(1.095)		468851	50.0000	44
76 Styrene	104		9.961	9.958	(1.097)		729098	50.0000	46
77 Bromoform	173		10.182	10.179	(1.122)		190295	50.0000	47
78 Isopropylbenzene	105		10.438	10.446	(1.150)		1154431	50.0000	44

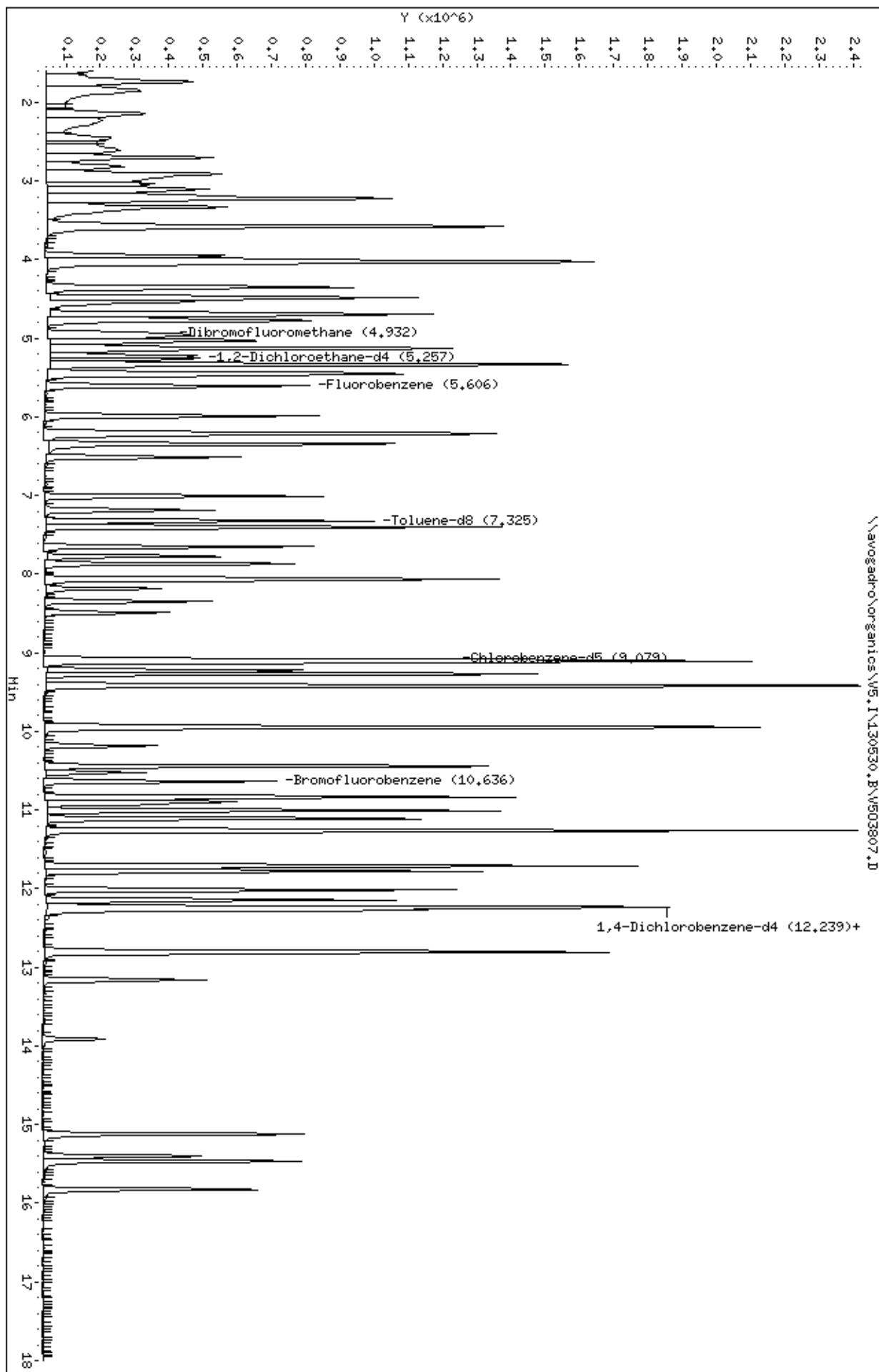
Compounds	QUANT SIG						AMOUNTS	
	MASS	RT	EXP RT	REL RT	RESPONSE		CAL-AMT (ug/L)	ON-COL (ug/L)
=====	=====	=====	=====	=====	=====		=====	=====
79 trans-1,4-Dichloro-2-butene	75	10.519	10.528	(1.159)	90244		50.0000	45
\$ 80 Bromofluorobenzene	95	10.635	10.632	(1.171)	295877		50.0000	49
81 Bromobenzene	156	10.833	10.830	(0.885)	299808		50.0000	43
82 1,1,2,2-Tetrachloroethane	83	10.844	10.841	(0.886)	418252		50.0000	43
83 1,2,3-Trichloropropane	75	10.891	10.888	(0.890)	399463		50.0000	46
84 n-Propylbenzene	120	11.007	11.004	(0.899)	281744		50.0000	44
85 2-Chlorotoluene	126	11.111	11.109	(0.908)	273548		50.0000	43
86 4-Chlorotoluene	126	11.262	11.260	(0.920)	283693		50.0000	43
87 1,3,5-Trimethylbenzene	105	11.262	11.260	(0.920)	907977		50.0000	44
88 tert-Butylbenzene	119	11.704	11.713	(0.956)	863054		50.0000	44
89 1,2,4-Trimethylbenzene	105	11.774	11.771	(0.962)	918274		50.0000	44
90 sec-Butylbenzene	105	12.018	12.015	(0.982)	1138751		50.0000	44
91 1,3-Dichlorobenzene	146	12.145	12.142	(0.992)	517171		50.0000	44
92 4-Isopropyltoluene	119	12.227	12.235	(0.999)	852985		50.0000	44
* 93 1,4-Dichlorobenzene-d4	152	12.238	12.247	(1.000)	258983		50.0000	(Q)
94 1,4-Dichlorobenzene	146	12.273	12.270	(1.003)	521029		50.0000	43
95 1,2-Dichlorobenzene	146	12.796	12.793	(1.046)	491562		50.0000	43
M 96 Xylene (Total)	106				1409280		150.000	130
97 n-Butylbenzene	91	12.819	12.816	(1.047)	827011		50.0000	43
98 Hexachloroethane	117	13.156	13.165	(1.075)	153483		50.0000	48
99 1,2-Dibromo-3-chloropropane	75	13.911	13.908	(1.137)	54869		50.0000	39
100 1,3,5-Trichlorobenzene	182	15.119	15.128	(2.697)	305952		50.0000	45
101 1,2,4-Trichlorobenzene	180	15.119	15.128	(1.235)	317598		50.0000	44
102 Hexachlorobutadiene	225	15.398	15.407	(1.258)	127370		50.0000	44
103 Naphthalene	128	15.468	15.465	(1.264)	745626		50.0000	44
104 1,2,3-Trichlorobenzene	180	15.828	15.825	(1.293)	276293		50.0000	44

QC Flag Legend

T - Target compound detected outside RT window.
 Q - Qualifier signal failed the ratio test.

Data File: \\avogadro\organics\W5,I\130530.B\W503807.D
Date : 30-May-2013 15:46
Client ID: VICV0505V
Sample Info: 5ML,VICV0505V,VICV0505V
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: ML
Column diameter: 0.25



Spectrum Analytical, Inc. RI Division

Data file : \\avogadro\organics\V5.I\130607.B\V503990.D
 Lab Smp Id: BFBD5 Client Smp ID: BFBD5
 Inj Date : 07-JUN-2013 13:58
 Operator : WL SRC: WL Inst ID: V5.i
 Smp Info : 2UL,BFBD5,BFBD5
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\bfb8260.m
 Meth Date : 30-May-2013 14:16 wluo Quant Type: ISTD
 Cal Date : 21-MAR-2011 17:17 Cal File: V5M6831.D
 Als bottle: 2 QC Sample: BFB
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: all.sub
 Target Version: 4.14 Sample Matrix: WATER
 Processing Host: TARGET103

Concentration Formula: Amt * DF * Uf * Vf * VI * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vf	1.000	Volumetric correction factor
VI	1.000	Injection Volume
Cpnd Variable		Local Compound Variable

CONCENTRATIONS									
				ON-COL		FINAL			
RT	EXP RT	REL RT	MASS	RESPONSE	(ug/L)	(ug/L)	TARGET	RANGE	RATIO
=====	=====	=====	=====	=====	=====	=====	=====	=====	=====
1 bfb				CAS #: 460-00-4					
10.635	11.000	(0.000)	95	139008			0.00-	100.00	100.00
10.635	11.000	(0.000)	50	34504			15.00-	40.00	24.82
10.635	11.000	(0.000)	75	62296			30.00-	60.00	44.81
10.635	11.000	(0.000)	96	8712			5.00-	9.00	6.27
10.635	11.000	(0.000)	173	0	0.0	0.0	0.00-	2.00	0.00
10.635	11.000	(0.000)	174	100312			50.00-	100.00	72.16
10.635	11.000	(0.000)	175	7993			5.00-	9.00	7.97
10.635	11.000	(0.000)	176	96984			95.00-	101.00	96.68
10.635	11.000	(0.000)	177	6155			5.00-	9.00	6.35

Date : 07-JUN-2013 13:58

Client ID: BFB05

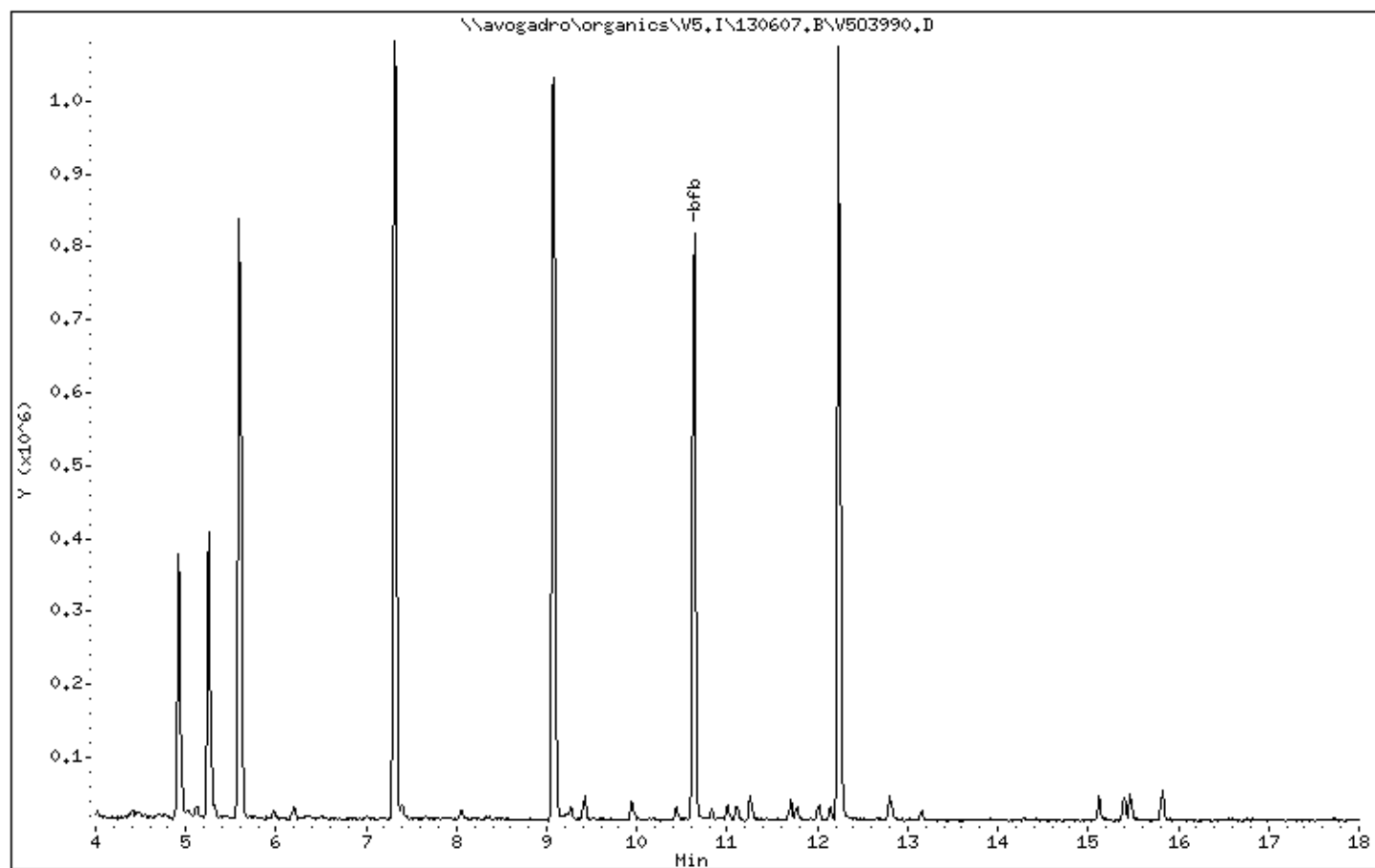
Instrument: V5.i

Sample Info: 2UL,BFB05,BFB05

Operator: WL SRC: WL

Column phase: DB-624

Column diameter: 0.25



Date : 07-JUN-2013 13:58

Client ID: BFBD5

Instrument: V5.i

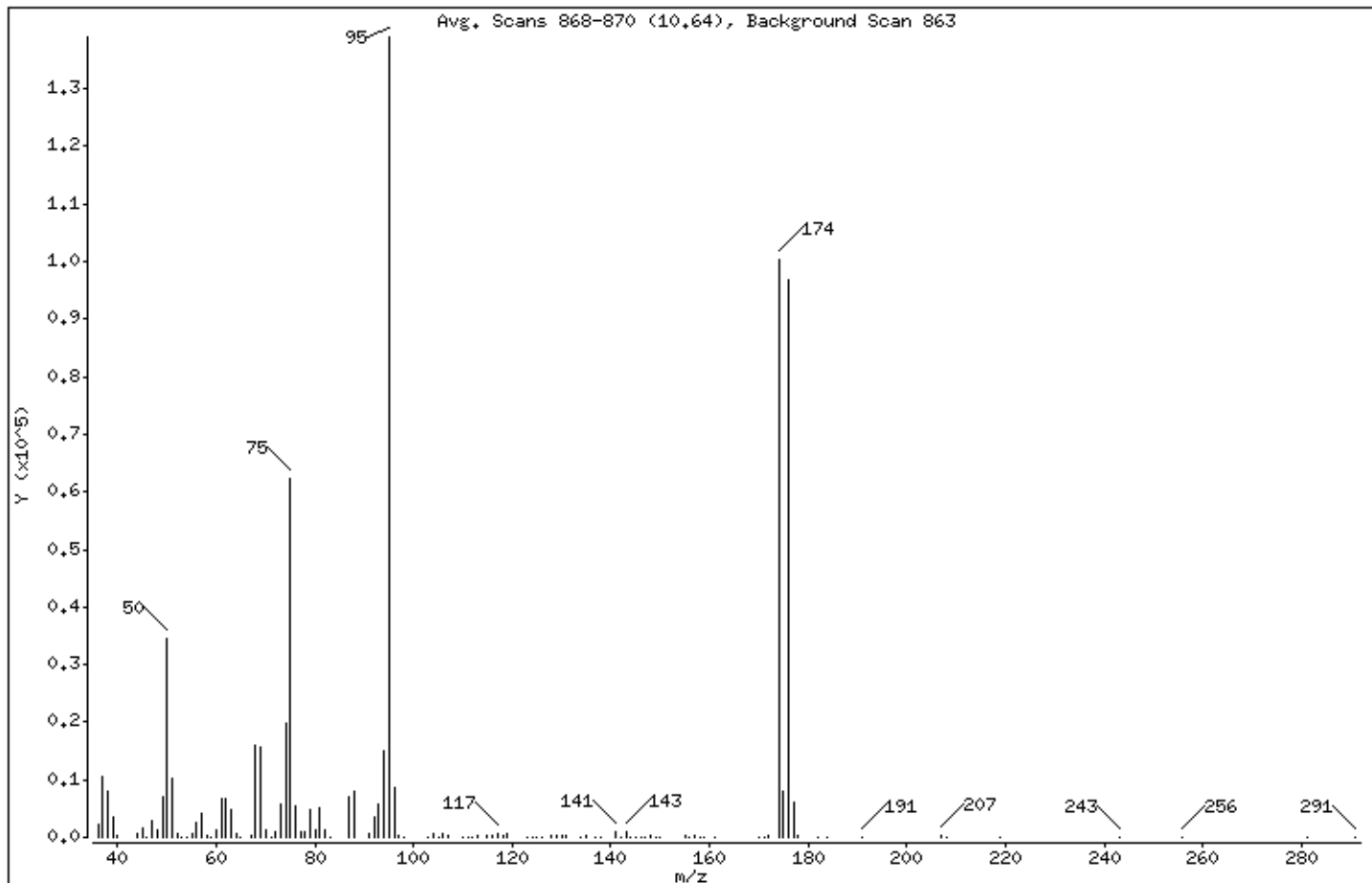
Sample Info: 2UL,BFBD5,BFBD5

Operator: WL SRC: WL

Column phase: DB-624

Column diameter: 0.25

1 bfb



m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
95	Base Peak, 100% relative abundance	100.00
50	15.00 - 40.00% of mass 95	24.82
75	30.00 - 60.00% of mass 95	44.81
96	5.00 - 9.00% of mass 95	6.27
173	Less than 2.00% of mass 174	0.00 (0.00)
174	50.00 - 100.00% of mass 95	72.16
175	5.00 - 9.00% of mass 174	5.75 (7.97)
176	95.00 - 101.00% of mass 174	69.77 (96.68)
177	5.00 - 9.00% of mass 176	4.43 (6.35)

Date : 07-JUN-2013 13:58

Client ID: BFBDS

Instrument: V5.i

Sample Info: 2UL,BFBDS,BFBDS

Operator: WL SRC: WL

Column phase: DB-624

Column diameter: 0.25

Data File: V503990.D

Spectrum: Avg. Scans 868-870 (10.64), Background Scan 863

Location of Maximum: 95.00

Number of points: 114

m/z	Y	m/z	Y	m/z	Y	m/z	Y

36.00	2228	69.00	15687	107.00	179	148.00	209
37.00	10645	70.00	1119	110.00	46	149.00	125
38.00	8146	71.00	69	111.00	112	150.00	51
39.00	3442	72.00	883	112.00	105	155.00	239
40.00	291	73.00	5749	113.00	163	156.00	60

44.00	775	74.00	19968	115.00	219	157.00	180
45.00	1664	75.00	62296	116.00	452	158.00	41
46.00	109	76.00	5319	117.00	763	159.00	146
47.00	3001	77.00	1046	118.00	463	161.00	54
48.00	1303	78.00	853	119.00	663	170.00	48

49.00	6946	79.00	4908	123.00	36	171.00	54
50.00	34504	80.00	1401	124.00	35	172.00	183
51.00	10215	81.00	5101	125.00	39	174.00	100312
52.00	548	82.00	1190	126.00	37	175.00	7993
53.00	34	83.00	69	128.00	448	176.00	96984

54.00	6	87.00	7159	129.00	264	177.00	6155
55.00	583	88.00	7890	130.00	426	178.00	168
56.00	2587	91.00	622	131.00	183	182.00	36
57.00	4300	92.00	3389	134.00	45	184.00	34
58.00	257	93.00	5689	135.00	220	191.00	88

59.00	34	94.00	14875	137.00	134	207.00	216
60.00	1343	95.00	139008	138.00	35	208.00	25
61.00	6837	96.00	8712	141.00	1029	219.00	27
62.00	6642	97.00	289	142.00	43	243.00	35
63.00	4867	98.00	36	143.00	846	256.00	45

64.00	488	103.00	139	144.00	122	281.00	19
65.00	94	104.00	580	145.00	80	291.00	39
67.00	369	105.00	107	146.00	95		
68.00	15875	106.00	571	147.00	93		

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130607.B\V503991.D
 Lab Smp Id: VSTD050D5 Client Smp ID: VSTD050D5
 Inj Date : 07-JUN-2013 14:23
 Operator : WL SRC: WL Inst ID: V5.i
 Smp Info : 5ML,VSTD050D5,VSTD050D5
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
 Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V503805.D
 Als bottle: 3 Continuing Calibration Sample
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: FULL.sub
 Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG						AMOUNTS	
		MASS	RT	EXP RT	REL RT	RESPONSE	CAL-AMT (ug/L)	ON-COL (ug/L)
1 Dichlorodifluoromethane	85		1.584	1.584 (0.283)		338761	50.0000	42
3 Chloromethane	50		1.735	1.735 (0.310)		796233	50.0000	52
4 Vinyl Chloride	62		1.839	1.839 (0.328)		694425	50.0000	54
5 Bromomethane	94		2.142	2.142 (0.382)		446582	50.0000	58
6 Chloroethane	64		2.234	2.234 (0.399)		391031	50.0000	57
7 Trichlorofluoromethane	101		2.444	2.444 (0.436)		558847	50.0000	57
8 Ethanol	46		2.618	2.618 (0.467)		213695	5000.00	3800(Q)
9 Ether	59		2.699	2.699 (0.482)		383762	50.0000	53(Q)
10 Acrolein	56		2.804	2.804 (0.500)		485003	250.000	260
11 1,1-Dichloroethene	96		2.908	2.908 (0.519)		375769	50.0000	51
12 1,1,2-Trichloro-1,2,2-Trifluo	101		2.920	2.920 (0.521)		364653	50.0000	57
13 Acetone	58		2.943	2.943 (0.525)		51175	50.0000	44
14 Iodomethane	142		3.036	3.036 (0.542)		617661	50.0000	51
15 Carbon Disulfide	76		3.106	3.106 (0.554)		1359002	50.0000	41
16 Acetonitrile	41		3.210	3.210 (0.573)		1657781	500.000	520
17 Allyl Chloride	39		3.210	3.210 (0.573)		710159	50.0000	57(Q)
18 Methyl Acetate	43		3.233	3.233 (0.577)		560760	50.0000	47
19 Methylene Chloride	84		3.326	3.326 (0.594)		450027	50.0000	51
20 tert-Butanol	59		3.431	3.431 (0.612)		70631	100.000	80
21 Acrylonitrile	53		3.536	3.536 (0.631)		174895	50.0000	52
22 trans-1,2-Dichloroethene	96		3.570	3.570 (0.637)		394950	50.0000	51
23 Methyl tert-butyl ether	73		3.570	3.570 (0.637)		991919	50.0000	49
24 1,1-Dichloroethane	63		3.942	3.942 (0.704)		859740	50.0000	55

Compounds	QUANT SIG						AMOUNTS	
	MASS	RT	EXP RT	REL RT	RESPONSE		CAL-AMT (ug/L)	ON-COL (ug/L)
=====	=====	=====	=====	=====	=====		=====	=====
25 Vinyl acetate	43	4.000	4.000	(0.714)	1645563		50.0000	53
26 Diisopropyl Ether	45	4.012	4.012	(0.716)	2089299		50.0000	57
27 2-Chloro-1,3-Butadiene	53	4.035	4.035	(0.720)	674178		50.0000	58
28 Ethyl tert-butyl ether	59	4.349	4.349	(0.776)	1378756		50.0000	53
29 cis-1,2-Dichloroethene	96	4.476	4.476	(0.799)	428272		50.0000	53
30 2,2-Dichloropropane	77	4.488	4.488	(0.801)	433366		50.0000	57
31 2-Butanone	72	4.488	4.488	(0.801)	45922		50.0000	48
32 Propionitrile	54	4.535	4.535	(0.809)	590500		500.000	470
33 Methacrylonitrile	41	4.686	4.686	(0.836)	651752		100.000	100
34 Bromochloromethane	128	4.697	4.697	(0.838)	202505		50.0000	52
35 Tetrahydrofuran	72	4.755	4.755	(0.849)	88505		100.000	90
36 Chloroform	83	4.778	4.778	(0.853)	702256		50.0000	55
\$ 37 Dibromofluoromethane	113	4.929	4.929	(0.880)	258246		50.0000	52
38 1,1,1-Trichloroethane	97	4.964	4.964	(0.886)	469246		50.0000	53
39 Cyclohexane	56	5.022	5.022	(0.896)	728135		50.0000	56
40 1,1-Dichloropropene	110	5.127	5.127	(0.915)	183643		50.0000	54
41 Carbon Tetrachloride	117	5.127	5.127	(0.915)	385530		50.0000	55
42 Isobutyl Alcohol	43	5.220	5.220	(0.932)	383487		1000.00	850
\$ 43 1,2-Dichloroethane-d4	102	5.255	5.255	(0.938)	51958		50.0000	47
44 Benzene	78	5.324	5.324	(0.950)	1496829		50.0000	54
45 1,2-Dichloroethane	62	5.336	5.336	(0.952)	553748		50.0000	56
46 tert-Amyl methyl ether	73	5.441	5.441	(0.971)	1023543		50.0000	52
* 47 Fluorobenzene	96	5.603	5.603	(1.000)	793954		50.0000	
48 Trichloroethene	130	5.987	5.987	(1.068)	348597		50.0000	50
49 Methylcyclohexane	83	6.196	6.196	(1.106)	486552		50.0000	58
50 1,2-Dichloropropane	63	6.219	6.219	(1.110)	485645		50.0000	56
51 Dibromomethane	93	6.335	6.335	(1.131)	284548		50.0000	54
52 Methyl Methacrylate	69	6.335	6.335	(1.131)	277723		50.0000	46
53 1,4-Dioxane	88	6.358	6.358	(1.135)	47679		1000.00	650
54 Bromodichloromethane	83	6.509	6.509	(1.162)	515231		50.0000	56
56 2-Chloroethyl vinyl ether	63	7.009	7.009	(1.251)	4616		50.0000	32(T)
57 cis-1,3-Dichloropropene	75	7.009	7.009	(1.251)	660752		50.0000	55
58 4-Methyl-2-pentanone	43	7.183	7.183	(1.282)	571235		50.0000	50
\$ 59 Toluene-d8	98	7.323	7.323	(0.807)	798371		50.0000	46
60 Toluene	91	7.404	7.404	(1.321)	1389711		50.0000	55
61 trans-1,3-Dichloropropene	75	7.648	7.648	(1.365)	563851		50.0000	55
62 Ethyl Methacrylate	69	7.776	7.776	(1.388)	314167		50.0000	45
63 1,1,2-Trichloroethane	97	7.868	7.868	(1.404)	323336		50.0000	50
64 Tetrachloroethene	164	8.054	8.054	(0.887)	258063		50.0000	41
65 1,3-Dichloropropane	76	8.066	8.066	(0.889)	590992		50.0000	45
66 2-Hexanone	43	8.182	8.182	(0.901)	398133		50.0000	44
67 Dibromochloromethane	129	8.345	8.345	(0.919)	332509		50.0000	44
68 1,2-Dibromoethane	107	8.484	8.484	(0.935)	363665		50.0000	43
* 69 Chlorobenzene-d5	117	9.077	9.077	(1.000)	665170		50.0000	
70 1-Chlorohexane	91	9.100	9.100	(1.003)	505622		50.0000	47
71 Chlorobenzene	112	9.111	9.111	(1.004)	899182		50.0000	46
72 1,1,1,2-Tetrachloroethane	131	9.216	9.216	(1.015)	303216		50.0000	44
73 Ethylbenzene	106	9.262	9.262	(1.020)	461777		50.0000	46
74 m,p-Xylene	106	9.413	9.413	(1.037)	1129278		100.000	91
75 o-Xylene	106	9.936	9.936	(1.095)	573873		50.0000	47
76 Styrene	104	9.959	9.959	(1.097)	883553		50.0000	49
77 Bromoform	173	10.180	10.180	(1.122)	195494		50.0000	42
78 Isopropylbenzene	105	10.436	10.436	(1.150)	1396096		50.0000	46
79 trans-1,4-Dichloro-2-butene	75	10.517	10.517	(1.159)	83298		50.0000	36

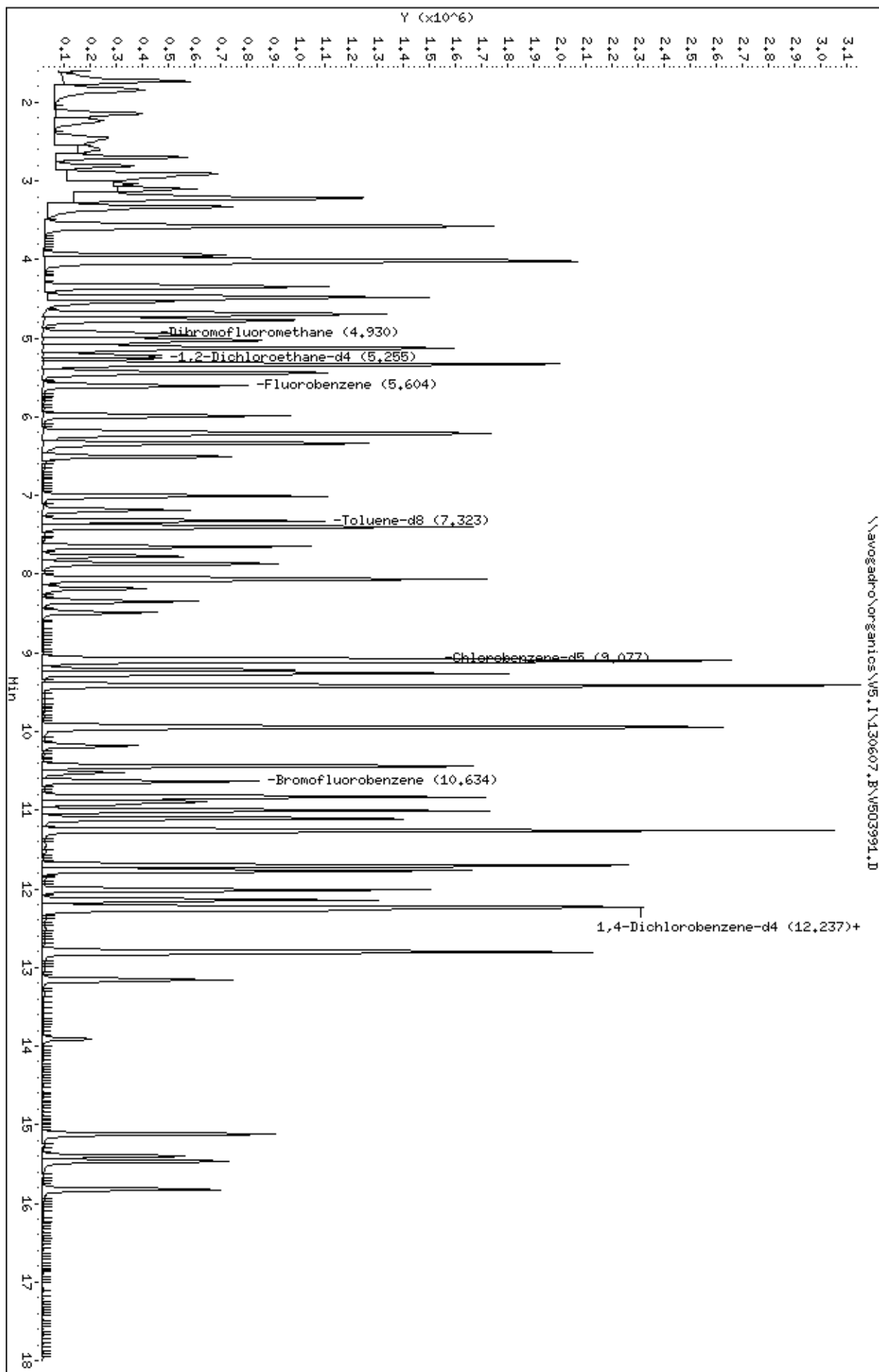
Compounds	QUANT SIG						AMOUNTS	
	MASS	RT	EXP RT	REL RT	RESPONSE		CAL-AMT (ug/L)	ON-COL (ug/L)
=====	=====	=====	=====	=====	=====		=====	=====
\$ 80 Bromofluorobenzene	95	10.633	10.633	(1.171)	367513		50.0000	53
81 Bromobenzene	156	10.831	10.831	(0.885)	348210		50.0000	42
82 1,1,2,2-Tetrachloroethane	83	10.842	10.842	(0.886)	468795		50.0000	40
83 1,2,3-Trichloropropane	75	10.889	10.889	(0.890)	418563		50.0000	40
84 n-Propylbenzene	120	11.005	11.005	(0.899)	339563		50.0000	44
85 2-Chlorotoluene	126	11.110	11.110	(0.908)	318430		50.0000	42
86 4-Chlorotoluene	126	11.261	11.261	(0.920)	344248		50.0000	43
87 1,3,5-Trimethylbenzene	105	11.261	11.261	(0.920)	1090956		50.0000	44
88 tert-Butylbenzene	119	11.702	11.702	(0.956)	1036050		50.0000	44
89 1,2,4-Trimethylbenzene	105	11.772	11.772	(0.962)	1114064		50.0000	44
90 sec-Butylbenzene	105	12.016	12.016	(0.982)	1357716		50.0000	44
91 1,3-Dichlorobenzene	146	12.143	12.143	(0.992)	596044		50.0000	42
92 4-Isopropyltoluene	119	12.225	12.225	(0.999)	1026274		50.0000	44
* 93 1,4-Dichlorobenzene-d4	152	12.236	12.236	(1.000)	311190		50.0000	
94 1,4-Dichlorobenzene	146	12.271	12.271	(1.003)	617182		50.0000	42
95 1,2-Dichlorobenzene	146	12.794	12.794	(1.046)	580966		50.0000	42
97 n-Butylbenzene	91	12.817	12.817	(1.047)	1053736		50.0000	46
98 Hexachloroethane	117	13.154	13.154	(1.075)	214938		50.0000	56
99 1,2-Dibromo-3-chloropropane	75	13.909	13.909	(1.137)	54682		50.0000	32
100 1,3,5-Trichlorobenzene	182	15.117	15.117	(2.698)	328024		50.0000	50
101 1,2,4-Trichlorobenzene	180	15.117	15.117	(1.235)	339542		50.0000	39
102 Hexachlorobutadiene	225	15.408	15.408	(1.259)	135098		50.0000	39
103 Naphthalene	128	15.466	15.466	(1.264)	713456		50.0000	35
104 1,2,3-Trichlorobenzene	180	15.826	15.826	(1.293)	272363		50.0000	36

QC Flag Legend

T - Target compound detected outside RT window.
 Q - Qualifier signal failed the ratio test.

Data File: \\avogadro\organics\W5.1\130607.B\W503991.D
Date : 07-JUN-2013 14:23
Client ID: WSTD050D5
Sample Info: 5HL,WSTD050D5,WSTD050D5
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: ML
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V503993.D
 Report Date: 11-Jun-2013 09:23

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130607.B\V503993.D
 Lab Smp Id: MB-72123 Client Smp ID: VBLKD5
 Inj Date : 07-JUN-2013 15:14
 Operator : WL SRC: LIMS Inst ID: V5.i
 Smp Info : 5ML,MB-72123,VBLKD5,72123
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
 Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V503805.D
 Als bottle: 5 QC Sample: BLANK
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: FULL.sub
 Target Version: 4.14

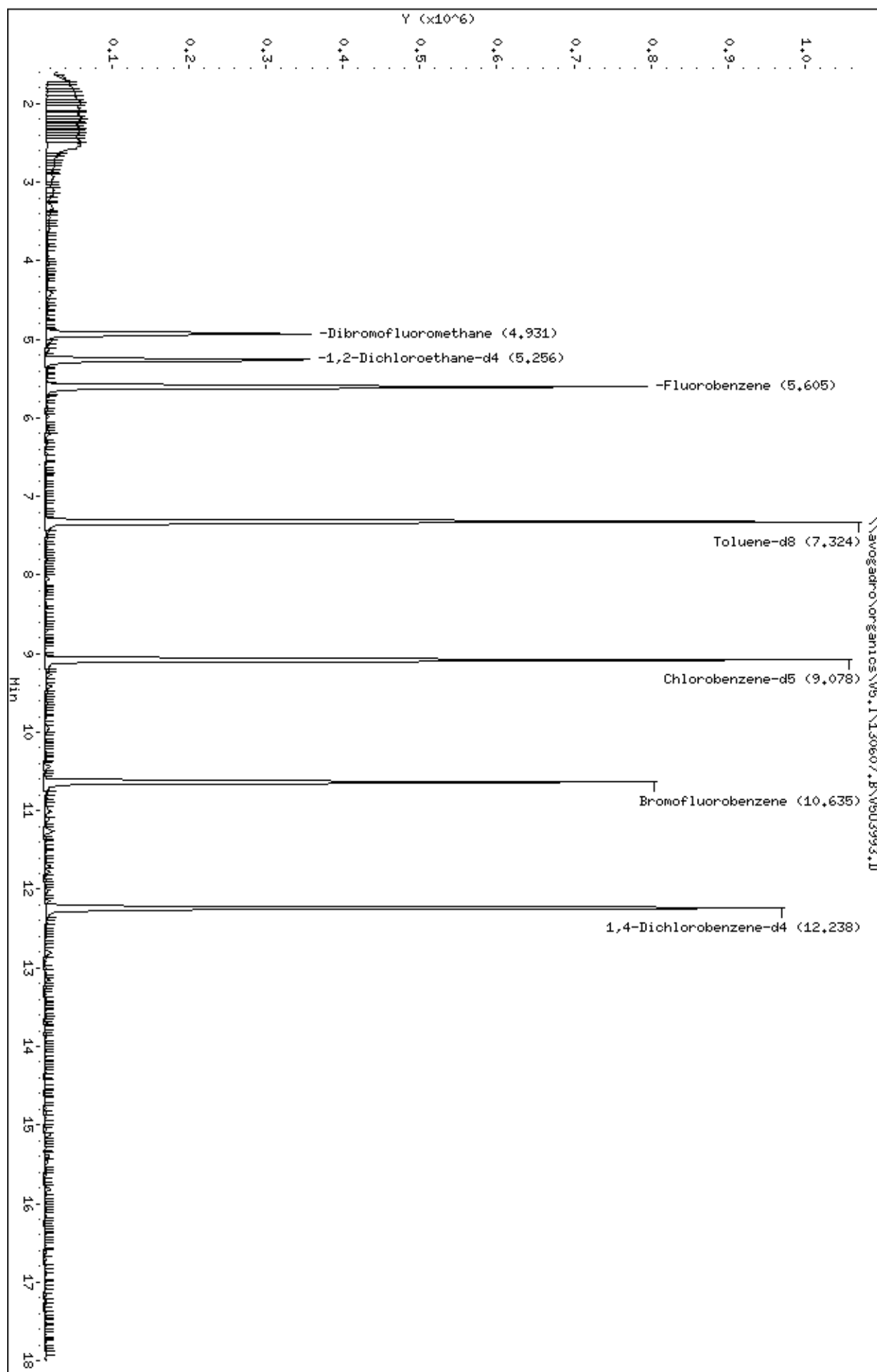
Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS	
						ON-COLUMN (ug/L)	FINAL (ug/L)
\$ 37 Dibromofluoromethane	113	4.930	4.929	(0.880)	257142	53.1459	53
\$ 43 1,2-Dichloroethane-d4	102	5.267	5.255	(0.940)	54625	50.6938	51
* 47 Fluorobenzene	96	5.604	5.603	(1.000)	779985	50.0000	
\$ 59 Toluene-d8	98	7.323	7.323	(0.807)	778971	46.4784	46
* 69 Chlorobenzene-d5	117	9.077	9.077	(1.000)	649391	50.0000	
\$ 80 Bromofluorobenzene	95	10.634	10.633	(1.171)	348858	51.8860	52
* 93 1,4-Dichlorobenzene-d4	152	12.237	12.236	(1.000)	281655	50.0000	

Data File: \\avogadro\organics\W5.1\130607.B\W503993.D
Date : 07-JUN-2013 15:14
Client ID: VBLKD5
Sample Info: 5ML,MB-72123,VBLKD5,72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
Data file : \\avogadro\organics\V5.I\130607.B\V503994.D
Lab Smp Id: LCS-72123 Client Smp ID: VLCSD5
Inj Date : 07-JUN-2013 15:39
Operator : WL SRC: LIMS Inst ID: V5.i
Smp Info : 5ML,LCS-72123,VLCSD5,72123
Misc Info :
Comment :
Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
Cal Date : 30-MAY-2013 14:32 Cal File: V503805.D
Als bottle: 6 Calibration Sample, Level: 3
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: FULL.sub
Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG						AMOUNTS	
		MASS	RT	EXP RT	REL RT	RESPONSE	CAL-AMT (ug/L)	ON-COL (ug/L)
1 Dichlorodifluoromethane	85		1.579	1.584 (0.282)		310501	50.0000	38
3 Chloromethane	50		1.718	1.735 (0.307)		762816	50.0000	50
4 Vinyl Chloride	62		1.835	1.839 (0.328)		605379	50.0000	48
5 Bromomethane	94		2.137	2.142 (0.382)		405931	50.0000	53
6 Chloroethane	64		2.230	2.234 (0.398)		351498	50.0000	51
7 Trichlorofluoromethane	101		2.439	2.444 (0.436)		484679	50.0000	50
8 Ethanol	46		2.613	2.618 (0.467)		235359	5000.00	4200(Q)
9 Ether	59		2.694	2.699 (0.481)		372267	50.0000	52(Q)
10 Acrolein	56		2.799	2.804 (0.500)		473153	250.000	260
11 1,1-Dichloroethene	96		2.892	2.908 (0.517)		330248	50.0000	46
12 1,1,2-Trichloro-1,2,2-Trifluo	101		2.915	2.920 (0.521)		323419	50.0000	51
13 Acetone	58		2.938	2.943 (0.525)		44322	50.0000	39
14 Iodomethane	142		3.031	3.036 (0.541)		574456	50.0000	47
15 Carbon Disulfide	76		3.089	3.106 (0.552)		1240740	50.0000	38
16 Acetonitrile	41		3.205	3.210 (0.573)		1611056	500.000	500
17 Allyl Chloride	39		3.205	3.210 (0.573)		667203	50.0000	54(Q)
18 Methyl Acetate	43		3.229	3.233 (0.577)		560159	50.0000	47
19 Methylene Chloride	84		3.322	3.326 (0.593)		432676	50.0000	49
20 tert-Butanol	59		3.426	3.431 (0.612)		73374	100.000	84
21 Acrylonitrile	53		3.531	3.536 (0.631)		170177	50.0000	51
22 trans-1,2-Dichloroethene	96		3.566	3.570 (0.637)		359567	50.0000	47
23 Methyl tert-butyl ether	73		3.566	3.570 (0.637)		982631	50.0000	49
24 1,1-Dichloroethane	63		3.937	3.942 (0.703)		812576	50.0000	52

Compounds	QUANT SIG						AMOUNTS	
	MASS	RT	EXP RT	REL RT	RESPONSE		CAL-AMT (ug/L)	ON-COL (ug/L)
=====	=====	=====	=====	=====	=====		=====	=====
25 Vinyl acetate	43	3.995	4.000 (0.714)		1599994		50.0000	52
26 Diisopropyl Ether	45	4.007	4.012 (0.716)		1998556		50.0000	55
27 2-Chloro-1,3-Butadiene	53	4.030	4.035 (0.720)		596067		50.0000	52
28 Ethyl tert-butyl ether	59	4.344	4.349 (0.776)		1346241		50.0000	52
29 cis-1,2-Dichloroethene	96	4.472	4.476 (0.799)		403093		50.0000	50
30 2,2-Dichloropropane	77	4.483	4.488 (0.801)		379598		50.0000	50
31 2-Butanone	72	4.483	4.488 (0.801)		45065		50.0000	48
32 Propionitrile	54	4.530	4.535 (0.809)		587980		500.000	470
33 Methacrylonitrile	41	4.681	4.686 (0.836)		669878		100.000	100
34 Bromochloromethane	128	4.692	4.697 (0.838)		195796		50.0000	51
35 Tetrahydrofuran	72	4.750	4.755 (0.849)		88970		100.000	91
36 Chloroform	83	4.774	4.778 (0.853)		669241		50.0000	53
\$ 37 Dibromofluoromethane	113	4.925	4.929 (0.880)		254800		50.0000	52
38 1,1,1-Trichloroethane	97	4.960	4.964 (0.886)		438570		50.0000	50
39 Cyclohexane	56	5.029	5.022 (0.898)		651777		50.0000	50
40 1,1-Dichloropropene	110	5.122	5.127 (0.915)		159253		50.0000	47
41 Carbon Tetrachloride	117	5.134	5.127 (0.917)		348219		50.0000	50
42 Isobutyl Alcohol	43	5.215	5.220 (0.932)		401801		1000.00	900
\$ 43 1,2-Dichloroethane-d4	102	5.262	5.255 (0.940)		53982		50.0000	50
44 Benzene	78	5.320	5.324 (0.950)		1427322		50.0000	52
45 1,2-Dichloroethane	62	5.331	5.336 (0.952)		552322		50.0000	56
46 tert-Amyl methyl ether	73	5.436	5.441 (0.971)		993205		50.0000	50
* 47 Fluorobenzene	96	5.598	5.603 (1.000)		788569		50.0000	
48 Trichloroethene	130	5.982	5.987 (1.068)		317305		50.0000	46
49 Methylcyclohexane	83	6.191	6.196 (1.106)		428831		50.0000	52
50 1,2-Dichloropropane	63	6.214	6.219 (1.110)		465016		50.0000	54
51 Dibromomethane	93	6.330	6.335 (1.131)		273149		50.0000	52
52 Methyl Methacrylate	69	6.330	6.335 (1.131)		269529		50.0000	45
53 1,4-Dioxane	88	6.354	6.358 (1.135)		55431		1000.00	760
54 Bromodichloromethane	83	6.505	6.509 (1.162)		494351		50.0000	54
M 55 1,2-Dichloroethene (Total)	96				762660		100.000	97
56 2-Chloroethyl vinyl ether	63	7.004	7.009 (1.251)		4043		50.0000	29 (T)
57 cis-1,3-Dichloropropene	75	7.004	7.009 (1.251)		627488		50.0000	53
58 4-Methyl-2-pentanone	43	7.190	7.183 (1.284)		557010		50.0000	49
\$ 59 Toluene-d8	98	7.318	7.323 (0.807)		791763		50.0000	46
60 Toluene	91	7.399	7.404 (1.322)		1299964		50.0000	52
61 trans-1,3-Dichloropropene	75	7.643	7.648 (1.365)		539643		50.0000	53
62 Ethyl Methacrylate	69	7.771	7.776 (1.388)		316906		50.0000	46
63 1,1,2-Trichloroethane	97	7.864	7.868 (1.405)		309298		50.0000	48
64 Tetrachloroethene	164	8.050	8.054 (0.887)		233196		50.0000	37
65 1,3-Dichloropropane	76	8.073	8.066 (0.890)		567696		50.0000	44
66 2-Hexanone	43	8.177	8.182 (0.901)		368913		50.0000	41
67 Dibromochloromethane	129	8.340	8.345 (0.919)		331020		50.0000	44
68 1,2-Dibromoethane	107	8.479	8.484 (0.935)		359640		50.0000	43
* 69 Chlorobenzene-d5	117	9.072	9.077 (1.000)		658540		50.0000	
70 1-Chlorohexane	91	9.095	9.100 (1.003)		449977		50.0000	42 (Q)
71 Chlorobenzene	112	9.118	9.111 (1.005)		844608		50.0000	44
72 1,1,1,2-Tetrachloroethane	131	9.223	9.216 (1.017)		289108		50.0000	42
73 Ethylbenzene	106	9.269	9.262 (1.022)		426414		50.0000	42
74 m,p-Xylene	106	9.420	9.413 (1.038)		1063411		100.000	86
75 o-Xylene	106	9.943	9.936 (1.096)		533577		50.0000	44
76 Styrene	104	9.955	9.959 (1.097)		828199		50.0000	46
77 Bromoform	173	10.187	10.180 (1.123)		187690		50.0000	41
78 Isopropylbenzene	105	10.443	10.436 (1.151)		1276254		50.0000	43

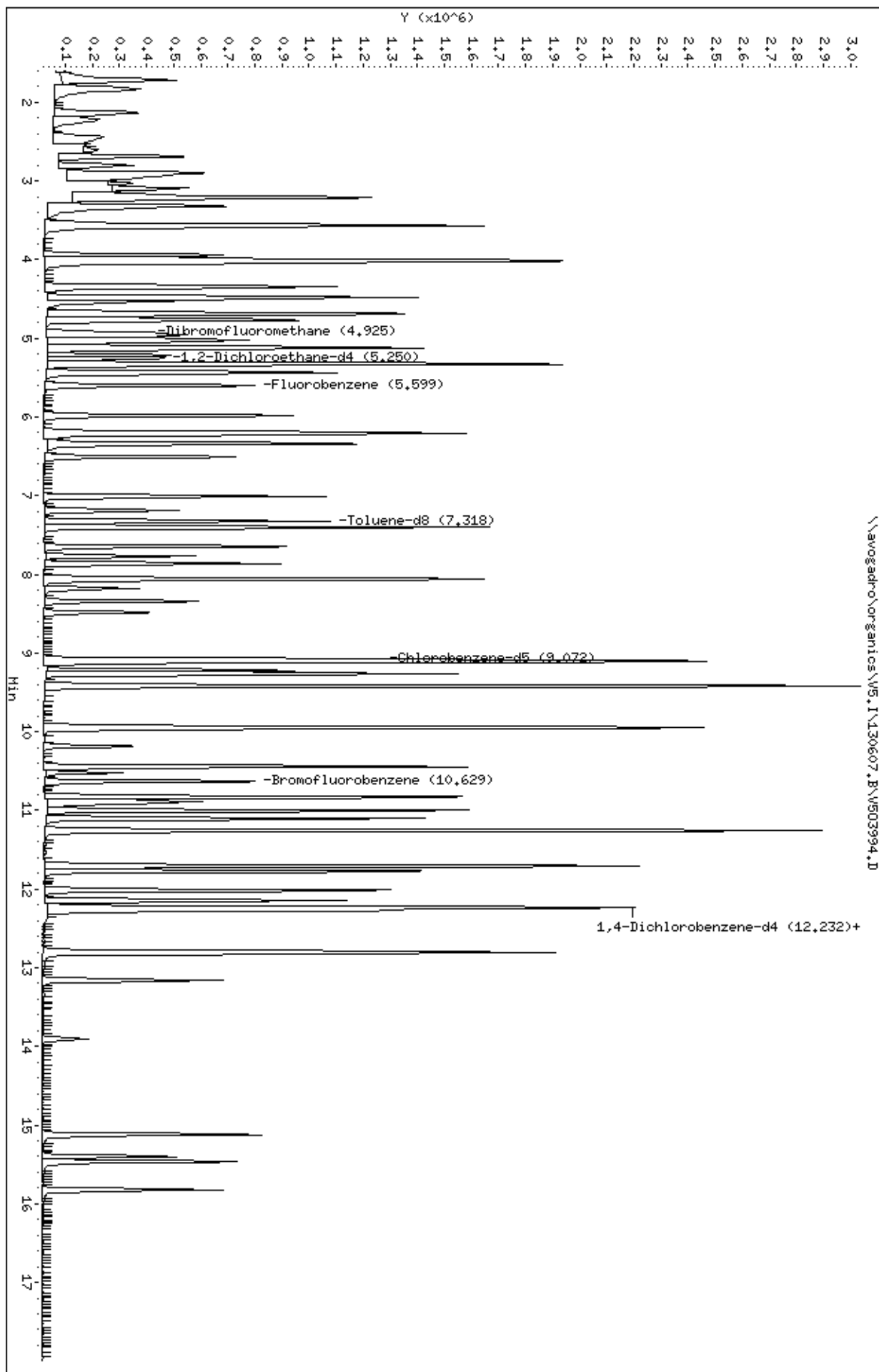
Compounds	QUANT SIG						AMOUNTS	
	MASS	RT	EXP RT	REL RT	RESPONSE		CAL-AMT (ug/L)	ON-COL (ug/L)
=====	=====	=====	=====	=====	=====		=====	=====
79 trans-1,4-Dichloro-2-butene	75	10.524	10.517	(1.160)	80667		50.0000	36
\$ 80 Bromofluorobenzene	95	10.628	10.633	(1.172)	367036		50.0000	54
81 Bromobenzene	156	10.826	10.831	(0.884)	332276		50.0000	40
82 1,1,2,2-Tetrachloroethane	83	10.838	10.842	(0.885)	459374		50.0000	40
83 1,2,3-Trichloropropane	75	10.896	10.889	(0.890)	409679		50.0000	40
84 n-Propylbenzene	120	11.000	11.005	(0.898)	311042		50.0000	41
85 2-Chlorotoluene	126	11.105	11.110	(0.907)	298208		50.0000	40
86 4-Chlorotoluene	126	11.256	11.261	(0.919)	313970		50.0000	40
87 1,3,5-Trimethylbenzene	105	11.256	11.261	(0.919)	1019062		50.0000	42
88 tert-Butylbenzene	119	11.709	11.702	(0.956)	950579		50.0000	41
89 1,2,4-Trimethylbenzene	105	11.778	11.772	(0.962)	1033449		50.0000	42
90 sec-Butylbenzene	105	12.011	12.016	(0.981)	1227904		50.0000	40
91 1,3-Dichlorobenzene	146	12.150	12.143	(0.992)	560875		50.0000	40
92 4-Isopropyltoluene	119	12.231	12.225	(0.999)	925062		50.0000	40
* 93 1,4-Dichlorobenzene-d4	152	12.243	12.236	(1.000)	306088		50.0000	
94 1,4-Dichlorobenzene	146	12.278	12.271	(1.003)	590355		50.0000	41
95 1,2-Dichlorobenzene	146	12.789	12.794	(1.045)	546584		50.0000	41
M 96 Xylene (Total)	106				1596988		150.000	130
97 n-Butylbenzene	91	12.812	12.817	(1.046)	937951		50.0000	42
98 Hexachloroethane	117	13.161	13.154	(1.075)	189385		50.0000	50
99 1,2-Dibromo-3-chloropropane	75	13.904	13.909	(1.136)	54283		50.0000	33
100 1,3,5-Trichlorobenzene	182	15.124	15.117	(2.701)	312851		50.0000	48
101 1,2,4-Trichlorobenzene	180	15.124	15.117	(1.235)	326317		50.0000	38
102 Hexachlorobutadiene	225	15.403	15.408	(1.258)	117682		50.0000	34
103 Naphthalene	128	15.461	15.466	(1.263)	677366		50.0000	34
104 1,2,3-Trichlorobenzene	180	15.821	15.826	(1.292)	258886		50.0000	35

QC Flag Legend

T - Target compound detected outside RT window.
 Q - Qualifier signal failed the ratio test.

Data File: \\avogadro\organics\W5,I\130607.B\W503994.D
Date : 07-JUN-2013 15:39
Client ID: WLCSD5
Sample Info: 5HL,LCS-72123,WLCSD5,72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V503995.D
Report Date: 11-Jun-2013 09:23

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
Data file : \\avogadro\organics\V5.I\130607.B\V503995.D
Lab Smp Id: M0903-06AMS Client Smp ID: MW17-060413MS
Inj Date : 07-JUN-2013 16:05
Operator : WL SRC: LIMS Inst ID: V5.i
Smp Info : 5ML,M0903-06AMS,,72123
Misc Info :
Comment :
Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
Cal Date : 30-MAY-2013 14:32 Cal File: V503805.D
Als bottle: 7 Calibration Sample, Level: 3
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: BEX.sub
Target Version: 4.14

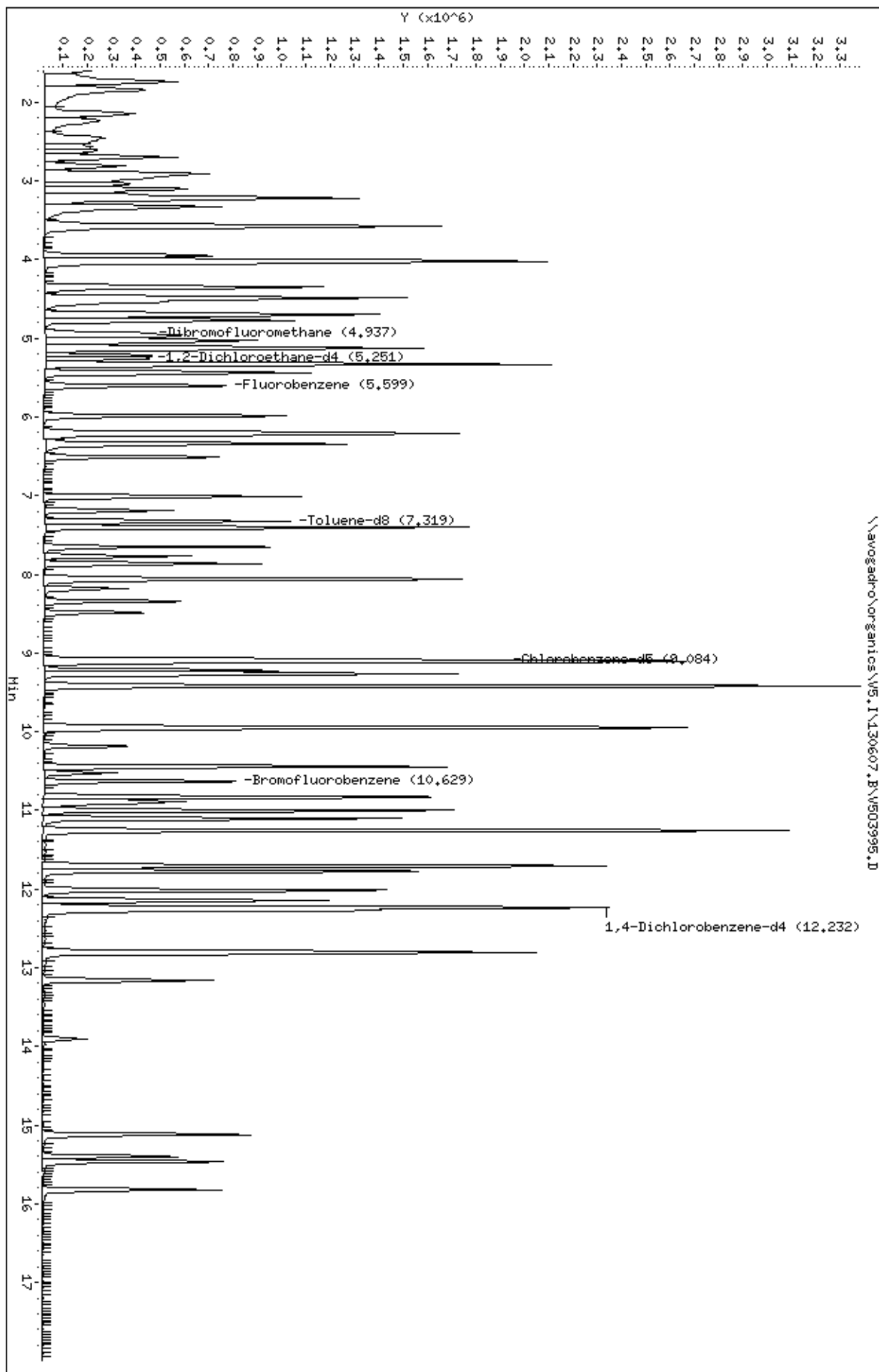
Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG	RT	EXP RT	REL RT	RESPONSE	AMOUNTS	
						CAL-AMT (ug/L)	ON-COL (ug/L)
\$ 37 Dibromofluoromethane	113	4.937	4.929	(0.880)	262271	50.0000	53
\$ 43 1,2-Dichloroethane-d4	102	5.262	5.255	(0.938)	51458	50.0000	47
44 Benzene	78	5.332	5.324	(0.950)	1521065	50.0000	56
* 47 Fluorobenzene	96	5.610	5.603	(1.000)	793004	50.0000	
\$ 59 Toluene-d8	98	7.318	7.323	(0.806)	800340	50.0000	46
* 69 Chlorobenzene-d5	117	9.084	9.077	(1.000)	665941	50.0000	
73 Ethylbenzene	106	9.270	9.262	(1.020)	471067	50.0000	46
74 m,p-Xylene	106	9.421	9.413	(1.037)	1148662	100.000	92
75 o-Xylene	106	9.943	9.936	(1.095)	586459	50.0000	48
\$ 80 Bromofluorobenzene	95	10.629	10.633	(1.170)	365270	50.0000	53
* 93 1,4-Dichlorobenzene-d4	152	12.244	12.236	(1.000)	312009	50.0000	
M 96 Xylene (Total)	106				1735121	150.000	140

Data File: \\avogadro\organics\W5.I\130607.B\W503995.D
Date : 07-JUN-2013 16:05
Client ID: MW17-060413MS
Sample Info: SWL.H0903-06AMS,,72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V503996.D
 Report Date: 11-Jun-2013 09:23

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130607.B\V503996.D
 Lab Smp Id: M0903-06AMSD Client Smp ID: MW17-060413MSD
 Inj Date : 07-JUN-2013 16:30
 Operator : WL SRC: LIMS Inst ID: V5.i
 Smp Info : 5ML,M0903-06AMSD,,72123
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
 Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V503805.D
 Als bottle: 8 Calibration Sample, Level: 3
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: BEX.sub
 Target Version: 4.14

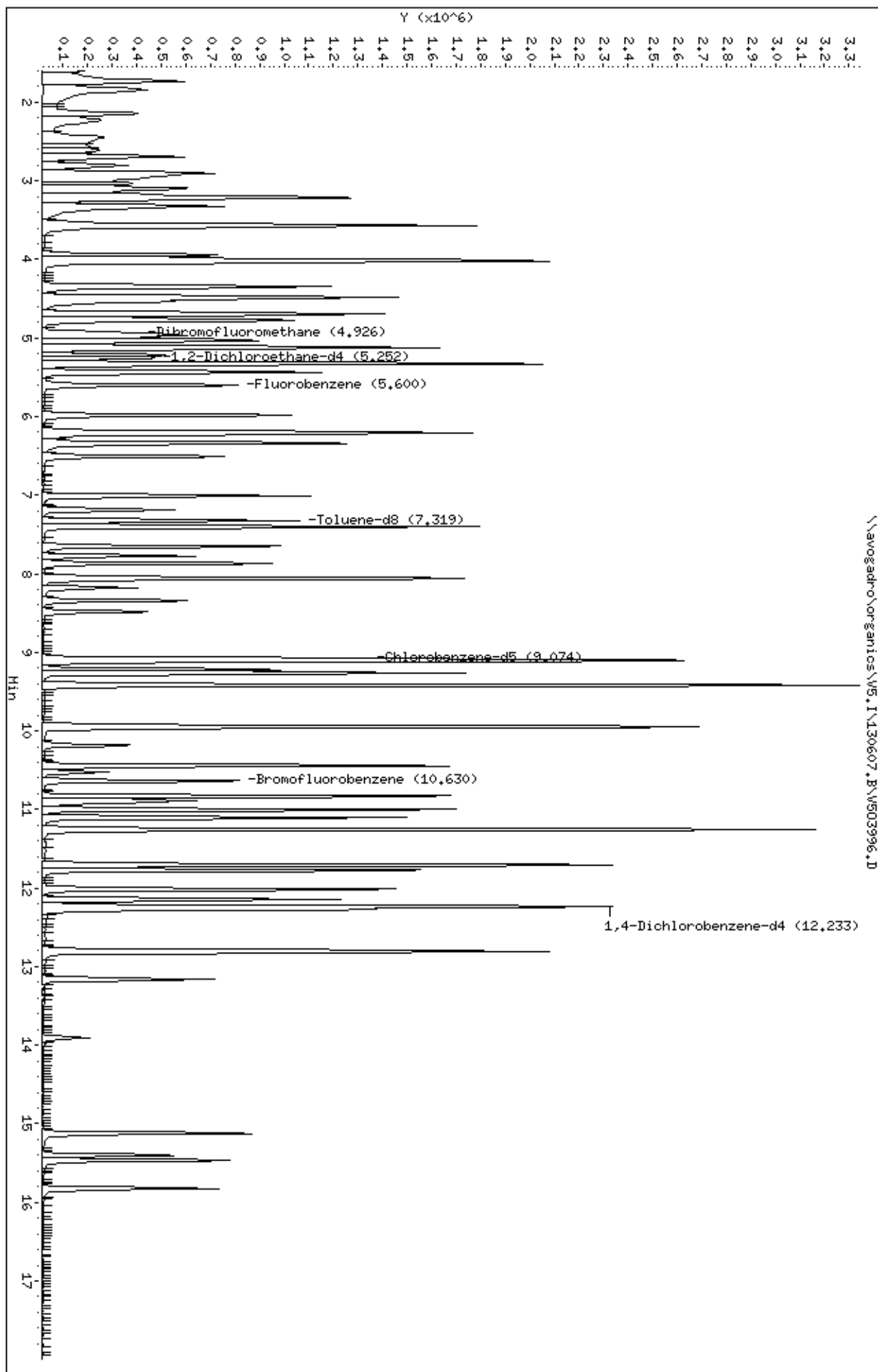
Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG	RT	EXP RT	REL RT	RESPONSE	AMOUNTS	
						CAL-AMT (ug/L)	ON-COL (ug/L)
\$ 37 Dibromofluoromethane	113	4.926	4.929	(0.880)	260694	50.0000	53
\$ 43 1,2-Dichloroethane-d4	102	5.263	5.255	(0.940)	54131	50.0000	50
44 Benzene	78	5.332	5.324	(0.952)	1523587	50.0000	56
* 47 Fluorobenzene	96	5.600	5.603	(1.000)	790614	50.0000	
\$ 59 Toluene-d8	98	7.319	7.323	(0.807)	796718	50.0000	46
* 69 Chlorobenzene-d5	117	9.073	9.077	(1.000)	667620	50.0000	
73 Ethylbenzene	106	9.270	9.262	(1.022)	472644	50.0000	46
74 m,p-Xylene	106	9.422	9.413	(1.038)	1158875	100.000	93
75 o-Xylene	106	9.944	9.936	(1.096)	574591	50.0000	46
\$ 80 Bromofluorobenzene	95	10.630	10.633	(1.172)	366940	50.0000	53
* 93 1,4-Dichlorobenzene-d4	152	12.244	12.236	(1.000)	304729	50.0000	
M 96 Xylene (Total)	106				1733466	150.000	140

Data File: \\avogadro\organics\W5.I\130607.B\W503996.D
Date : 07-JUN-2013 16:30
Client ID: MM17-060413MSD
Sample Info: SHL, H0903-06MSD, 72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V5O4002.D
 Report Date: 11-Jun-2013 09:23

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130607.B\V5O4002.D
 Lab Smp Id: M0903-01A Client Smp ID: MW112-060413
 Inj Date : 07-JUN-2013 19:04
 Operator : WL SRC: LIMS Inst ID: V5.i
 Smp Info : 5ML,M0903-01A,,72123
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
 Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V5O3805.D
 Als bottle: 14
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: BEX.sub
 Target Version: 4.14

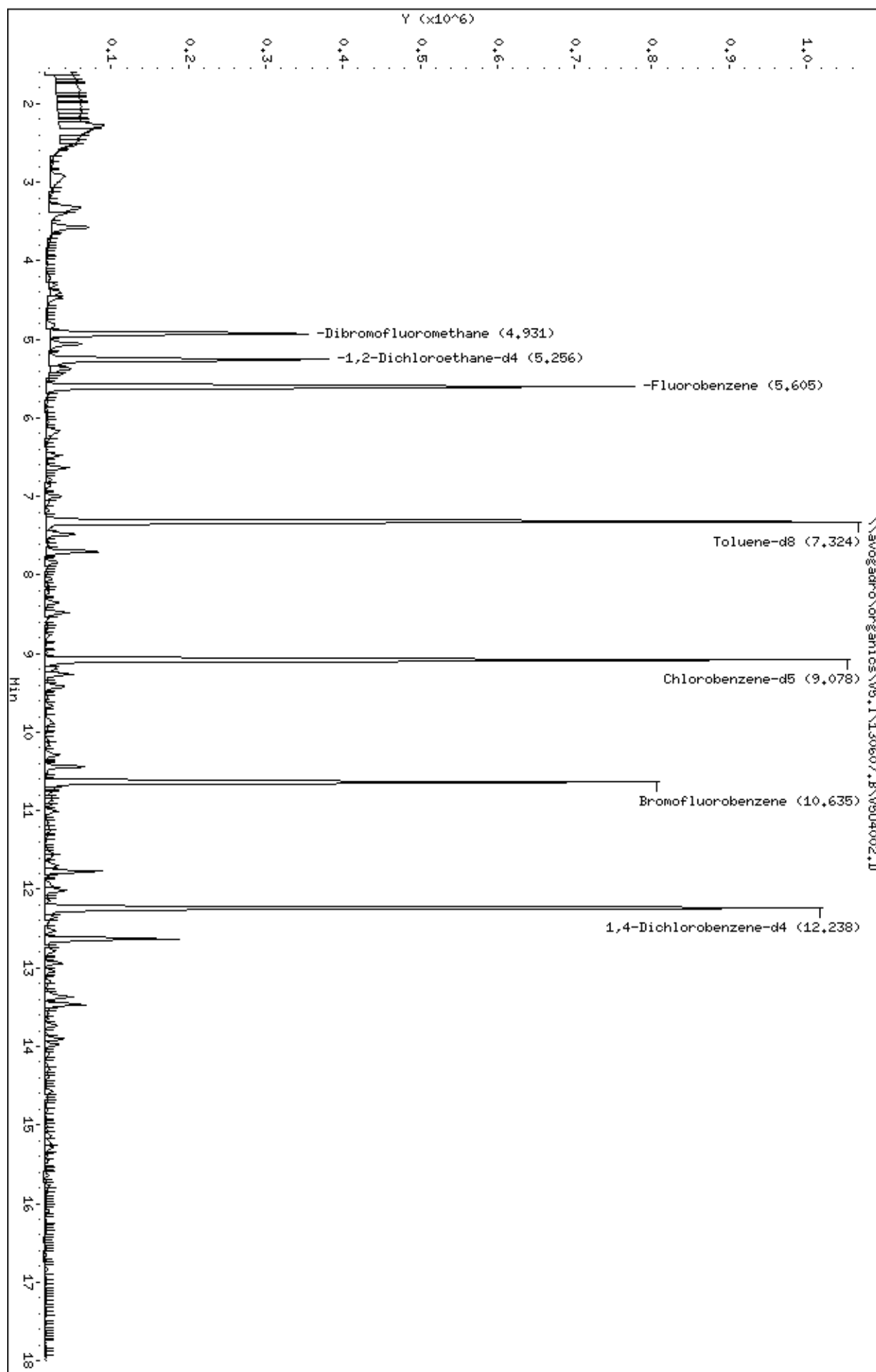
Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS	
						ON-COLUMN (ug/L)	FINAL (ug/L)
\$ 37 Dibromofluoromethane	113	4.930	4.929	(0.880)	264287	53.4753	53
\$ 43 1,2-Dichloroethane-d4	102	5.256	5.255	(0.938)	55884	50.7729	51
* 47 Fluorobenzene	96	5.604	5.603	(1.000)	796720	50.0000	
\$ 59 Toluene-d8	98	7.324	7.323	(0.807)	791782	47.0192	47
* 69 Chlorobenzene-d5	117	9.078	9.077	(1.000)	652480	50.0000	
\$ 80 Bromofluorobenzene	95	10.634	10.633	(1.171)	357595	52.9337	53
* 93 1,4-Dichlorobenzene-d4	152	12.237	12.236	(1.000)	284647	50.0000	

Data File: \\avogadro\organics\W5.1\130607.B\W504002.D
Date : 07-JUN-2013 19:04
Client ID: MM12-060413
Sample Info: SHL.M0903-01A,72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V5O4003.D
 Report Date: 11-Jun-2013 09:23

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130607.B\V5O4003.D
 Lab Smp Id: M0903-02A Client Smp ID: MW106-060413
 Inj Date : 07-JUN-2013 19:30
 Operator : WL SRC: LIMS Inst ID: V5.i
 Smp Info : 5ML,M0903-02A,,72123
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
 Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V5O3805.D
 Als bottle: 15
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: BEX.sub
 Target Version: 4.14

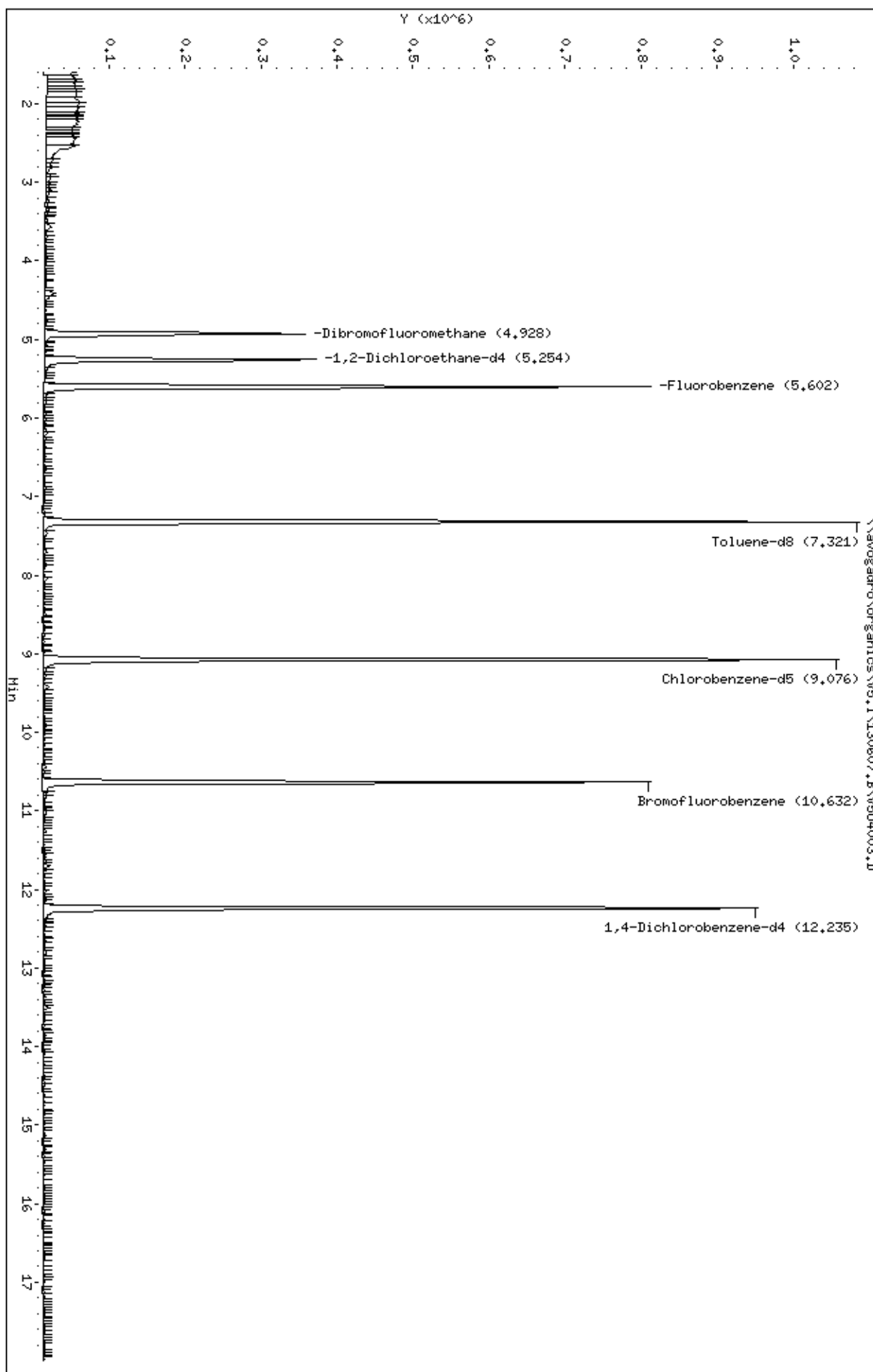
Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS	
						ON-COLUMN (ug/L)	FINAL (ug/L)
\$ 37 Dibromofluoromethane	113	4.928	4.929	(0.880)	255948	51.4857	51
\$ 43 1,2-Dichloroethane-d4	102	5.253	5.255	(0.938)	54986	49.6654	50
* 47 Fluorobenzene	96	5.602	5.603	(1.000)	801398	50.0000	
\$ 59 Toluene-d8	98	7.321	7.323	(0.807)	795684	46.8504	47
* 69 Chlorobenzene-d5	117	9.075	9.077	(1.000)	658057	50.0000	
\$ 80 Bromofluorobenzene	95	10.632	10.633	(1.172)	356390	52.3082	52
* 93 1,4-Dichlorobenzene-d4	152	12.235	12.236	(1.000)	280142	50.0000	

Data File: \\avogadro\organics\W5, I\130607.B\W504003.D
Date : 07-JUN-2013 19:30
Client ID: MM106-060413
Sample Info: SHL, M0903-02H, 72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V504004.D
 Report Date: 11-Jun-2013 09:23

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130607.B\V504004.D
 Lab Smp Id: M0903-03A Client Smp ID: DUP-060413
 Inj Date : 07-JUN-2013 19:56
 Operator : WL SRC: LIMS Inst ID: V5.i
 Smp Info : 5ML,M0903-03A,,72123
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
 Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V503805.D
 Als bottle: 16
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: BEX.sub
 Target Version: 4.14

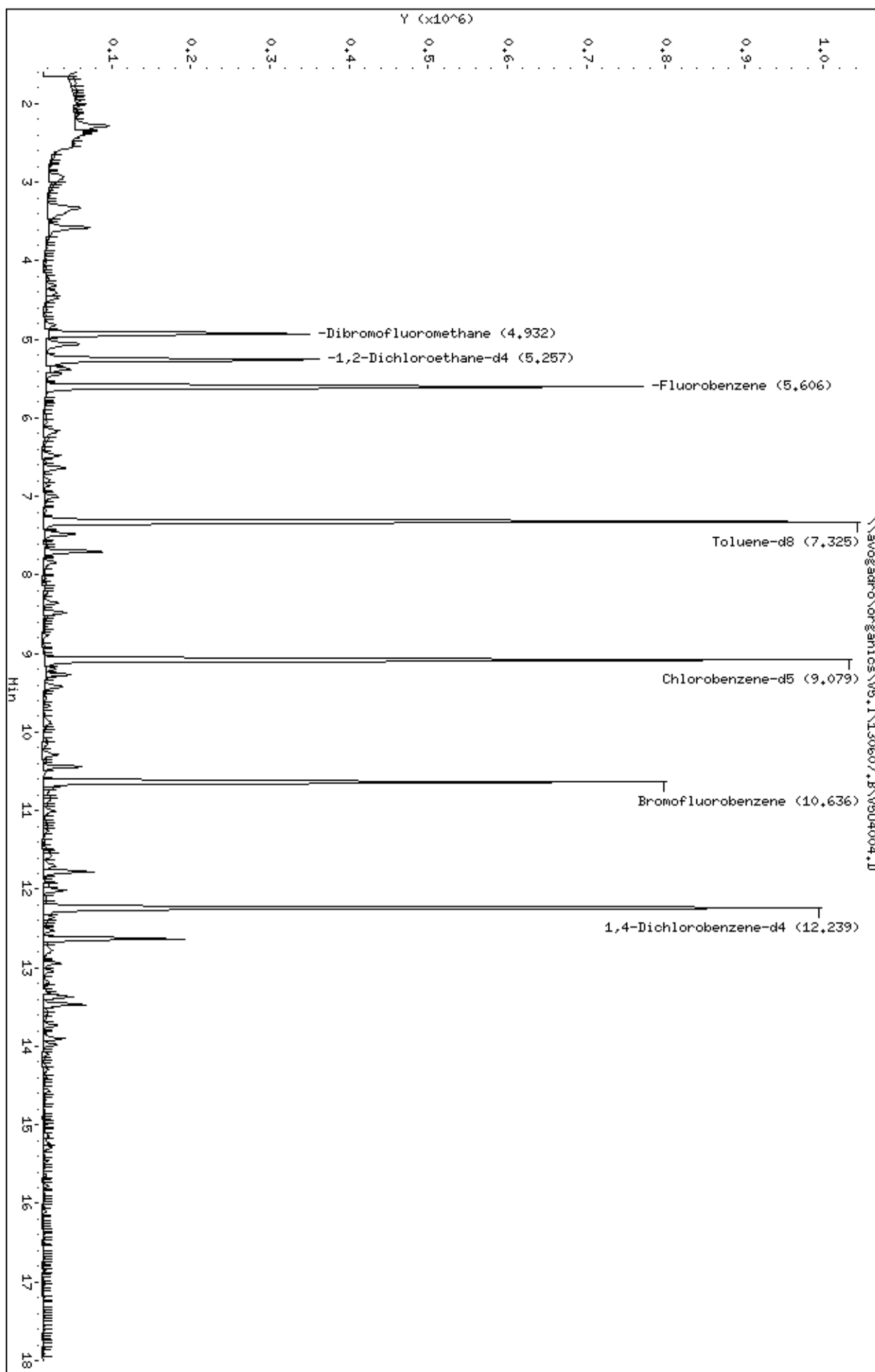
Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS	
						ON-COLUMN (ug/L)	FINAL (ug/L)
\$ 37 Dibromofluoromethane	113	4.931	4.929	(0.880)	253232	52.5839	52
\$ 43 1,2-Dichloroethane-d4	102	5.268	5.255	(0.940)	51725	48.2283	48
* 47 Fluorobenzene	96	5.605	5.603	(1.000)	776334	50.0000	
\$ 59 Toluene-d8	98	7.324	7.323	(0.807)	774103	47.0149	47
* 69 Chlorobenzene-d5	117	9.078	9.077	(1.000)	637969	50.0000	
\$ 80 Bromofluorobenzene	95	10.635	10.633	(1.171)	346273	52.4236	52
* 93 1,4-Dichlorobenzene-d4	152	12.238	12.236	(1.000)	274792	50.0000	

Data File: \\avogadro\organics\W5.1\130607.B\W504004.D
Date : 07-JUN-2013 19:56
Client ID: DUP-060413
Sample Info: SHL.M0903-03H,72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V504005.D
 Report Date: 11-Jun-2013 09:23

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130607.B\V504005.D
 Lab Smp Id: M0903-04A Client Smp ID: TB-060413
 Inj Date : 07-JUN-2013 20:22
 Operator : WL SRC: LIMS Inst ID: V5.i
 Smp Info : 5ML,M0903-04A,,72123
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
 Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V503805.D
 Als bottle: 17
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: BEX.sub
 Target Version: 4.14

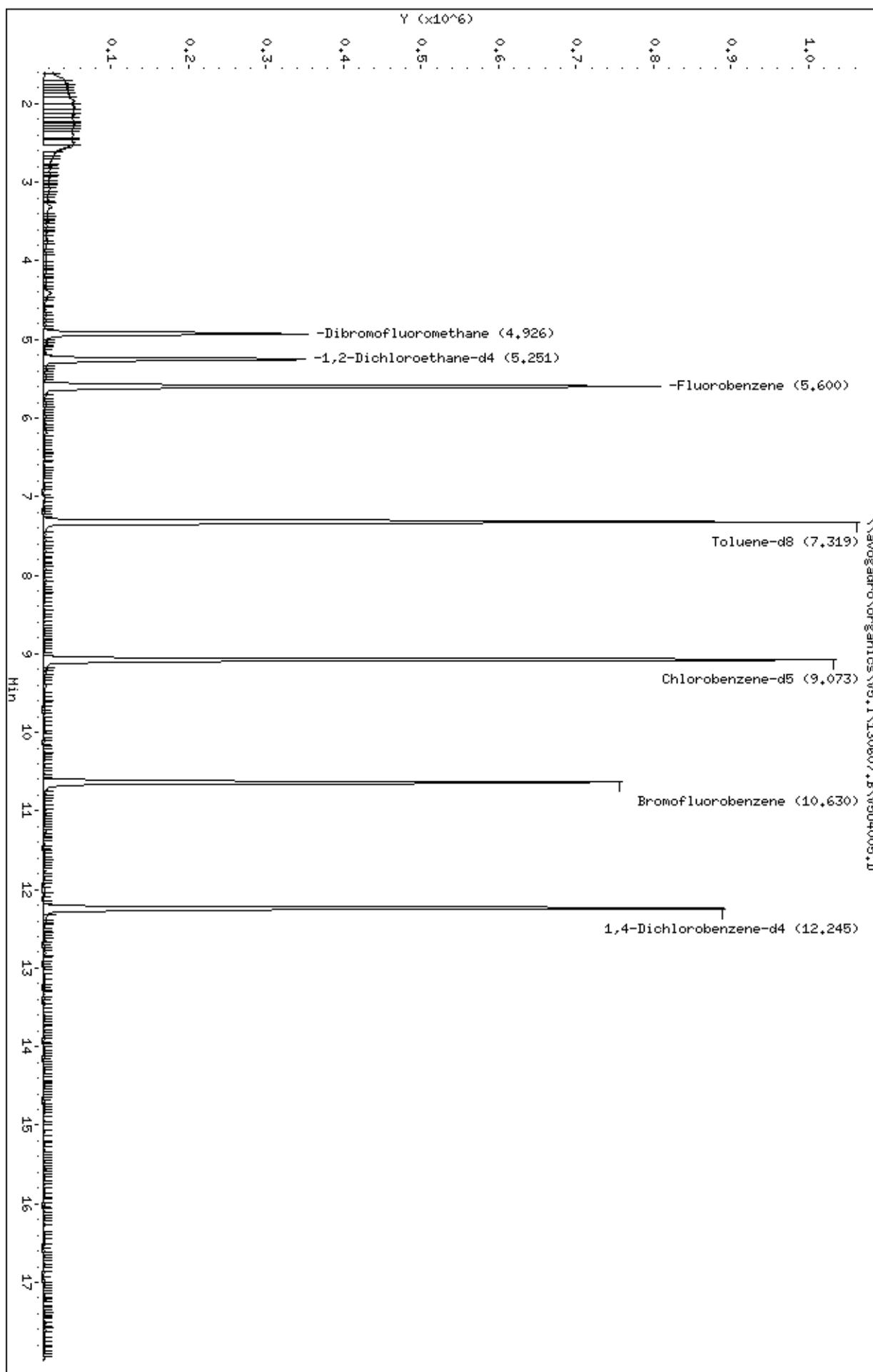
Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS	
						ON-COLUMN (ug/L)	FINAL (ug/L)
\$ 37 Dibromofluoromethane	113	4.926	4.929	(0.880)	251271	52.0223	52
\$ 43 1,2-Dichloroethane-d4	102	5.251	5.255	(0.938)	51422	47.8038	48
* 47 Fluorobenzene	96	5.599	5.603	(1.000)	778639	50.0000	
\$ 59 Toluene-d8	98	7.319	7.323	(0.807)	764993	46.1438	46
* 69 Chlorobenzene-d5	117	9.073	9.077	(1.000)	642363	50.0000	
\$ 80 Bromofluorobenzene	95	10.629	10.633	(1.172)	340570	51.2075	51
* 93 1,4-Dichlorobenzene-d4	152	12.244	12.236	(1.000)	274583	50.0000	

Data File: \\avogadro\organics\W5,I\130607.B\W504005.D
Date : 07-JUN-2013 20:22
Client ID: TB-060413
Sample Info: 5HL,M0903-044,72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V5O4006.D
 Report Date: 11-Jun-2013 09:23

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130607.B\V5O4006.D
 Lab Smp Id: M0903-05A Client Smp ID: MW19-060413
 Inj Date : 07-JUN-2013 20:48
 Operator : WL SRC: LIMS Inst ID: V5.i
 Smp Info : 5ML,M0903-05A,,72123
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
 Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V5O3805.D
 Als bottle: 18
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: BEX.sub
 Target Version: 4.14

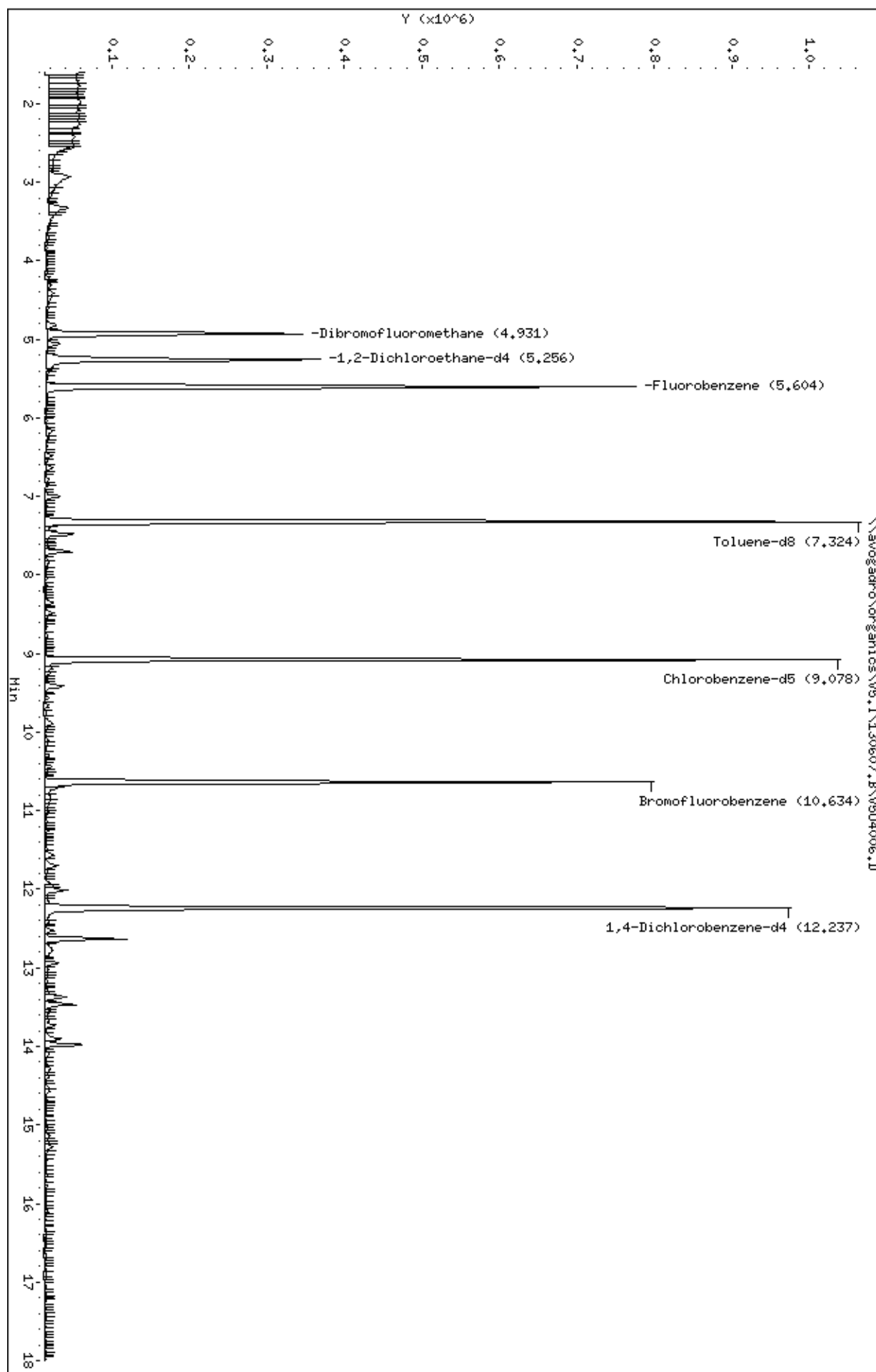
Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS	
						ON-COLUMN (ug/L)	FINAL (ug/L)
\$ 37 Dibromofluoromethane	113	4.930	4.929	(0.880)	254418	52.6519	53
\$ 43 1,2-Dichloroethane-d4	102	5.255	5.255	(0.938)	52540	48.8229	49
* 47 Fluorobenzene	96	5.604	5.603	(1.000)	778963	50.0000	
\$ 59 Toluene-d8	98	7.323	7.323	(0.807)	766790	47.1352	47
* 69 Chlorobenzene-d5	117	9.077	9.077	(1.000)	630329	50.0000	
\$ 80 Bromofluorobenzene	95	10.634	10.633	(1.171)	342377	52.4620	52
* 93 1,4-Dichlorobenzene-d4	152	12.237	12.236	(1.000)	278498	50.0000	

Data File: \\avogadro\organics\W5.1\130607.B\W504006.D
Date : 07-JUN-2013 20:48
Client ID: MM19-060413
Sample Info: 5HL, M0903-05H, 72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V504007.D
 Report Date: 11-Jun-2013 09:23

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130607.B\V504007.D
 Lab Smp Id: M0903-06A Client Smp ID: MW17-060413
 Inj Date : 07-JUN-2013 21:13
 Operator : WL SRC: LIMS Inst ID: V5.i
 Smp Info : 5ML,M0903-06A,,72123
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
 Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V503805.D
 Als bottle: 19
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: BEX.sub
 Target Version: 4.14

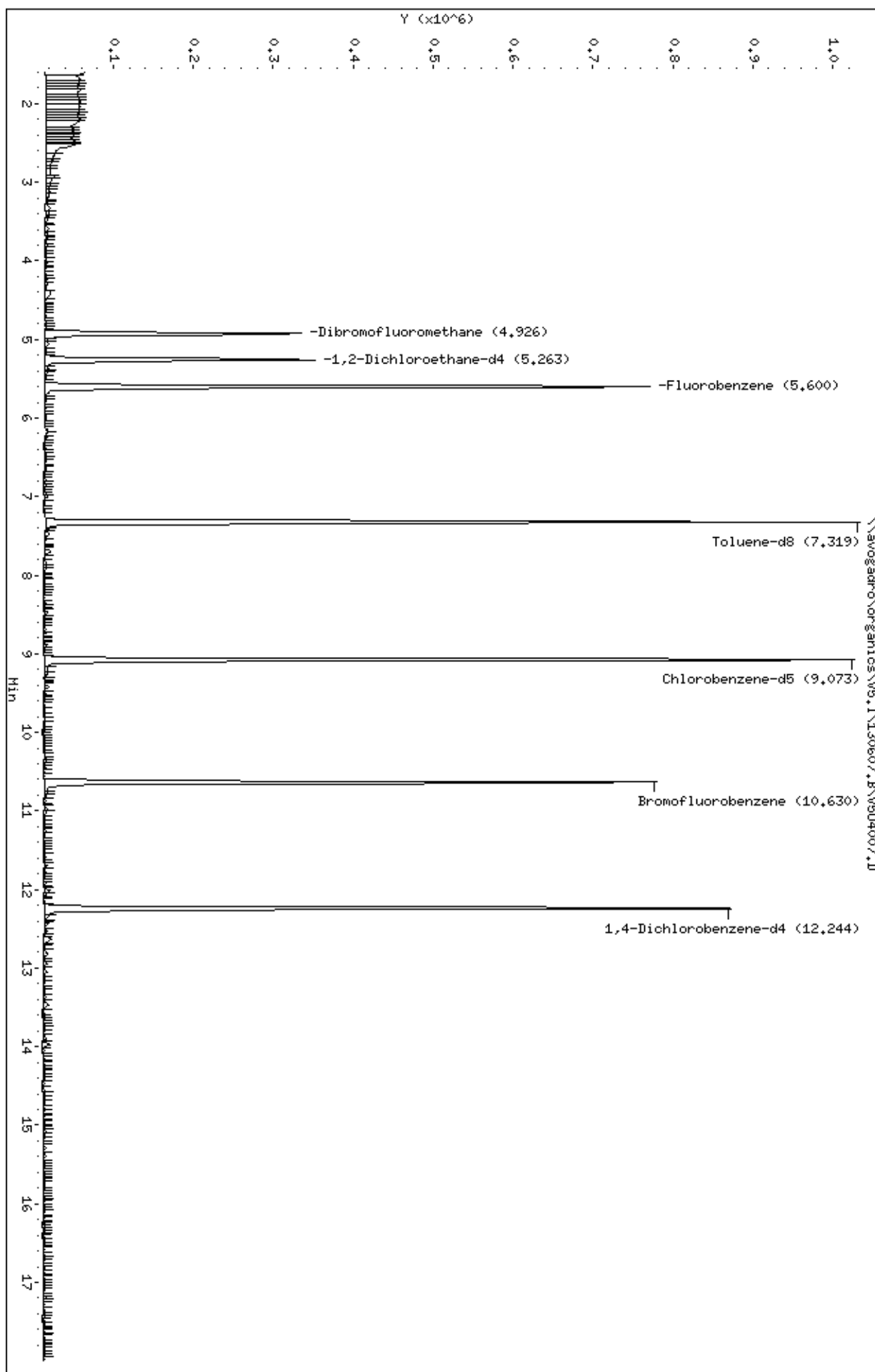
Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS	
						ON-COLUMN (ug/L)	FINAL (ug/L)
\$ 37 Dibromofluoromethane	113	4.925	4.929	(0.880)	246721	52.1832	52
\$ 43 1,2-Dichloroethane-d4	102	5.262	5.255	(0.940)	52435	49.7981	50
* 47 Fluorobenzene	96	5.599	5.603	(1.000)	762182	50.0000	
\$ 59 Toluene-d8	98	7.318	7.323	(0.807)	750595	46.6596	47
* 69 Chlorobenzene-d5	117	9.073	9.077	(1.000)	623305	50.0000	
\$ 80 Bromofluorobenzene	95	10.629	10.633	(1.172)	338019	52.3779	52
* 93 1,4-Dichlorobenzene-d4	152	12.244	12.236	(1.000)	264311	50.0000	

Data File: \\avogadro\organics\W5.1\130607.B\W504007.D
Date : 07-JUN-2013 21:13
Client ID: MW17-060413
Sample Info: SWL, M0903-06H, 72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V504008.D
 Report Date: 11-Jun-2013 09:23

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130607.B\V504008.D
 Lab Smp Id: M0903-07A Client Smp ID: MW15-060413
 Inj Date : 07-JUN-2013 21:39
 Operator : WL SRC: LIMS Inst ID: V5.i
 Smp Info : 5ML,M0903-07A,,72123
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
 Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V503805.D
 Als bottle: 20
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: BEX.sub
 Target Version: 4.14

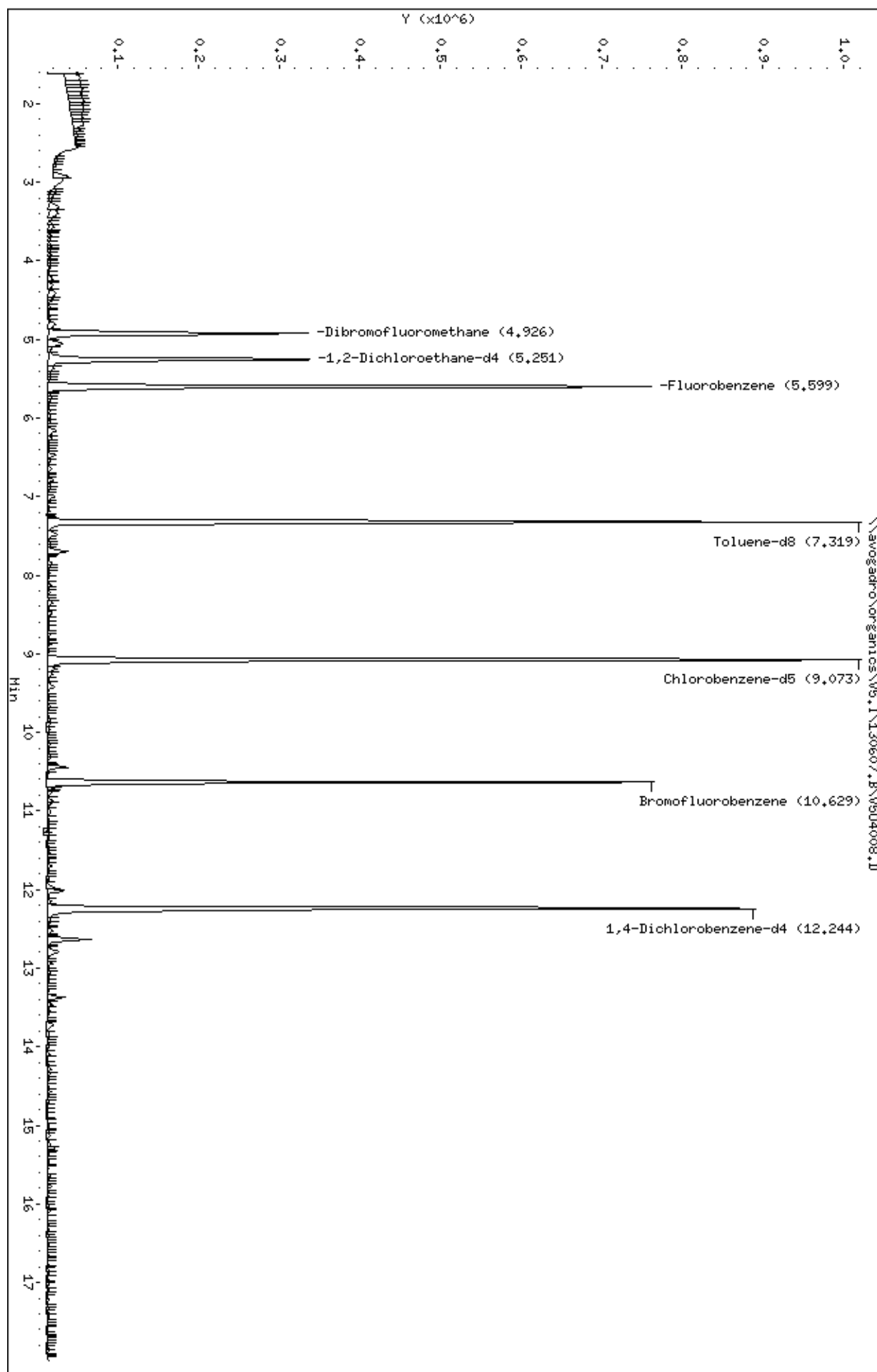
Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS	
						ON-COLUMN (ug/L)	FINAL (ug/L)
\$ 37 Dibromofluoromethane	113	4.925	4.929	(0.880)	243192	52.1566	52
\$ 43 1,2-Dichloroethane-d4	102	5.262	5.255	(0.940)	50823	48.9426	49
* 47 Fluorobenzene	96	5.599	5.603	(1.000)	751663	50.0000	
\$ 59 Toluene-d8	98	7.318	7.323	(0.807)	736902	45.6317	46
* 69 Chlorobenzene-d5	117	9.072	9.077	(1.000)	625719	50.0000	
\$ 80 Bromofluorobenzene	95	10.629	10.633	(1.172)	337942	52.1640	52
* 93 1,4-Dichlorobenzene-d4	152	12.244	12.236	(1.000)	264861	50.0000	

Data File: \\avogadro\organics\W5,I\130607.B\W504008.D
Date : 07-JUN-2013 21:39
Client ID: MM15-060413
Sample Info: SHL,M0903-07H,72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V504009.D
 Report Date: 11-Jun-2013 09:23

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130607.B\V504009.D
 Lab Smp Id: M0903-08A Client Smp ID: MW11-060413
 Inj Date : 07-JUN-2013 22:05
 Operator : WL SRC: LIMS Inst ID: V5.i
 Smp Info : 5ML,M0903-08A,,72123
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
 Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V503805.D
 Als bottle: 21
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: BEX.sub
 Target Version: 4.14

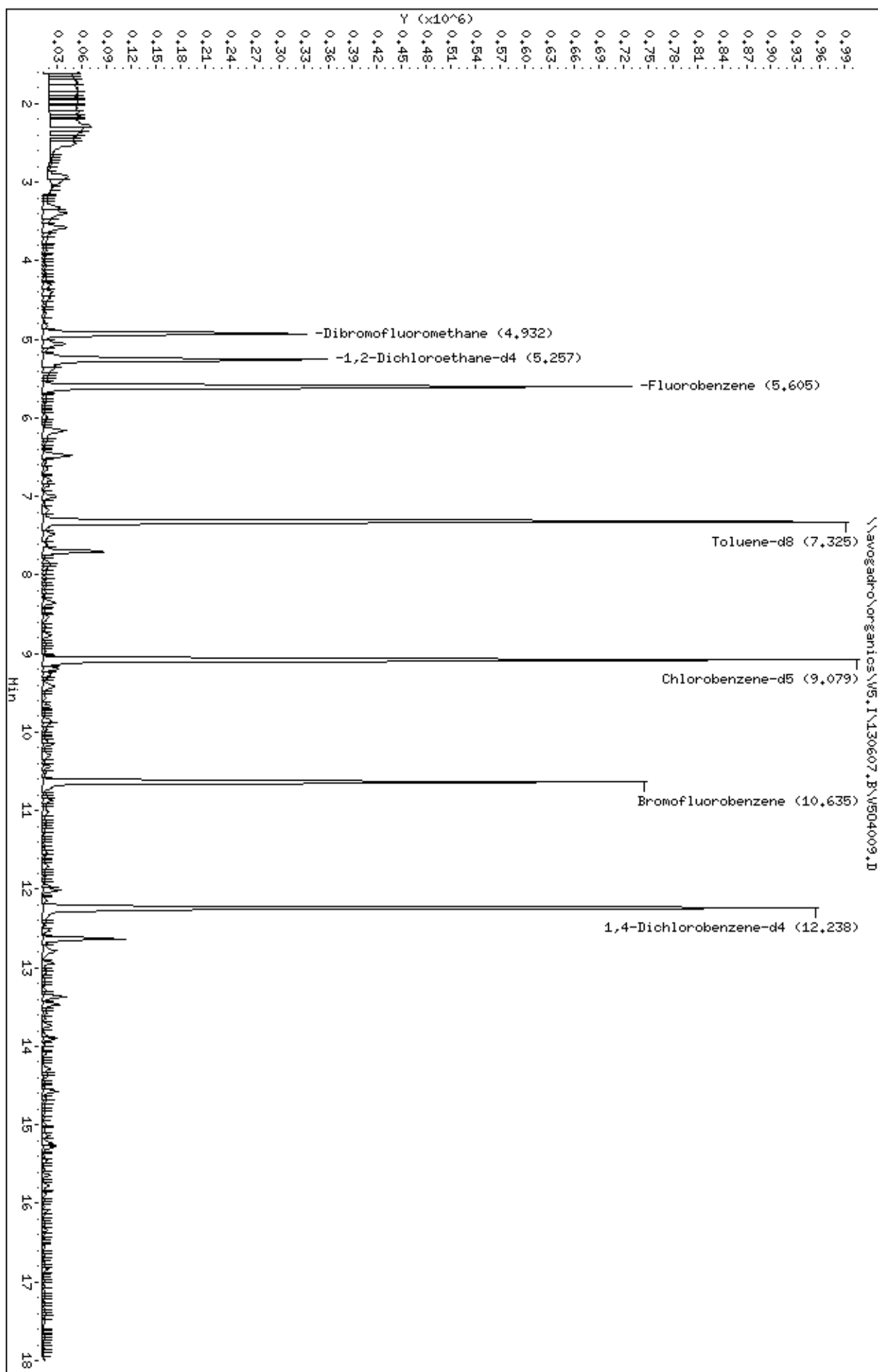
Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS	
						ON-COLUMN (ug/L)	FINAL (ug/L)
\$ 37 Dibromofluoromethane	113	4.931	4.929	(0.880)	244252	53.3612	53
\$ 43 1,2-Dichloroethane-d4	102	5.256	5.255	(0.938)	48462	47.5396	48
* 47 Fluorobenzene	96	5.605	5.603	(1.000)	737897	50.0000	
\$ 59 Toluene-d8	98	7.324	7.323	(0.807)	736926	46.3178	46
* 69 Chlorobenzene-d5	117	9.078	9.077	(1.000)	616471	50.0000	
\$ 80 Bromofluorobenzene	95	10.635	10.633	(1.171)	329155	51.5698	52
* 93 1,4-Dichlorobenzene-d4	152	12.238	12.236	(1.000)	261716	50.0000	

Data File: \\avogadro\organics\W5.1\130607.B\W504009.D
Date : 07-JUN-2013 22:05
Client ID: MW11-060413
Sample Info: SWL, M0903-08H, 72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V5O4010.D
 Report Date: 11-Jun-2013 09:23

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130607.B\V5O4010.D
 Lab Smp Id: M0903-09A Client Smp ID: MW105-060513
 Inj Date : 07-JUN-2013 22:30
 Operator : WL SRC: LIMS Inst ID: V5.i
 Smp Info : 5ML,M0903-09A,,72123
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
 Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V5O3805.D
 Als bottle: 22
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: BEX.sub
 Target Version: 4.14

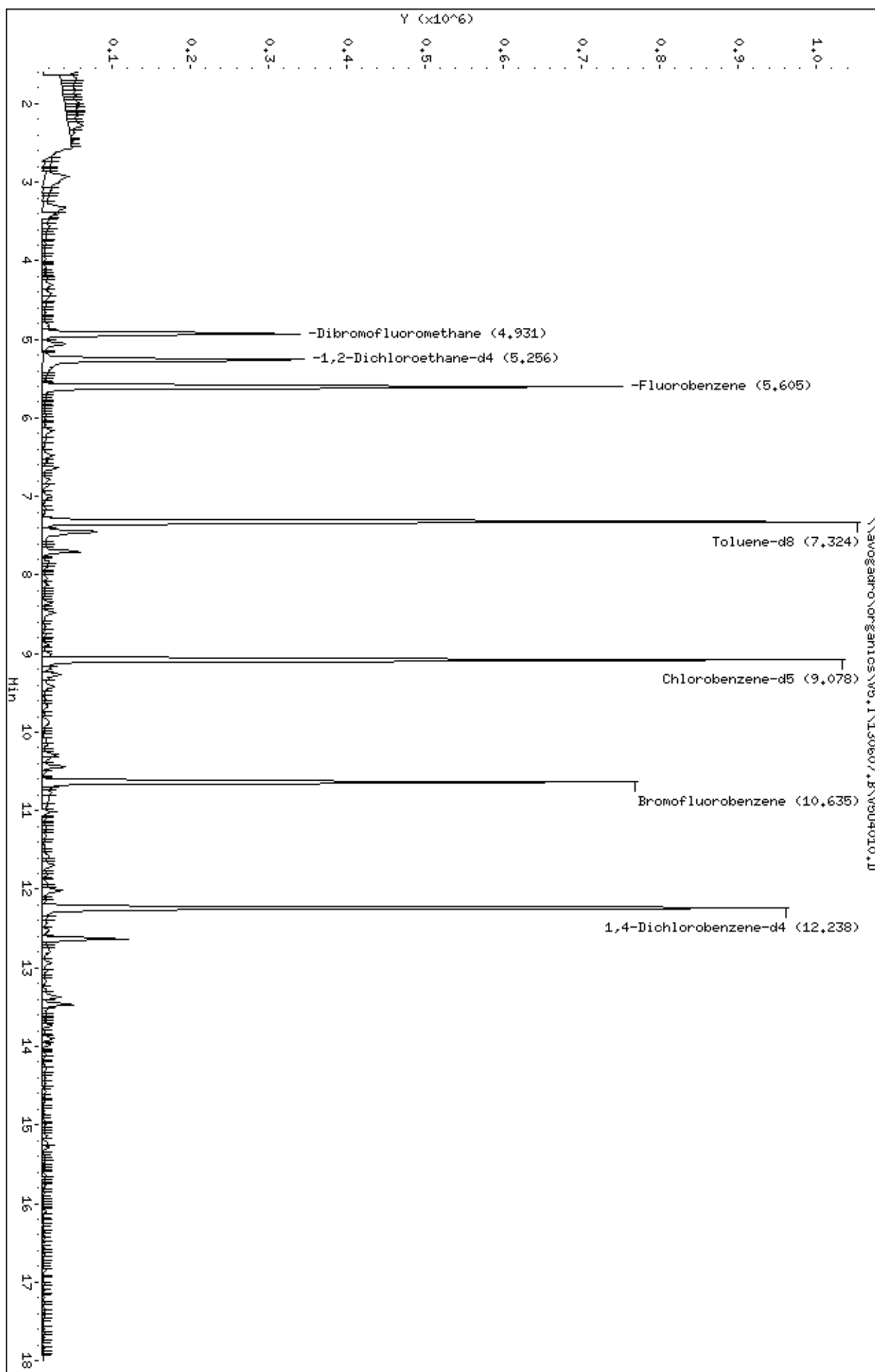
Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS	
						ON-COLUMN (ug/L)	FINAL (ug/L)
\$ 37 Dibromofluoromethane	113	4.930	4.929	(0.880)	245456	52.9913	53
\$ 43 1,2-Dichloroethane-d4	102	5.256	5.255	(0.938)	50615	49.0656	49
* 47 Fluorobenzene	96	5.604	5.603	(1.000)	746710	50.0000	
\$ 59 Toluene-d8	98	7.323	7.323	(0.807)	754903	47.1343	47
* 69 Chlorobenzene-d5	117	9.078	9.077	(1.000)	620570	50.0000	
\$ 80 Bromofluorobenzene	95	10.634	10.633	(1.171)	336878	52.4312	52
* 93 1,4-Dichlorobenzene-d4	152	12.237	12.236	(1.000)	269318	50.0000	

Data File: \\avogadro\organics\W5, I\130607.B\W504010.D
Date : 07-JUN-2013 22:30
Client ID: MM105-060513
Sample Info: 5HL, M0903-09H, 72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V504011.D
 Report Date: 11-Jun-2013 09:23

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130607.B\V504011.D
 Lab Smp Id: M0903-10A Client Smp ID: MW103-060513
 Inj Date : 07-JUN-2013 22:56
 Operator : WL SRC: LIMS Inst ID: V5.i
 Smp Info : 5ML,M0903-10A,,72123
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
 Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V503805.D
 Als bottle: 23
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: BEX.sub
 Target Version: 4.14

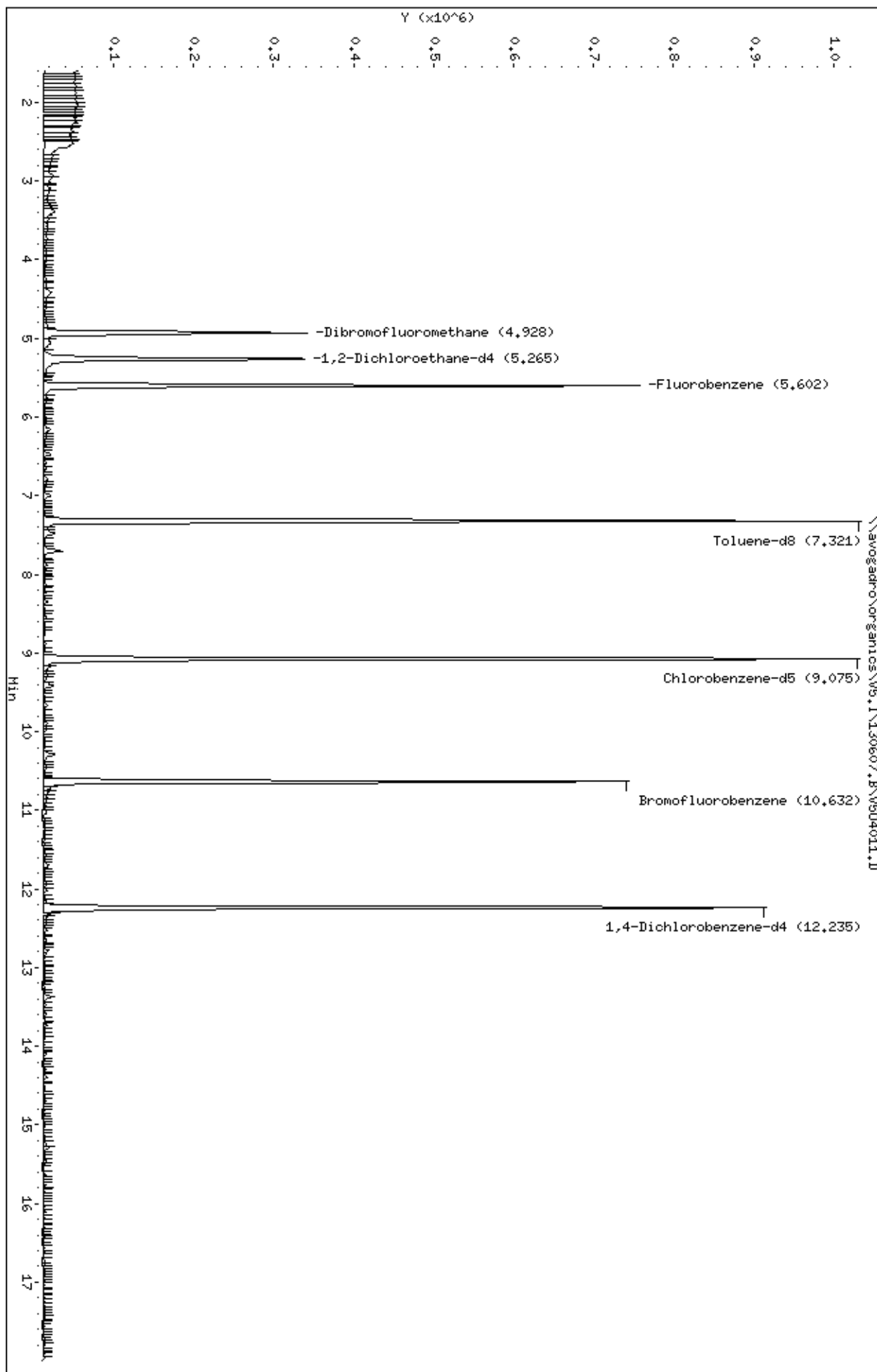
Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS	
						ON-COLUMN (ug/L)	FINAL (ug/L)
\$ 37 Dibromofluoromethane	113	4.927	4.929	(0.880)	242943	53.2175	53
\$ 43 1,2-Dichloroethane-d4	102	5.264	5.255	(0.940)	50384	49.5575	50
* 47 Fluorobenzene	96	5.601	5.603	(1.000)	735924	50.0000	
\$ 59 Toluene-d8	98	7.320	7.323	(0.807)	747166	47.1280	47
* 69 Chlorobenzene-d5	117	9.074	9.077	(1.000)	614291	50.0000	
\$ 80 Bromofluorobenzene	95	10.631	10.633	(1.172)	326867	51.3931	51
* 93 1,4-Dichlorobenzene-d4	152	12.234	12.236	(1.000)	257071	50.0000	

Data File: \\avogadro\organics\W5, I\130607.B\W504011.D
Date : 07-JUN-2013 22:56
Client ID: MM103-060513
Sample Info: 5HL, M0903-10H, 72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V5O4012.D
 Report Date: 11-Jun-2013 09:23

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130607.B\V5O4012.D
 Lab Smp Id: M0903-11A Client Smp ID: RW01-060513
 Inj Date : 07-JUN-2013 23:21
 Operator : WL SRC: LIMS Inst ID: V5.i
 Smp Info : 5ML,M0903-11A,,72123
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
 Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V5O3805.D
 Als bottle: 24
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: BEX.sub
 Target Version: 4.14

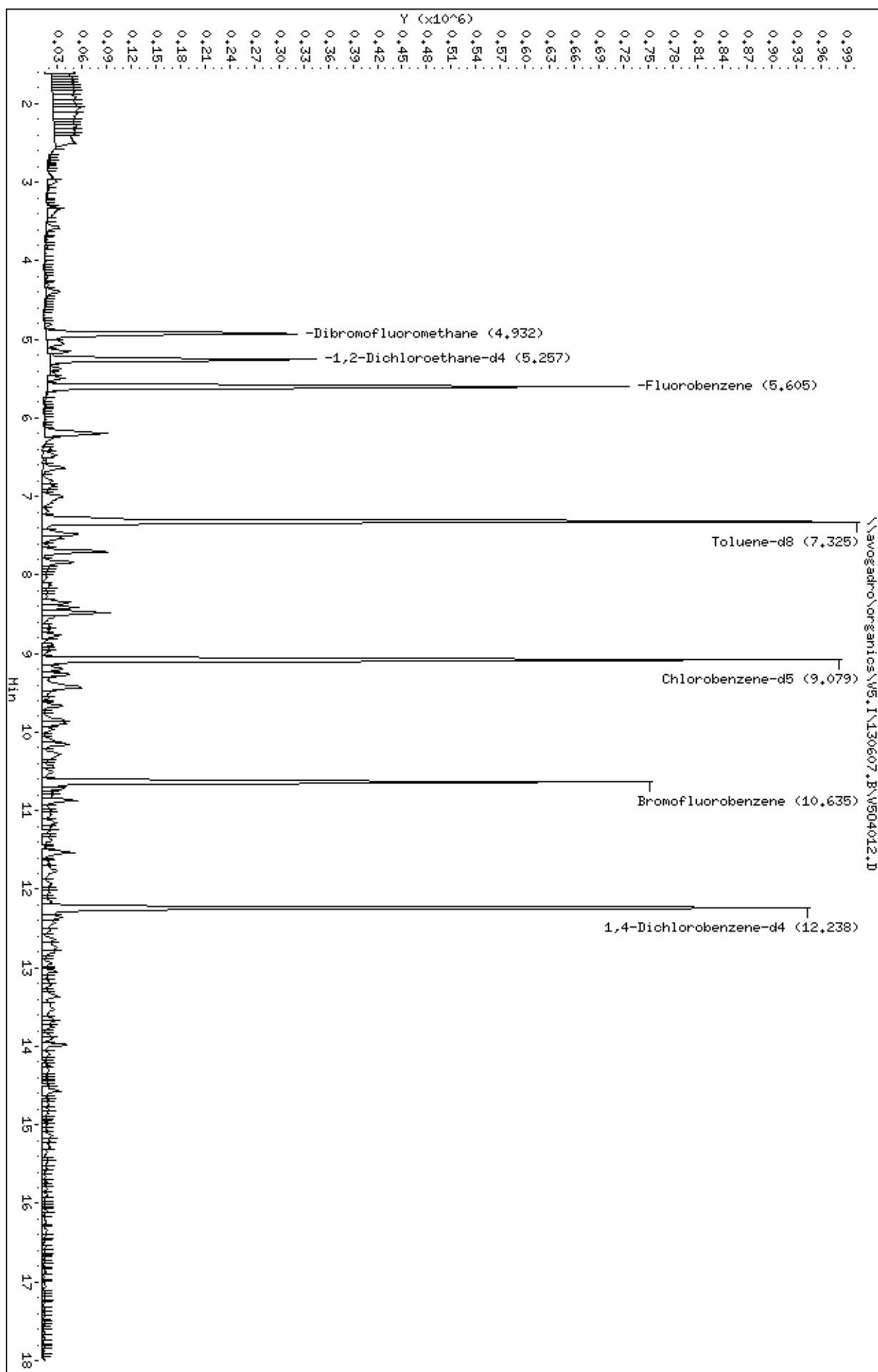
Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS	
						ON-COLUMN (ug/L)	FINAL (ug/L)
\$ 37 Dibromofluoromethane	113	4.931	4.929	(0.880)	243768	53.2538	53
\$ 43 1,2-Dichloroethane-d4	102	5.256	5.255	(0.938)	49193	48.2552	48
* 47 Fluorobenzene	96	5.605	5.603	(1.000)	737920	50.0000	
\$ 59 Toluene-d8	98	7.324	7.323	(0.807)	732035	46.9255	47
* 69 Chlorobenzene-d5	117	9.078	9.077	(1.000)	604449	50.0000	
\$ 80 Bromofluorobenzene	95	10.635	10.633	(1.171)	330499	52.8103	53
* 93 1,4-Dichlorobenzene-d4	152	12.238	12.236	(1.000)	260408	50.0000	

Data File: \\avogadro\organics\W5.1\130607.B\W504012.D
Date : 07-JUN-2013 23:21
Client ID: RM01-060513
Sample Info: 5HL, M0903-11A, 72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V504013.D
 Report Date: 11-Jun-2013 09:23

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130607.B\V504013.D
 Lab Smp Id: M0903-12A Client Smp ID: MW14-060513
 Inj Date : 07-JUN-2013 23:47
 Operator : WL SRC: LIMS Inst ID: V5.i
 Smp Info : 5ML,M0903-12A,,72123
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
 Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V503805.D
 Als bottle: 25
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: BEX.sub
 Target Version: 4.14

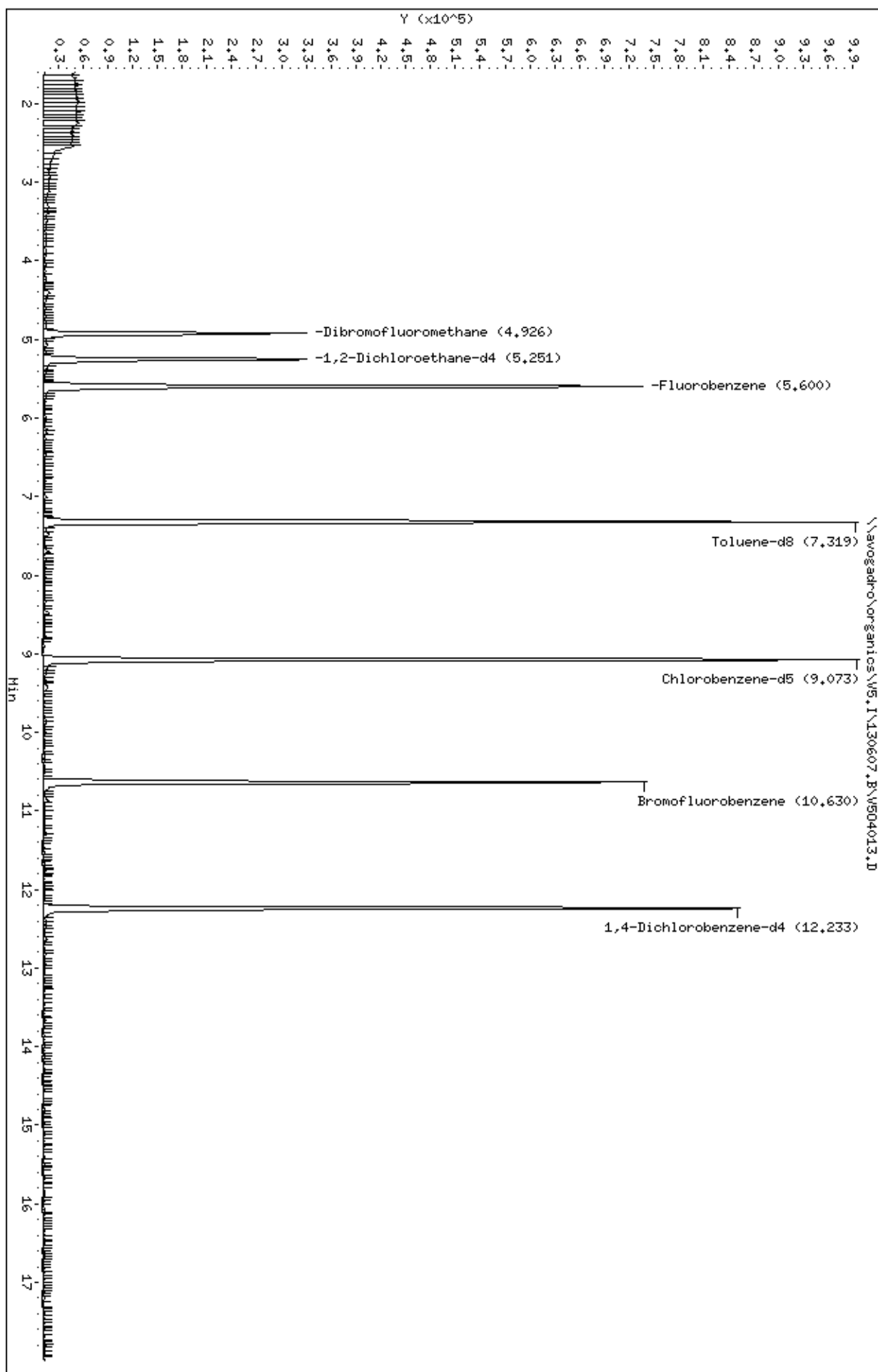
Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

Compounds	QUANT SIG	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS	
						ON-COLUMN (ug/L)	FINAL (ug/L)
\$ 37 Dibromofluoromethane	113	4.926	4.929	(0.880)	238720	52.8196	53
\$ 43 1,2-Dichloroethane-d4	102	5.262	5.255	(0.940)	49647	49.3248	49
* 47 Fluorobenzene	96	5.599	5.603	(1.000)	728580	50.0000	
\$ 59 Toluene-d8	98	7.319	7.323	(0.807)	725325	46.5121	46
* 69 Chlorobenzene-d5	117	9.073	9.077	(1.000)	604231	50.0000	
\$ 80 Bromofluorobenzene	95	10.629	10.633	(1.172)	320223	51.1867	51
* 93 1,4-Dichlorobenzene-d4	152	12.244	12.236	(1.000)	255556	50.0000	

Data File: \\avogadro\organics\W5, I\130607.B\W504013.D
Date : 07-JUN-2013 23:47
Client ID: MW14-060513
Sample Info: SWL, M0903-12H, 72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V5O4014.D
 Report Date: 11-Jun-2013 09:23

Spectrum Analytical, Inc. RI Division

Method 8260 Water and Medium Soil
 Data file : \\avogadro\organics\V5.I\130607.B\V5O4014.D
 Lab Smp Id: M0903-13A Client Smp ID: MW101-060513
 Inj Date : 08-JUN-2013 00:12
 Operator : WL SRC: LIMS Inst ID: V5.i
 Smp Info : 5ML,M0903-13A,,72123
 Misc Info :
 Comment :
 Method : \\avogadro\organics\V5.I\130607.B\v5_8260W.m
 Meth Date : 10-Jun-2013 15:40 V5.i Quant Type: ISTD
 Cal Date : 30-MAY-2013 14:32 Cal File: V5O3805.D
 Als bottle: 26
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: BEX.sub
 Target Version: 4.14

Concentration Formula: Amt * DF * Uf * 5/Vo * CpndVariable

Name	Value	Description
DF	1.000	Dilution Factor
Uf	1.000	ng unit correction factor
Vo	5.000	Sample Volume purged (mL)
Cpnd Variable		Local Compound Variable

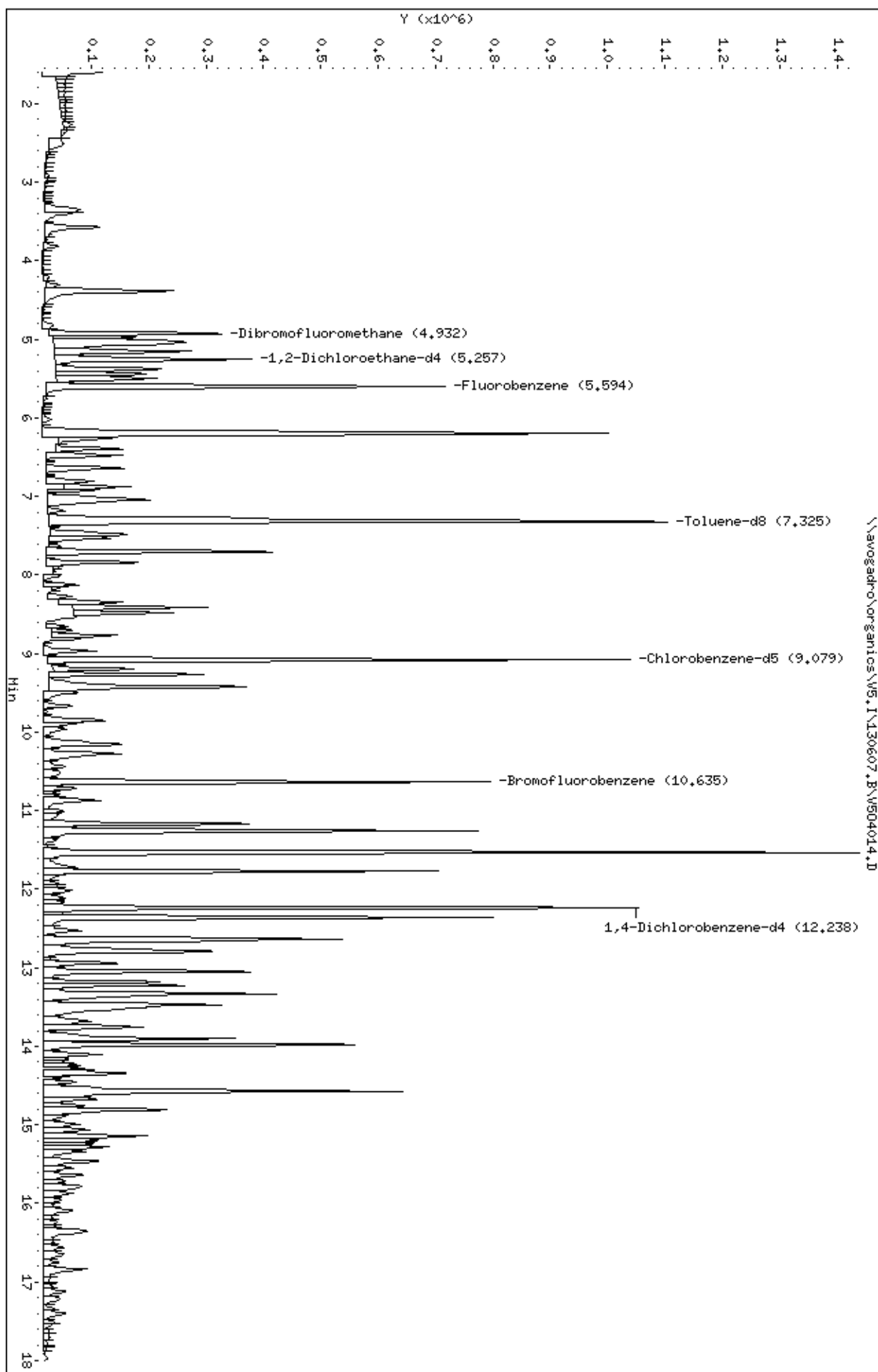
Compounds	QUANT SIG	RT	EXP RT	REL RT	RESPONSE	CONCENTRATIONS	
						ON-COLUMN	FINAL
	MASS					(ug/L)	(ug/L)
\$ 37 Dibromofluoromethane	113	4.931	4.929	(0.880)	243118	53.5297	54
\$ 43 1,2-Dichloroethane-d4	102	5.256	5.255	(0.938)	49892	49.3260	49
* 47 Fluorobenzene	96	5.605	5.603	(1.000)	732159	50.0000	
\$ 59 Toluene-d8	98	7.324	7.323	(0.807)	750763	47.5320	48
* 69 Chlorobenzene-d5	117	9.078	9.077	(1.000)	612002	50.0000	
73 Ethylbenzene	106	9.264	9.262	(1.020)	53679	5.74941	6(Q)
74 m,p-Xylene	106	9.427	9.413	(1.038)	81862	7.16467	7
\$ 80 Bromofluorobenzene	95	10.635	10.633	(1.171)	333298	52.6002	53
* 93 1,4-Dichlorobenzene-d4	152	12.238	12.236	(1.000)	272002	50.0000	
M 96 Xylene (Total)	106				81862	7.16467	7

QC Flag Legend

Q - Qualifier signal failed the ratio test.

Data File: \\avogadro\organics\W5.1\130607.B\W504014.D
Date : 08-JUN-2013 00:12
Client ID: MM101-060513
Sample Info: 5HL, M0903-13H, 72123
Purge Volume: 5.0
Column phase: DB-624

Instrument: W5.i
Operator: ML SRC: LIMS
Column diameter: 0.25



Data File: \\avogadro\organics\V5.I\130607.B\V504014.D

Date : 08-JUN-2013 00:12

Client ID: MW101-060513

Instrument: V5.i

Sample Info: 5HL,M0903-13A,,72123

Purge Volume: 5.0

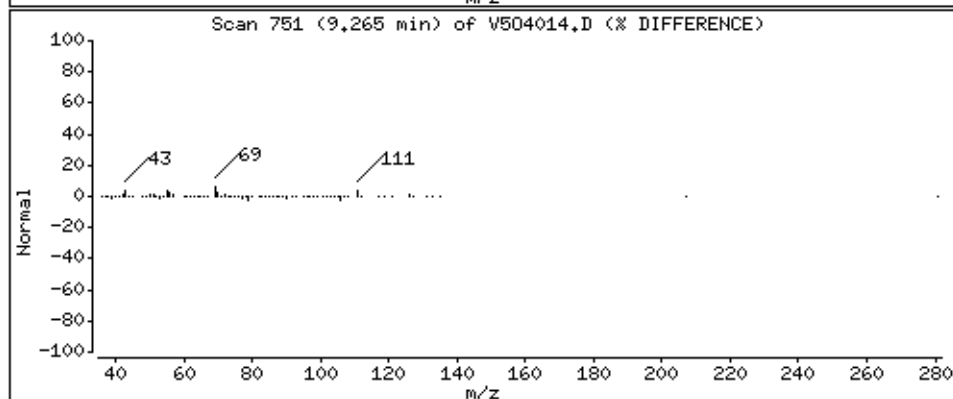
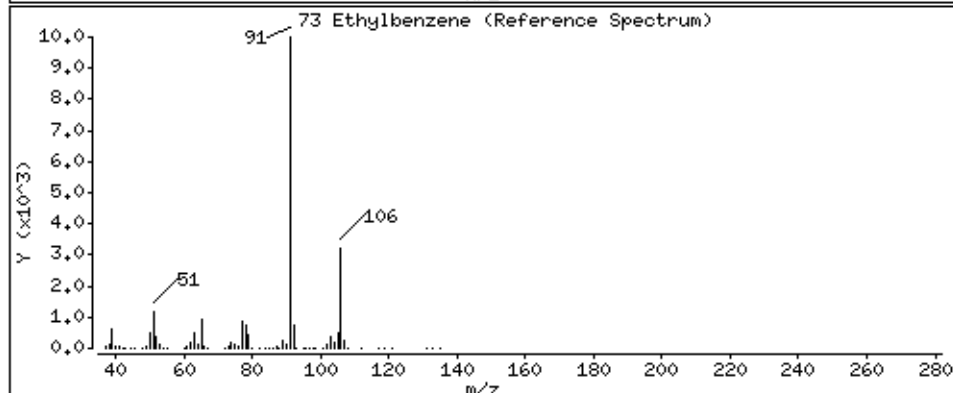
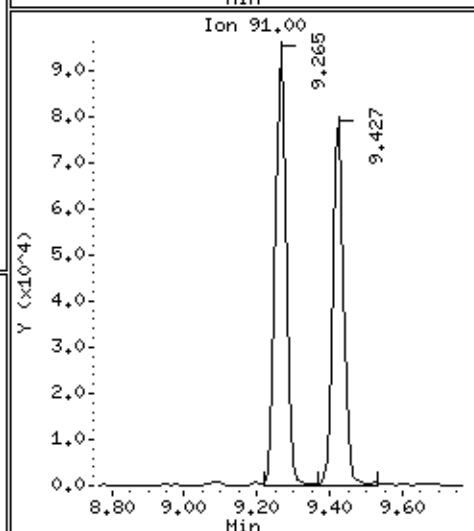
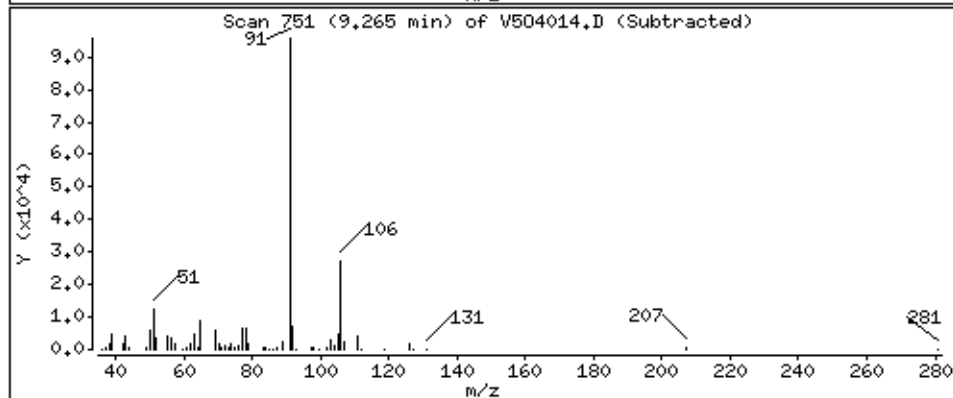
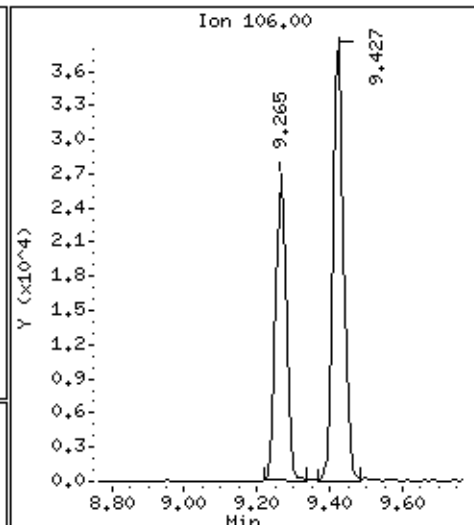
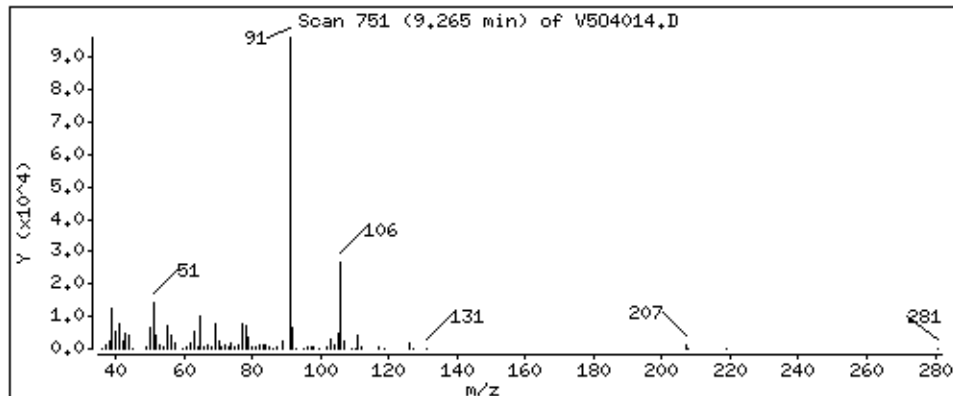
Operator: WL SRC: LIMS

Column phase: DB-624

Column diameter: 0.25

73 Ethylbenzene

Concentration: 6 ug/L



Data File: \\avogadro\organics\V5.I\130607.B\V504014.D

Date : 08-JUN-2013 00:12

Client ID: MW101-060513

Instrument: V5.i

Sample Info: 5HL,M0903-13A,,72123

Purge Volume: 5.0

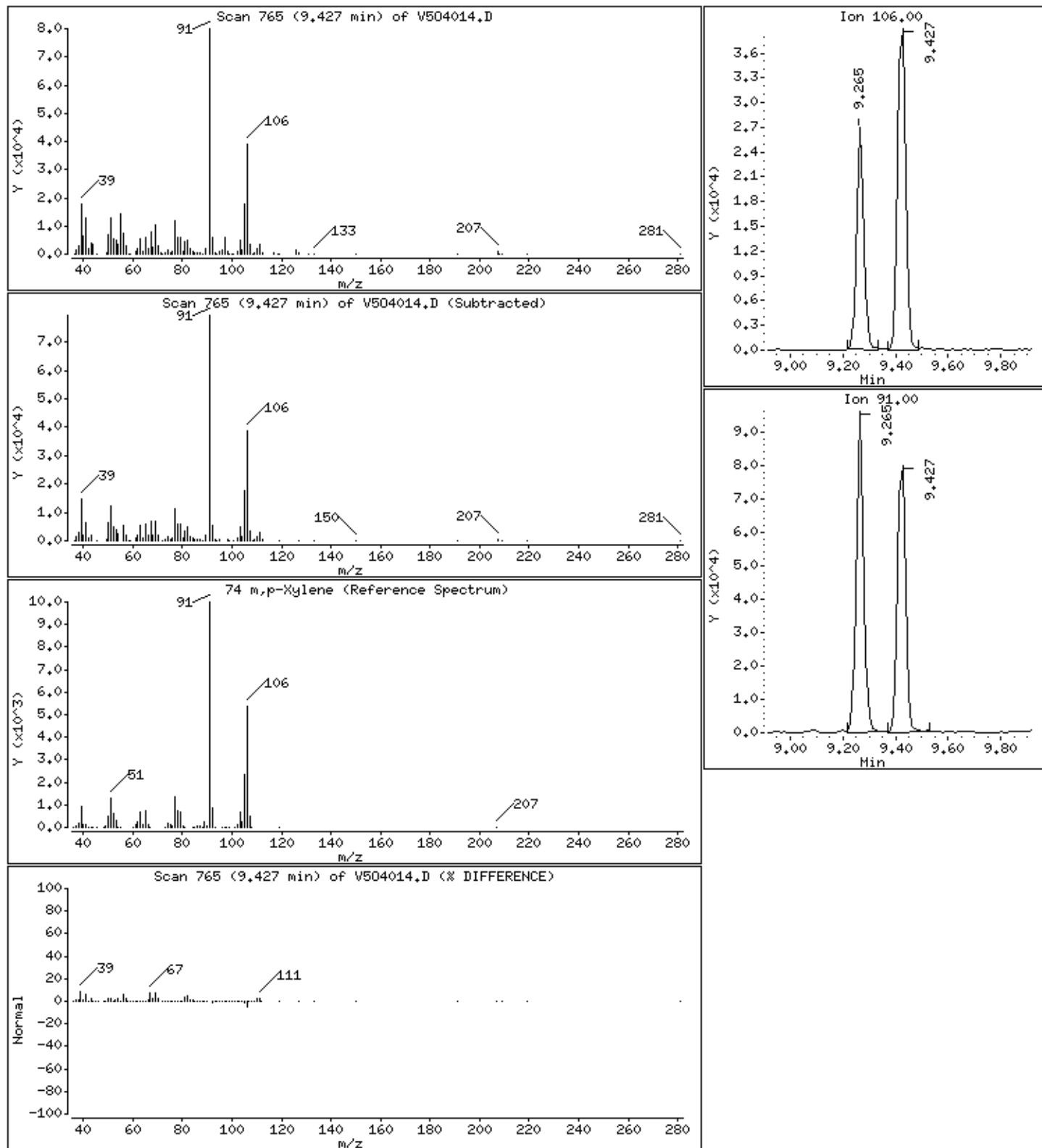
Operator: WL SRC: LIMS

Column phase: DB-624

Column diameter: 0.25

74 m,p-Xylene

Concentration: 7 ug/L



INJECTION LOG

M0903

<u>METHOD:</u>		<u>CAL ID:</u>	<u>ANALYST:</u>
<u>INITIAL CAL:</u>		<u>IS/SS ID:</u>	<u>ARCHIVE:</u>
<u>COMMENTS:</u>		<u>ICV ID:</u>	
SAMPLE			

Spectrum Analytical, Inc. RI Division V5 Injection Log Volatiles Laboratory
 Start: 30-MAY-13 12:22
 End: 30-MAY-13 15:46
 BATCH: 130530.B
 METHOD: 8260-W
 ANALYST: WL
 ICAL DATE: 5/30/13

Comments:

Standards: 2555-VW130521A uL
170-VW130521B uL
170-VW130521A uL
170-VW130521A uL

Reviewed By: 6-13 Manual Integration: NA MI Review: WL

FILE	TIME	LAB ID	CLIENT ID	PREP	MT	BN	INTERNAL	STDS	FBZ	CBZ	DCB	DFM	DCE	TOL	BFB	DILN	FLG	COMMENTS	pH
V503800	12:22	BFB5V		AQ															
V503801	12:48	VSTD0015V		AQ			99	97									1	OK	
V503802	13:14	VSTD0055V		AQ			99	98	97								1	OK	
V503803	13:39	VSTD0205V		AQ			100	98	99								1	OK	
V503804	14:05	VSTD0505V		AQ			100	100	100								1	OK	
V503805	14:32	VSTD1005V		AQ			64	62	59								1	OK	
V503806	14:58	VSTD2005V		AQ			61	60	60								1	OK	
V503807	15:46	VICV0505V		AQ			107	107	106			100	97	98	98		1	OK	

* - Internal Standard or Surrogate outside of control limits

E - One or more target compounds are above the calibration range

T - Sample was injected outside of the 12 hour sequence

R - One or more spike compounds are outside of control limits

D - Surrogates are diluted

1 WL 5/31/13

M0903

METHOD:

Spectrum Analytical, Inc. RI Division
Volatiles Laboratory

CAL ID:

METHOD: 8260-W
ICAL DATE: 5/30/13

ANALYST:

ANALYST: WL
BATCH: 130607.B

Start: 07-JUN-13 13:58
End: 08-JUN-13 01:32

Comments:

Standards: 25/55-WL130521A uL
5TD-WL130521B uL
BFB-WL130521A uL
uL

Reviewed By: ALW2
Manual Integration: NA
MI Review: WL

FILE	TIME	LAB ID	CLIENT ID	PREP	MT	EN	INTERNAL	STDS	SURROGATES	DILN	FLG	COMMENTS	pH
V503990	13:58	BFBDS	BFBDS										
V503991	14:23	VSTD050D5	VSTD050D5										
V503992	14:49	NB-72123	VBLKDS	72123	AQ		100	100				OK	
V503993	15:14	NB-72123	VBLKDS	72123	AQ		101	100	93	105	102	94	103
V503994	15:39	LCS-72123	VLCSDS	72123	AQ		98	98	91	106	101	93	104
V503995	16:05	M0903-06AMS	MW17-060413MS	72123	AQ	2	100	100	100	107	94	93	106
V503996	16:30	M0903-06AMS	MW17-060413MSD	72123	AQ	3	100	100	98	106	99	92	106
V503997	16:56	M0901-01A	SB70980-01	72123	AQ	1	102	99	95	106	99	95	106
V503998	17:21	M0902-01A	SB70984-01	72123	AQ	1	103	102	94	102	104	91	104
V503999	17:47	M0904-01A	016A	72124	AQ	1	103	102	94	102	104	91	104
V504000	18:13	M0907-01A	INFLUENT	72123	AQ	1	100	99	92	105	105	92	106
V504001	18:39	M0907-02A	EFFLUENT	72123	AQ	1	101	98	93	104	101	94	107
V504002	19:04	M0903-01A	MW112-060413	72123	AQ	1	100	98	91	107	102	94	106
V504003	19:30	M0903-02A	MW106-060413	72123	AQ	1	101	99	90	103	99	94	105
V504004	19:56	M0903-03A	DUP-060413	72123	AQ	1	98	96	88	105	96	94	105
V504005	20:22	M0903-04A	TB-060413	72123	AQ	1	98	97	88	104	96	92	102
V504006	20:48	M0903-05A	MW19-060413	72123	AQ	1	98	95	89	105	98	94	105
V504007	21:13	M0903-06A	MW17-060413	72123	AQ	1	96	94	85	104	100	93	105
V504008	21:39	M0903-07A	MW15-060413	72123	AQ	1	95	94	85	104	98	91	104
V504009	22:05	M0903-08A	MW11-060413	72123	AQ	1	93	93	84	107	95	93	103
V504010	22:30	M0903-09A	MW105-060513	72123	AQ	1	94	93	87	106	98	94	105
V504011	22:56	M0903-10A	MW103-060513	72123	AQ	1	93	92	83	106	99	94	103
V504012	23:21	M0903-11A	RW01-060513	72123	AQ	1	93	91	84	106	96	94	106
V504013	23:47	M0903-12A	MW14-060513	72123	AQ	1	92	91	82	106	99	93	102
V504014	00:12	M0903-13A	MW101-060513	72123	AQ	1	92	92	87	107	99	95	105
V504015	00:38	M0910-01A	IDW-AQ-03	72123	AQ	1	100	97	90	103	97	93	104
V504017	01:32	M0907-01A	INFLUENT	72123	AQ	2	131	102	87	102	100	102	99

Internal Standard or Surrogate outside of control limits
One or more target compounds are above the calibration range
Sample was injected outside of the 12 hour sequence

R - One or more spike compounds are outside of control limits
D - Surrogates are diluted

WL 6/10/13

M0903

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Reviewed By: W 6/11/13

UA = Unpreserved Aqueous

M = MeOH

A = Air

$$H = HCL$$

US = Unpreserved Soil

F = Freeze

$$N = \text{NaHSO}_4$$

T = Trace, HCL

Last Page of Data Report