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# **SITE CHARACTERIZATION REPORT FOR THE FORMER HUNTS POINT GAS WORKS HALLECK STREET SIDEWALK AREA**

**VCA Index No. D2-0003-02-08**

**Site #V00554**

**Bronx, New York**

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*Prepared For:*



**Consolidated Edison Company of New York, Inc.**

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**JULY 2014**

*"I, Shane Blauvelt, certify that I am currently a Qualified Environmental Professional as defined in 6 NYCRR Part 375 and that this Report was prepared in accordance with all applicable statutes and regulations and in substantial conformance with the DER Technical Guidance for Site Investigation and Remediation (DER-10) and that all activities were performed in full accordance with the DER-approved work plan and any DER-approved modifications."*



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# **SECTION 1**

## **INTRODUCTION**

### **1.1 SITE CHARACTERIZATION STUDY OBJECTIVES**

Site Characterization of the Halleck Street Sidewalk Area was conducted to (1) assess the potential presence of MGP-related impacts; and (2) to ascertain the potential need for further investigation or remediation. This Site Characterization Report (SCR) addresses only the sidewalk and right of way of Halleck Street (Site), also known as Hunts Point MGP OU-3. The specific objectives of this SCR are to assess whether hazardous substances have been released to the environment and may be present onsite, if they may have migrated offsite, and whether they may have impacted human health or the environment. If no potential impacts are identified, a “no further action” conclusion may be warranted. If potential impacts are verified, additional sampling may be needed to determine the nature and extent of those impacts, or the need for remediation and interim measures to address the impacts. These objectives are consistent with those of the New York State Department of Environmental Conservation’s (NYSDEC) comprehensive remedial investigation process, specifically Chapter 3 of the NYSDEC DER-10 Technical Guidance for Site Investigation and Remediation (NYSDEC, 2010).

### **1.2 REPORT ORGANIZATION**

The Site Characterization was conducted by Parsons in October and November 2013. The field investigation activities and results are documented in this report in the following sections and appendices:

- Section 1: Introduction
  - Section 2: Site Background
  - Section 3: Site Characterization Activities
  - Section 4: Site Characterization Results
  - Section 5: Exposure Assessment
  - Section 6: Conclusions
  - Section 7: References
- 
- Appendix A: Soil Boring and Monitoring Well Logs
  - Appendix B: Groundwater Sampling Logs
  - Appendix C: Data Usability Summary Report
  - Appendix D: Hydrocarbon Fingerprint Results

## **SECTION 2**

### **SITE BACKGROUND**

#### **2.1 SITE OVERVIEW**

The Consolidated Edison Company of New York, Inc. (Con Edison) has entered into a Voluntary Cleanup Agreement with the NYSDEC to investigate, and if necessary, remediate potential impacts at former manufactured gas plant (MGP) properties. One of these facilities is the former Hunts Point Gas Works (VCA Index No. D2-003-02-08; Site #V00554), located in a commercial area of the Bronx, New York ([Figure 1](#)). The former Hunts Point Gas Works (Gas Works) was a MGP operated by Con Edison between 1926 and 1962. Currently, the former Hunts Point Gas Works property is owned by the City of New York for use as the Hunts Point Food Distribution Center. The Site includes the street and sidewalk area of Halleck Street and is approximately 2,500 linear feet long, within the east right of way of Halleck Street, also known as Hunts Point MGP OU-3([Figure 2](#)).

Historical research conducted for the former Hunts Point Gas Works is further described in Section 2.5. Results of those research efforts are documented in the *Hunts Point Offsite Manufactured Gas Plant Site History Report, Bronx, New York* (Parsons, 2003). Portions of the former Hunts Point Gas Works were previously investigated and/or remediated.

#### **2.2 ADJOINING PROPERTY DESCRIPTION**

The Site is bound by Food Center Drive to the north and a Department of Corrections facility to the south. The Site is bordered to the east by the New Fulton Fish Market Cooperative and Hunts Point Cooperative Market Complex. Commercial buildings border the Site to the west. The properties east of the Site are currently owned by the City of New York and are managed by the New York City Economic Development Corporation.

At the northeastern end of the Site is the Con Edison Hunts Point Gas Compressor Station (HPGCS).

#### **2.3 SITE HISTORY**

Historical research was previously conducted and documented in the *Hunts Point Offsite Manufactured Gas Plant Site History Report, Bronx, New York* (Parsons, 2003). Based on this report, the Gas Works was owned and/or operated as an MGP and gas holder station by Con Edison between 1926 and 1962. A large gas holder on the northern portion of the Gas Works was used until 1968, after which it was removed from service and demolished.

The City of New York acquired the majority of the former Hunts Point Gas Work property in the late 1960s. The property was then transitioned into warehouse space for a wholesale food cooperative. One of the former facility roads (Hunts Point Avenue) was removed from service as part of the property redevelopment. Reviews of historical aerial photos reveal that Halleck Street was not modified as part of redevelopment. It was extended to the south sometime after 1980 for use as an access road to the current Department of Corrections facility, located at the southwest portion of the Hunts Point Peninsula.

## **2.4 TOPOGRAPHY, REGIONAL GEOLOGY, AND HYDROGEOLOGY**

Hunts Point is a peninsula on the East River and Bronx River that is surrounded by brackish or salty tidal water. The former Hunts Point Gas Works covered an area of approximately 182 acres. Halleck Street is the western boundary of the former site operations and extends for approximately 2,500 linear feet.

The site-specific geology and subsurface site conditions can vary depending on the local site history and the specific activities conducted (construction, excavation, filling, etc.). Prior to significant construction and development, Hunts Point was drained by small creeks, which emptied into the Bronx and East rivers. Most of these creeks are now filled in and are covered by buildings and streets. However, the movement of the shallow groundwater is somewhat influenced by the old channels as well as ancient former stream deposits and the areas of sewer lines and former piping. Groundwater in the area has been reported to occur within the shallow subsurface, and flows in a southerly direction toward the Hunts Point promontory, and the confluence of the Bronx and East Rivers (Hygienetics, 1997b).

During the PSA at the HPGCS, shallow water was encountered within the fill materials on the eastern portion of the property at depths between two and five feet bgs. This shallow water was not encountered at drilling locations on the western half of the Hunts Point Gas Works and appears to be perched water within the fill and sand materials above the clay. A deeper water-bearing zone was encountered during the subsurface investigation within a deeper sand layer at approximately 12 to 15 feet bgs. Boring logs from the Hygienetics report indicate the presence of groundwater ranged from 3 to 9 feet bgs depending on the proximity to the Bronx River (Hygienetics, 1997b and LMS, 1999a and b).

## **2.5 PREVIOUS INVESTIGATIONS/REMEDIAL MEASURES**

Three of the previously investigated parcels (A, B, and E) are immediately adjacent to Halleck Street. From 1997 through 1999, Hygienetics Environmental Services Inc. (HES) performed investigations on various parcels of the Gas Works, results of which were summarized in the following report:

- *Phase I Environmental Site Assessment Report of Hunts Point Produce Market Complex Parcels A, B, C, D, and E* (HES, 1997a)
- *Phase II Environmental Investigation Report of Hunts Point Market Complex Parcels A, B, C, D, and E* (HES, 1997b)

In addition, Lawler, Matusky, and Skelly Engineers LLP (LMS) conducted site investigations at selected parcels, results of which were summarized in the following reports:

- *Investigation Report for the Operating Unit Portion of Parcel A* (LMS, 1999a)
- *Investigation Report for the Operating Unit Portion of Parcel E* (LMS, 1999b)
- *Investigation Report for the Operating Unit Portion of Parcel C* (LMS, 1999c)
- *Investigation Report for the Operating Unit Portion of Parcel B* (LMS, 2001)
- *Hunts Point Cooperative Market Redevelopment Plan, Investigation Report for Parcel D, Bronx, New York* (LMS 2005)

- *Hunts Point Cooperative Market Redevelopment Plan, Investigation Report, Operable Unit 3 of Parcel E, Bronx, New York, Final* (HDR/LMS 2007a)
- *Hunts Point Cooperative Market Redevelopment Plan, Investigation Report for Parcel F, Bronx, New York, Final* (HDR/LMS 2007b)

Parsons conducted remedial investigation at the HPCGS, also located immediately adjacent to Halleck Street, in 2003. Results of that investigation were summarized in the following report:

- *Site Investigation Report for the Hunts Point Gas Regulator Station Site Investigation* (Parsons, 2003)

A brief description of the investigation on those Parcels follows here. For detailed sampling locations refer to the listed reports.

## **PARCEL A**

Previous investigations on Parcel A were divided into two operable units: Operable Unit-1 (OU-1) and Operable Unit-2 (OU-2). The OU-1 investigation was completed in 1999 and included five test pits, one soil boring, and two groundwater grab samples from two of the test pits.

Test pits encountered mixed soils, sand, gravel, garbage, structural materials, coal slag, incinerator ash, and coal tar residues. Dense coal tar was detected in one test pit (Test Pit 2). No free phase oil and no purifier material were observed during this investigation. Several polycyclic aromatic hydrocarbon (PAH) compounds exceeded TAGM concentrations in soils collected during this investigation. Metals concentrations were reportedly consistent with urban fill materials but arsenic, barium, beryllium, cadmium, chromium, copper, iron, mercury, nickel, selenium, and zinc exceeded the TAGM Recommended Soil Cleanup Objectives in at least one soil sample. LMS concluded the constituents present should not preclude development of Parcel A. New pavement would cap the materials below (LMS, 1999a).

The OU-2 investigation was completed in 2003 and included the installation of five test pits, three temporary piezometers, three deep soil borings, and thirteen shallow soil borings. Coal tar was observed at OU-2 between 3 and 4 feet bgs. Tar boils were also observed in thicknesses ranging from inches to several feet. Purifier waste, consisting of wood shavings with a blue/green color, was encountered at the ground surface to a maximum depth of five feet bgs within the eastern portion of Parcel A. The recommended remedial action of Parcel A was excavation of coal tar and purifier wastes (LMS, 2003).

## **PARCEL B**

Two investigations were completed at Parcel B, the first in 1998 and the second in 2001. The 1998 investigation included the installation of one monitoring well and nine soil borings. The 2001 investigation consisted of 11 test trenches, 14 temporary piezometers, 1 monitoring well, 9 soil borings, 5 test pits, 4 surface soil samples, and 40 geotechnical borings. The investigations encountered wood shavings attributed to residual purifier materials, incinerator ash, coal tar, coke gravel, and slag throughout Parcel B. Tar boils were also present. Tar, where encountered, was several inches to several feet in thickness. One tar boil, located in the northwestern part of Parcel B, was 62 feet wide, by 68 feet long, by 3 feet deep. A smaller tar boil was located north of the NYDOC parking area.

Exceedances of TAGM values in soil, fill, and groundwater were generally coincident with areas where gross constituents were observed. Select SVOCs and metal were detected above the soil cleanup criteria in the surface soil samples of Parcel B. Total VOCs, BTEX, SVOCs were detected above NYSDEC Class GA Standards. One instance of LNAPL was observed and most groundwater contained either sheens and/or odors.

Since the investigations described above occurred, Parcel B was remediated and developed as the new center of the Fulton Fish Market, a 325,000 square foot building with a concrete slab on-grade construction. The surrounding areas have been developed as a parking lot.

## **PARCEL E**

Previous investigations of Parcel E were conducted in 1997 and 2007. Parcel E is located in the area of the former 15,000,000 cubic foot gasholder, coke filters, and substation structures. Parcel E was subdivided into three Operable Units (OU-1, OU-2 and OU-3). The 1997 investigation included the installation of 5 soil borings and 1 monitoring well. The 2007 investigation included the installation of 21 soil borings, 4 trenches, 5 test pits, 5 piezometers, and a GPR survey. The 2007 investigation also included the collection and analysis of soil, soil gas, and groundwater samples (one from Trench 2 and one from Trench 4B). The investigations of Parcel E encountered historical utilities, piping, and tank structures during test pit excavations and trenching. Some of the tank structures and piping contained oily product and the adjacent soils were impacted by the product, other structures and piping contained water. The test pits and trenches encountered an upper layer of fill material consisting of mixed soils, structural materials, coal slag, C&D debris, garbage, brick, and glass. Tar and oil-saturated soil with strong naphthalene odors was encountered in places. Groundwater in the trenches had sheens and small globules of product.

Benzene, total xylenes, SVOCs and metals were detected above recommended soil cleanup objectives in samples collected from the trenches (Trench 2, 3, and 4) and metals were detected above cleanup objectives in all four trenches. Only cyanide exceeded groundwater standards in one sample. The recommended remedy for Parcel E was excavation of impacted soil and wastewater, installation of an engineered cap, and 1 foot of imported fill for landscaped areas.

## **HPGCS**

A total of 47 soil samples were collected from soil borings drilled at the gas regulator station during the initial site investigation (SI). Four groundwater samples were collected. An additional 11 soil samples, 2 groundwater samples, and 3 tar samples were collected during supplemental sampling.

Three to eleven feet of fill was encountered across this parcel, consisting of sand, gravel, coal ash/cinders, and debris (wood, bricks, glass, etc.). Evidence of non-aqueous phase liquid (NAPL) and tar were observed during this investigation. Metals were detected throughout the HPGCS in soil and groundwater at concentrations exceeding the NYSDEC recommended TAGM soil cleanup objectives and/or Eastern U.S. background levels. The elevated metal concentrations are consistent with the historical industrial use of the HPGCS and surrounding area.

Analytical results indicated that elevated PAH and PCB levels were present in the HPCGS soils. As such, this parcel was remediated in 2003 to:

- Excavate soils with total PAH concentrations greater than 500 ppm.
- Remediate soils visibly impacted with MGP-related tar or free product; and
- Mitigate the potential exposure to soils remaining following the remedial actions

Remedial activities for this parcel are documented in the *Final Engineering Report, Hunts Point Gas Compressor Station Site, Bronx, NY* (Parsons, 2004).

## SECTION 3

### SITE CHARACTERIZATION ACTIVITIES

The following sections describe the field investigation activities conducted as part of the Site Characterization. Parsons personnel mobilized to the Site on October 7, 2013, and the field investigation activities were conducted between October and November 2013 in accordance with the NYSDEC approved *Site Characterization Work Plan* (work plan) (Parsons, 2011). The scope of field investigation activities included the installation of soil borings and monitoring wells. Soil, groundwater, and non-aqueous phase liquid (NAPL) samples were collected for laboratory analysis. During all intrusive activities, a Community Air Monitoring Plan (CAMP) was implemented in accordance with the approved work plan. Sample locations are shown on Figure 3. Table 1 provides a summary of the samples and analyses.

#### **3.1 SITE INSPECTION AND PRELIMINARY INVESTIGATION ACTIVITIES**

On October 7, 2013, a Site inspection was conducted to refine the locations of the proposed investigation points. The proposed scope of work was reviewed with Con Edison. Proposed locations and proposed methods were altered in the field, as necessary, based on Site conditions, access, utilities, and safety. The modifications to the sampling locations are further discussed below. Sampling location changes were made in consultation with Con Edison and the NYSDEC.

#### **3.2 UTILITY CLEARANCE**

A geophysical survey was conducted to identify potential/possible underground conduits/utilities in the area of the proposed soil boring and monitoring well locations. The geophysical survey was completed by Naeva Geophysics Inc. (Naeva), of Congers, New York prior to start of Site work.

Once the initial geophysical survey was completed, utility clearance keyhole test pits were hand or vacuum excavated at each proposed soil boring and monitoring well location for subsurface utilities. Utility clearance test pits were completed by Aquifer Drilling & Testing, Inc. (ADT) of New Hyde Park, New York in October 2013. The typical utility clearance test pit excavation consisted of saw-cutting and jack-hammering the surface pavement (as necessary), and excavating using a Vactron, an air knife, and hand tools (as necessary) to a minimum depth of 5 feet below ground surface (bgs). During these excavation activities, soils were screened for VOCs using a photoionization detector (PID), their physical characteristics (e.g., soil type, grain size, color, etc.) were described, and notes of any evidence of physical impacts observed (staining, odor, sheen, non-aqueous phase liquid (NAPL), etc.) were recorded. When a utility clearance test pit could not be completed to a depth of 5 ft bgs due to the presence of underground utilities or subsurface obstructions, the location was moved approximately 5 to 10 feet away from the original location and re-excavated. Following completion of the utility clearance test pits, each test pit was backfilled prior to drilling or excavation.

### **3.3 SOIL BORING INSTALLATION**

A total of twelve (12) soil borings (MW-1, SB-1 through SB-11) were advanced during the Site Characterization activities to characterize subsurface conditions. The soil borings were completed from October 2013 through November 2013. Advancement of the soil borings was conducted by ADT under the supervision of a Parsons geologist. Soil borings were completed to depths ranging from approximately 15 to 25 feet bgs, depending on observed impacts and refusals. [Figure 3](#) shows the soil boring locations and the corresponding boring logs are presented in [Appendix A](#). Based on Site conditions and observations made during implementation of the Site Characterization activities, the following additions/modifications were made to the work plan:

- The SB-02 location was shifted 3 feet south after the hand-clearance experienced refusal at 2.5 feet bgs. SB-02 hand-clearance experienced refusal again at 2.5 feet bgs and was moved 3 feet north of the original location. The third hand-clearance attempt was successful.

Soil borings were advanced using a track-mounted Geoprobe ® Direct Push Drill Rig. Soil samples were collected continuously to the bottom of the boring using 5-foot long, 2.25-inch outer diameter Macro-core ® barrels containing a 2-inch acetate lined sampling sleeve. Each sample was screened for the presence of VOCs using a PID. Soil was also logged for physical characteristics of each sample (e.g., soil type, color, texture, moisture content, etc.), along with physical evidence of any impacted material (e.g., oil-like or tar-like NAPL, staining, sheens, odors, etc.).

Soil samples were submitted to Chemtech and analyzed for TCL VOCs, TCL SVOCs, cyanide, and TAL metals. A summary of the soil samples collected and the analyses performed is provided in [Table 1](#). Soil samples were collected from selected zones within the borings (as described below) and were submitted for laboratory analysis:

- One sample was collected from the zone with the highest PID readings or visual impacts. If visual impacts or elevated PID readings were not observed, a sample was collected from the upper portion of the boring or directly above the water table (if present).
- One sample was collected below the impacted zone (if present) or near the base of the boring to identify the vertical extent of any impacts at the location.
- Where applicable, NAPL samples were collected and submitted to META Environmental, Inc. of Watertown, MA for hydrocarbon fingerprint analysis.

Upon completion, the boring locations were grouted with Portland cement and bentonite grout using a tremie pipe. Drilling equipment was decontaminated between each boring. Drill cuttings and decontamination water were containerized in 55-gallon steel drums and handled as described in Section 3.7.

### **3.4 MONITORING WELL INSTALLATION/DEVELOPMENT**

A total of three (3) monitoring wells (MW-1 through MW-3) were installed during the Site Characterization activities. Monitoring wells were installed in November 2013. The original MW-1 location did not reveal any water or moisture in the soil boring and the soils containing native clays were completely dry. It was proposed by Parsons to move MW-1 approximately 240 feet to the south to be installed at the SB-1 location, which showed moist soils in the 8 to 12 ft

bgs range. As discussed in the e-mail correspondence from November 5, 2013, this proposed change was approved by the New York State DEC since the change still satisfied the objective in the work plan to characterize the groundwater at the edge of Parcel D, along Halleck Street.

The monitoring well borings were advanced to varying depths before later being converted into monitoring wells with the well installation portion of the drilling work. The MW-1 boring was advanced to 15 ft bgs at the initial location. Since no water was produced at the initial location, MW-1 was moved to the SB-1 location, which was advanced to 20 ft bgs. The MW-1 monitoring well was installed to a depth of 17 ft bgs at the SB-1 location. The SB-5/MW-2 boring location was advanced to 15 ft bgs and then converted with the installation of the MW-2 monitoring well, which was installed to 16 ft bgs. The SB-11/MW-3 boring location was advanced to 25 ft bgs and then converted with the installation of the MW-3 monitoring well, which was installed to 17 ft bgs. Monitoring well borings were completed with 4.25-inch outside diameter hollow stem augers and a track-mounted Geoprobe ® drill rig. The monitoring well screens were set at varying depths: MW-1 was screened from 7 to 17 ft bgs, MW-2 was screened from 6 to 16 ft bgs, and MW-3 was screened from 7 to 17 ft bgs with the top of the screen approximately 2 feet above the observed groundwater table.

The monitoring wells were constructed with 2-inch inner diameter, threaded, flush-joint, PVC casing and 10-foot lengths of 0.02-inch slot screen. The annular space around each well screen was backfilled with a No. 2 sand filter pack extending from the bottom of the well to at least 2 feet above the top of the screen. The annular space around the well riser was sealed with at least 2 feet of hydrated bentonite pellets on top of the sand pack. The remainder of the boring was backfilled with cement-bentonite grout to approximately 4 to 5 feet below grade. Monitoring wells MW-1, MW-2, and MW-3 were all finished with a locking, flush-mount box set in concrete.

Monitoring well development was conducted in November 2013 a minimum of 24 hours after installation. Monitoring wells were developed until reasonably free of sediment (less than 50 NTU if possible) or until the pH, temperature, Oxygen Reduction Potential (ORP), and conductivity stabilized. Monitoring well development was monitored approximately every 5 minutes by reviewing water quality indicator measurements. Well development continued until turbidity was less than 50 nephelometric turbidity units (NTUs) for three successive readings or until water quality indicators stabilized, whichever occurred first in each monitoring well. The stabilization criteria were based on water quality indicators of three successive readings within 10%.

Non-disposable drilling equipment was decontaminated between monitoring well locations. Monitoring well drill cuttings, well development water, and decontamination water were containerized in 55-gallon steel drums and handled as described in Section 3.7.

### **3.5 SURVEYING**

At the conclusion of drilling activities, Chazen, a licensed New York state land surveyor, mobilized to the Site and identified the horizontal and vertical location of each new soil boring and monitoring well. Additionally, the survey included locating Site features such as manholes, bollards, hydrants, telephone poles, and more. Two elevation measurements were taken at each well location to identify the top of the PVC casing and the grade elevation. The survey elevations

were measured to an accuracy of 0.01 feet above the National Geodetic Vertical Datum of 1988 (NGVD 1988).

### **3.6 GROUNDWATER SAMPLING**

On November 27, 2013, groundwater samples were collected from the three (3) monitoring wells (MW-1 through MW-3). Prior to collecting samples, the depth to groundwater and thickness of any free product (if present) was measured in the monitoring wells using an electronic oil/water interface probe attached to a measuring tape accurate to 0.01 feet. [Table 2](#) provides a summary of the groundwater level measurements and elevations.

Prior to purging, the headspace within each well was measured with a PID. Each well was purged using a submersible pump and low-flow purging techniques to stabilize the following water quality parameters: temperature, conductivity, pH, dissolved oxygen, conductivity, oxidation reduction potential (ORP), and turbidity; which were measured approximately every five minutes.

Once stabilization was achieved, groundwater samples were collected using a low-flow submersible pump with dedicated tubing. Water quality parameter measurements and observations recorded during sampling activities are documented on the groundwater sampling records provided in [Appendix B](#). Groundwater samples were submitted to Chemtech Laboratories for the following analysis: TCL VOCs, TCL SVOCs, TAL Metals, and total cyanide. Non-dedicated sampling equipment (e.g., oil/water interface probe, submersible pump) was decontaminated between sampling locations. Decontamination water was placed in 55-gallon drums and handled as described in Section 3.7.

### **3.7 MANAGEMENT OF INVESTIGATION-DERIVED WASTE**

Investigation-derived waste (IDW), which included decontamination wash and rinse water, soil cuttings, purge water, debris, and used personal protective equipment (PPE), was containerized in Department of Transportation (DOT)-approved 55-gallon drums. The drums were sealed at the end of each work day and labeled with the date, the well or boring number(s), and the type of waste (e.g., drill cuttings, purge water). Parsons collected representative waste characterization samples of the IDW and coordinated transportation and disposal. Clean Earth of North Jersey, Inc. from Kearny, New Jersey disposed of the IDW at an offsite Con Edison-approved location in accordance with applicable local, state, and federal regulations.

### **3.8 DATA VALIDATION AND REPORTING**

Data validation was performed in accordance with the USEPA Region II standard operating procedures (SOPs) for organic and inorganic data review which were in effect at the time of data validation (USEPA 2006; 2008a; 2008b). These validation guidelines are regional modifications to the National Functional Guidelines for organic and inorganic data review (USEPA, 1999 and 2004). Validation included the following:

- Verification of 100% of all quality control (QC) sample results (both qualitative and quantitative);
- Verification of the identification of 100% of all sample results (both positive hits and non-detects);
- Recalculation of 10% of all investigative sample results; and

- Preparation of a Data Usability Summary Report (DUSR).

The quality of the data has been assessed and is documented in the DUSR provided in [Appendix C](#). In summary, the results of the data usability assessment show that the collected analytical data for soil, groundwater and soil gas are valid for the intended purposes of the RI.

## SECTION 4

### SITE CHARACTERIZATION RESULTS

This section presents the results of the Site Characterization. Analytical results for the soil and groundwater samples collected during the Site Characterization have been summarized in [Tables 3](#) and [4](#) and on [Figures 4, 5, 6](#) and [7](#).

#### **4.1 SITE GEOLOGY**

The geology encountered in the soil borings during the Site Characterization is summarized in the logs provided in [Appendix A](#). The boring logs show that the upper 5 to 20 feet contained fill materials (generally sand, gravel and cobble with trace amounts of brick, concrete, wood and silt). Deposits of fine to coarse-grained sand with a clay layer underneath the sand were encountered beneath the fill. Clay was encountered from approximately 1 to 6 feet in thickness. Peat was also encountered at a thickness of less than two feet and a depth of 18 ft bgs in the SB-1 boring. Bedrock was not encountered during the Site Characterization activities. Soil boring logs generated during the Site Characterization were used to develop the representative cross section A to A' of the Site as shown on [Figure 7](#).

#### **4.2 FORMER GAS WORKS STRUCTURES**

Remnants of former gas works structures were not encountered within any soil boring or monitoring well installed during the Site Characterization.

#### **4.3 SITE HYDROGEOLOGY**

The depth to groundwater was gauged in the three new monitoring wells (MW-1 through MW-3) on November 27, 2013. Groundwater was encountered beneath the Site at 5.6 to 12.05 feet below ground surface and at elevations ranging from 1.56 feet above MSL at MW-3 to 5.09 feet above MSL at MW-2. The groundwater levels and corresponding elevations are summarized in [Table 2](#). Due to the straight line location of the three monitoring wells, neither groundwater contours nor groundwater flow direction could be calculated. However, it is anticipated that the groundwater flow direction is toward the Hunts Point promontory, and the confluence of the Bronx and East Rivers ([Hygienetics, 1997b](#)).

#### **4.4 SOIL SAMPLE RESULTS**

A total of 30 soil samples, which includes 2 duplicate, 2 matrix spike, and 2 matrix spike duplicate samples, were collected from the soil borings and monitoring well borings as part of the Site Characterization. Soil samples were submitted to Chemtech Laboratories and analyzed for TCL VOCs, TCL SVOCs, TAL metals, and cyanide as described in Section 3. The analytical results of the soil samples are summarized in [Table 3](#) and presented on [Figures 4, 5, and 6](#). The soil sample results have been compared to the Unrestricted Soil Cleanup Objectives (USCOs) provided by NYSDEC in 6 NYCRR Part 375 (NYSDEC, 2006). The USCOs assume there are no imposed restrictions on the use of the Site; however, the Site functions solely as a public sidewalk and right-of-way infrastructure. Therefore, a comparison of soil sample results to the

USCOs is conservative. PID readings, visual observation, and analytical results from the subsurface soil investigation are summarized below.

#### PID Readings/NAPL/Hydrocarbon Fingerprinting Results

PID readings for soil samples collected during soil boring/monitoring well installations ranged from 0.0 to 1,260 ppm above background. The highest PID reading of 1,260 ppm was observed in soil boring SB-4 at a depth interval of 15 to 17 ft bgs. PID readings in the remaining eleven (11) soil borings ranged from 0.0 to 10.7 ppm. Non-aqueous-phase-liquid (NAPL) was observed in one soil boring, SB-4 at approximately 11-18 feet bgs, during the Site Characterization activities.

A sample of soil containing NAPL from soil boring SB-4 was collected and submitted to META Environmental, Inc. of Watertown, MA for forensic hydrocarbon fingerprint analysis. The fingerprinting sample was analyzed by GC/FID (EPA 8100M) for fingerprinting and by GC/MS/SIM (EPA 8270M) for mono- and polycyclic aromatic hydrocarbons (MAHs and PAHs), alkyl PAH homologues and other selected compounds. The laboratory report for this fingerprint analysis is provided in [Appendix D](#). The report indicates that the sample from SB-4 (15-17') contained pyrogenic material and exhibited fluoranthene to pyrene ratios indicative of tars formed from MGPs utilizing carbureted water gas (CWG) processes.

#### VOCs

Sixteen (16) individual VOCs were detected at least once in the soil samples collected during the Site Characterization. Of these, seven (7) (acetone, methylene chloride, benzene, ethylbenzene, toluene, m/p-xylene, and o-xylene) were detected at concentrations exceeding the USCOs. Acetone, which is considered a laboratory contaminant, was detected above its USCO in four soil samples [SB-6 (10-15'), SB-7 (8-10'), SB-8 (15-16'), and SB-9(10-15')]. The remaining six VOCs: methylene chloride, benzene, ethylbenzene, toluene, m/p-xylene, and o-xylene were detected above their respective USCOs only in one soil sample [SB-4 (15-17')].

Total VOC concentrations in all soil samples ranged from non-detect to 1,277 milligrams/kilogram (mg/kg), which was detected in soil sample SB-4 collected at a depth of 15 to 17 ft bgs. Total VOC concentrations in the remaining twenty three (23) samples ranged from non-detect to 0.195 mg/kg.

The vertical extent of VOC impacts was delineated at each soil boring and monitoring well boring sample location with the exception of acetone, which is a common laboratory contaminant.

#### SVOCs

Twenty nine (29) individual SVOCs were detected in soil samples collected during the Site Characterization. Fourteen (14) PAHs [acenaphthene, anthracene, benzo(a)anthracene, benzo(a)pyrene, benzo(b)fluoranthene, benzo(k)fluoranthene, chrysene, dibenz(a,h)anthracene, fluoranthene, flourene, indeno(1,2,3-cd)pyrene, naphthalene, phenanthrene, and pyrene] were detected at concentrations exceeding the USCOs in at least one soil sample. Total SVOC concentrations ranged from 0.44 to 4,635 mg/kg, which was detected in soil sample SB-4 at a depth of 15 to 17 ft bgs. Total SVOC concentrations in the remaining twenty three (23) samples ranged from 0.44 to 238 mg/kg. With the exception of SB-4, PAH concentrations detected during the Halleck Street SC were below the 95<sup>th</sup> percentile of background soil concentrations,

detected within urban fill soils, encountered within street excavation soils located within New York City (RETEC, 2007).

SVOCs exceeding USCOs were detected at nine (9) soil borings: SB-2, SB-3, SB-4, SB-5, SB-6, SB-7, SB-8, SB-9, and SB-10 and the vertical extent of SVOC impacts was delineated at seven (7) of these soil borings: MW-1, SB-1, SB-3, SB-4, SB-5, SB-8, and SB-11 (i.e. no USCOs were exceeded in the deepest sample collected).

### **Inorganics**

A total of 24 inorganic constituents were detected in soil samples collected during the Site Characterization. Eleven of these exceeded the USCOs (arsenic, barium, cadmium, copper, lead, mercury, nickel, selenium, silver, zinc, and cyanide). The vertical extent of impacts from inorganics was delineated at SB-1.

## **4.5 GROUNDWATER SAMPLE RESULTS**

A total of three (3) groundwater samples and 1 duplicate/MS/MSD were collected during the Site Characterization and analyzed for TCL VOCs, TCL SVOCs, TAL Metals, and total cyanide. Laboratory analytical results for constituents detected in the groundwater samples are summarized in [Table 4](#). For evaluation purposes, analytical results were compared with ambient water quality standards (AWQS) and guidance values contained in NYSDEC Technical and Operational Guidance Series (TOGS) 1.1.1 (NYSDEC, 1998). These standards and guidance values are protective of groundwater quality assuming that groundwater is used as a source of drinking water. That assumption is not applicable to the Site because groundwater is not used now, nor will it be used in the future as a source of drinking water. Accordingly, the use of Class GA standards and guidance values for comparison to Site groundwater data is conservative. The analytical results of the groundwater samples collected from each well are presented on [Figure 6](#). Field measurements and observations as well as analytical results from the groundwater investigation are summarized below.

### **Field Measurements**

Each monitoring well was sampled upon reaching parameter stability and turbidity levels below 50 NTU. During groundwater sampling activities, each monitoring well was monitored for the presence of NAPL. No NAPL or sheens were noted in any of the wells. Visual descriptions and observations made during the groundwater sampling activities are presented on the groundwater sampling records provided in [Appendix B](#).

### **VOCs**

Only one VOC was detected in the groundwater samples collected during the Site Characterization. Methyl tert-butyl ether, which is not a MGP-related compound, was detected in two samples at concentrations below its guidance value. Groundwater analytical results for VOCs are summarized in [Table 4](#).

### **SVOCs**

Only one SVOC was detected in the groundwater samples collected during the Site Characterization. Dimethyl phthalate, which is considered a laboratory contaminant, was detected in one sample at a concentration below its guidance value. Groundwater analytical results for SVOCs are summarized in [Table 4](#).

## **Inorganics**

Seventeen (17) inorganic compounds (aluminum, arsenic, barium, cadmium, calcium, total chromium, iron, lead, magnesium, manganese, nickel, potassium, selenium, sodium, thallium, zinc, and cyanide) were detected at least once in the groundwater samples collected during the Site Characterization. Of these, eight (barium, iron, lead, magnesium, manganese, sodium, thallium, and cyanide) were detected at concentrations exceeding the AWQS and guidance values. Groundwater analytical results for inorganics are summarized in [Table 4](#) and on [Figure 7](#).

## **SECTION 5**

### **EXPOSURE ASSESSMENT**

Information collected during the Site Characterization at the former Hunts Point Gas Works Halleck Street Sidewalk Area Site has been used to qualitatively assess potential exposure pathways for the various detected compounds in Site soils and groundwater. The Site is approximately 2,500 linear feet long, located within a commercial district, and includes the street and sidewalk area of Halleck Street.

Analytical results from the soil samples collected during the Site Characterization activities indicate the presence of possible MGP-related contaminants in the soil, primarily at the SB-4 soil boring sample location where MGP-related NAPL was detected. VOCs, PAHs, and inorganic constituents were detected at concentrations above the USCOs at the Site in soil ranging from 5 to 20 ft bgs.

Shallow impacted soils on the Site may be encountered during intrusive maintenance activities (e.g., repair of underground utilities); however, it is unlikely that these materials would be encountered during day-to-day Site operations. It should be noted that the Site soils are covered by asphalt and concrete surfaces.

Groundwater analytical results indicated the presence of inorganic concentrations in the monitoring wells at the Site above the AWQS and guidance values. None of the three monitoring wells (MW-1, MW-2, MW-3) exceeded AWQS for possible MGP-related VOCs or SVOCs.

Groundwater at the Site is currently not used for a potable water source and there are no plans for future use of potable or commercial/industrial groundwater at the Site. Groundwater flow direction could not be determined. The depth to groundwater at the Site is approximately 5.6 to 12 feet bgs. Therefore, there is limited potential for exposure to groundwater during intrusive subsurface activities (e.g., repair of underground utilities) at the Site. Surface water and sediment are not present at the Site.

## **SECTION 6**

### **CONCLUSIONS**

The following conclusions have been made based on the results of the Site Characterization presented herein:

- Remnants of the former MGP structures were not encountered during the Site Characterization.
- NAPL was encountered and sampled at only one soil boring location, SB-4. Forensic hydrocarbon fingerprint analysis indicates that the sample contained materials indicative of tars formed from MGP processes. Soil boring SB-4 is located immediately to the west of Parcel A, OU-2, which is being addressed as part of a separate investigation and remediation. Areas to the west of soil boring SB-4 are covered by sidewalks and an active roadway and NAPL extents have been delineated to the north and south.
- Soil samples collected during the Site Characterization activities indicate the presence of possibly MGP-related VOCs, PAHs and inorganic constituents in the Site's soil at a depth ranging from 5 feet to 20 ft bgs. Elevated levels of these constituents were detected primarily at soil boring SB-4, where NAPL was observed.
- Possibly MGP-related VOCs: methylene chloride, benzene, ethylbenzene, toluene, m/p-xylene, and o-xylene were detected above their respective USCOs only in only one soil sample [SB-4 (15-17')].
- With the exception of SB-4, PAH concentrations detected in soil samples were below the 95<sup>th</sup> percentile of background soil concentrations, detected within urban fill soils, encountered within street excavation soils located within New York City (RETEC, 2007).
- Although metals such as lead, mercury, and sodium were found in Site's soil at levels exceeding USCOs, these constituents are commonly found in fill materials in urban settings. It should be noted, based on historical reports, many areas of the Hunts Point peninsula were subject to historical filling, raising, and land grading to reach their existing heights and configuration.
- In comparing exceedances of inorganics of USCOs with Commercial Restricted Use Soil Cleanup Criteria for the Protection of Public Health, only three constituents, barium (at SB-6), copper (at SB-6 and SB-7), and arsenic (at SB-11) exceeded the commercial criteria.
- No VOCs or SVOCs related to former MGP activities were detected in groundwater.
- Although eight inorganic constituents (barium, iron, lead, magnesium, manganese, sodium, thallium, and cyanide) were detected in groundwater at concentrations exceeding the AWQS and guidance values, it should be noted that groundwater at the Site is not used as a source of drinking water.

## **SECTION 7**

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## **TABLES**

**Table 1**  
**Sample Summary**  
**Former Hunts Point Gas Works**  
**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

Location	Sample ID	Depth (bgs)	TCL VOCs	TCL SVOCs	TAL Metals	Cyanide	Hydrocarbon Fingerprint
<b>SOIL SAMPLES</b>							
MW-1	MW-1(5-10)-20131104	5-10'	X	X	X	X	
	MW-11(5-10)-20131104*	5-10'	X	X	X	X	
	MW-1(10-15)-20131104	10-15'	X	X	X	X	
SB-1	SB-1(10-12)-20131101	10-12'	X	X	X	X	
	SB-1(15-17)-20131101	15-17'	X	X	X	X	
SB-2	SB-2(5-10)-20131101	5-10'	X	X	X	X	
	SB-2(10-15)-20131101	10-15'	X	X	X	X	
SB-3	SB-3(5-10)-20131101	5-10'	X	X	X	X	
	SB-3(12-14)-20131101	12-14'	X	X	X	X	
SB-4	SB-4(15-17)-20131031	15-17'	X	X	X	X	
	SB-4(19-20)-20131031	19-20'	X	X	X	X	
	SB-4(15-17)NAPL	15-17'					X
SB-5	SB-5(5-10)-20131031	5-10'	X	X	X	X	
	SB-5(10-15)-20131031	10-15'	X	X	X	X	
SB-6	SB-6(8-10)-20131031	8-10'	X	X	X	X	
	SB-6(10-15)-20131031	10-15'	X	X	X	X	
SB-7	SB-7(8-10)-20131031	8-10'	X	X	X	X	
	SB-17(8-10)-20131031*	8-10'	X	X	X	X	
	SB-7(16-18)-20131031	16-18'	X	X	X	X	
SB-8	SB-8(15-16)-20131030	15-16'	X	X	X	X	
	SB-8(17-19)-20131030	17-19'	X	X	X	X	
SB-9	SB-9(5-10)-20131030	5-10'	X	X	X	X	
	SB-9(10-15)-20131030	10-15'	X	X	X	X	
SB-10	SB-10(5-10)-20131030	5-10'	X	X	X	X	
	SB-10(10-12)-20131030	10-12'	X	X	X	X	
SB-11	SB-11(5-10)-20131030	5-10'	X	X	X	X	
	SB-11(15-20)-20131030	15-20'	X	X	X	X	
<b>GROUNDWATER SAMPLES</b>							
MW-1	MW-1-20131127	NA	X	X	X	X	
MW-2	MW-2-20131127	NA	X	X	X	X	
MW-3	MW-3-20131127	NA	X	X	X	X	
	MW-3A-20131127*	NA	X	X	X	X	

X - Indicates sample was analyzed

\* - Indicates a duplicate sample.

**Table 2**  
**Summary of Groundwater Elevations**  
**Former Hunts Point Gas Works**  
**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

Monitoring Well Number	Total Well Depth (feet)	Top of Casing Elevation (feet AMSL)	Depth to Water (feet) <sup>(1)</sup>	Groundwater Elevation (feet AMSL)
MW-1	17.15	15.64	12.05	3.59
MW-2	16.10	14.56	9.47	5.09
MW-3	17.25	7.16	5.60	1.56

Notes:

(1) Measured from top of PVC casing on November 27, 2013

AMSL = Above Mean Sea Level

Elevations are based on the North American Vertical Datum of 1988 (NAVD88).

**Table 3**  
**Summary of Soil Analytical Data**  
**Former Hunts Point Gas Works**  
**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	MW-1 MW-1(5-10)-20131104	MW-1 MW-11(5-10)-20131104	MW-1 MW-1(10-15)-20131104
CAS NO.	COMPOUND		E4340-01 5 - 10 ft CTECH E4340 SOIL 11/4/2013 10:50	E4340-05 5 - 10 ft CTECH E4340 SOIL 11/4/2013 10:40	E4340-02 10 - 15 ft CTECH E4340 SOIL 11/4/2013 11:00
	<b>VOLATILES</b>				
67-64-1	ACETONE	0.05	mg/kg	0.0077 J	ND
75-15-0	CARBON DISULFIDE	NS	mg/kg	ND	ND
98-82-8	ISOPROPYLBENZENE (CUMENE)	NS	mg/kg	ND	ND
108-10-1	METHYL ISOBUTYL KETONE	NS	mg/kg	ND	ND
108-87-2	METHYLCYCLOHEXANE	NS	mg/kg	ND	ND
75-09-2	METHYLENE CHLORIDE	0.05	mg/kg	0.0068 J	ND
100-42-5	STYRENE	NS	mg/kg	ND	ND
1634-04-4	TERT-BUTYL METHYL ETHER	0.93	mg/kg	ND	ND
87-61-6	1,2,3-TRICHLOROBENZENE	NS	mg/kg	ND	ND
120-82-1	1,2,4-TRICHLOROBENZENE	NS	mg/kg	ND	ND
106-46-7	1,4-DICHLOROBENZENE	1.8	mg/kg	ND	ND
	<b>BTEX</b>				
71-43-2	BENZENE	0.06	mg/kg	ND	ND
100-41-4	ETHYLBENZENE	1	mg/kg	ND	ND
108-88-3	TOLUENE	0.7	mg/kg	ND	ND
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	0.26	mg/kg	ND	ND
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	0.26	mg/kg	ND	ND
	<b>SEMOVOLATILES</b>				
100-52-7	BENZALDEHYDE	NS	mg/kg	R	R
85-68-7	BENZYL BUTYL PHTHALATE	NS	mg/kg	ND	ND
92-52-4	BIPHENYL (DIPHENYL)	NS	mg/kg	ND	ND
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	NS	mg/kg	ND	ND
86-74-8	CARBAZOLE	NS	mg/kg	ND	ND
132-64-9	DIBENZOFURAN	NS	mg/kg	ND	ND
84-66-2	DIETHYL PHTHALATE	NS	mg/kg	ND	ND
131-11-3	DIMETHYL PHTHALATE	NS	mg/kg	0.61	0.8
105-67-9	2,4-DIMETHYLPHENOL	NS	mg/kg	ND	ND
95-48-7	2-METHYLPHENOL (O-CRESOL)	0.33	mg/kg	ND	ND
MEPH3MEPH4	3- AND 4- METHYLPHENOL (TOTAL)	0.33	mg/kg	ND	ND
108-95-2	PHENOL	0.33	mg/kg	ND	ND

**Table 3**  
**Summary of Soil Analytical Data**  
**Former Hunts Point Gas Works**  
**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

						Dup of MW-1(5-10)-20131104
Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Unrestricted Use Soil Cleanup Objectives	Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	MW-1 MW-11(5-10)-20131104 E4340-01 5 - 10 ft CTECH E4340 SOIL 11/4/2013 10:50	MW-1 MW-11(5-10)-20131104 E4340-05 5 - 10 ft CTECH E4340 SOIL 11/4/2013 10:40	MW-1 MW-1(10-15)-20131104 E4340-02 10 - 15 ft CTECH E4340 SOIL 11/4/2013 11:00
CAS NO.	COMPOUND		UNITS:	2/4/2014	2/4/2014	2/4/2014
	PAHs					
83-32-9	ACENAPHTHENE	20	mg/kg	ND	0.0905 J	ND
208-96-8	ACENAPHTHYLENE	100	mg/kg	0.24 J	0.26 J	ND
120-12-7	ANTHRACENE	100	mg/kg	0.19 J	0.24 J	0.12 J
56-55-3	BENZO(A)ANTHRACENE	1	mg/kg	0.36 J	0.56	0.34 J
50-32-8	BENZO(A)PYRENE	1	mg/kg	0.44 J	0.63	0.29 J
205-99-2	BENZO(B)FLUORANTHENE	1	mg/kg	0.48 J	0.72	0.36 J
191-24-2	BENZO(G,H,I)PERYLENE	100	mg/kg	0.3 J	0.42 J	0.19 J
207-08-9	BENZO(K)FLUORANTHENE	0.8	mg/kg	0.15 J	0.24 J	0.11 J
218-01-9	CHRYSENE	1	mg/kg	0.4 J	0.63	0.26 J
53-70-3	DIBENZ(A,H)ANTHRACENE	0.33	mg/kg	ND	0.11 J	ND
206-44-0	FLUORANTHENE	100	mg/kg	0.65	0.99	0.52
86-73-7	FLUORENE	30	mg/kg	0.1 J	0.12 J	ND
193-39-5	INDENO(1,2,3-C,D)PYRENE	0.5	mg/kg	0.21 J	0.33 J	0.17 J
91-57-6	2-METHYLNAPHTHALENE	NS	mg/kg	ND	0.14 J	ND
91-20-3	NAPHTHALENE	12	mg/kg	0.23 J	0.3 J	ND
85-01-8	PHENANTHRENE	100	mg/kg	0.45 J	0.62	0.32 J
129-00-0	PYRENE	100	mg/kg	0.69	1	0.54
	INORGANICS					
7429-90-5	ALUMINUM	NS	mg/kg	14900 J	13800 J	13000 J
7440-36-0	ANTIMONY	NS	mg/kg	ND	ND	ND
7440-38-2	ARSENIC	13	mg/kg	10.8	6.94	8.22
7440-39-3	BARIUM	350	mg/kg	134 J	108 J	80.4 J
7440-41-7	BERYLLIUM	7.2	mg/kg	0.435	0.361	0.52
7440-43-9	CADMIUM	2.5	mg/kg	2.6	2.18	ND
7440-70-2	CALCIUM	NS	mg/kg	4200 J	4200 J	4300 J
7440-47-3	CHROMIUM, TOTAL	NS	mg/kg	36.1 J	30.1 J	28.3 J
7440-48-4	COBALT	NS	mg/kg	12.4 J	10.47 J	14.1 J
7440-50-8	COPPER	50	mg/kg	99.3 J	53.8 J	49.1 J
7439-89-6	IRON	NS	mg/kg	36900	33900	34700
7439-92-1	LEAD	63	mg/kg	493	352	390
7439-95-4	MAGNESIUM	NS	mg/kg	5500 J	5000 J	6800 J
7439-96-5	MANGANESE	1600	mg/kg	323 J	270 J	641 J
7439-97-6	MERCURY	0.18	mg/kg	2.33	1.54	0.268
7440-02-0	NICKEL	30	mg/kg	29.4	25.3	29.6
7440-09-7	POTASSIUM	NS	mg/kg	2900 J	2700 J	3300 J
7782-49-2	SELENIUM	3.9	mg/kg	2.56	2.62	2.74
7440-22-4	SILVER	2	mg/kg	1.43	1.18	1.82
7440-23-5	SODIUM	NS	mg/kg	271 J	245 J	607 J
7440-28-0	THALLIUM	NS	mg/kg	ND	ND	ND
7440-62-2	VANADIUM	NS	mg/kg	50.5 J	44.9 J	50.5 J
7440-66-6	ZINC	109	mg/kg	346	329	227
57-12-5	CYANIDE	27	mg/kg	0.649	0.672	0.244 J

Notes:

(1) 6NYCRR Part 375 Environmental Remediation Programs (December 14, 2006)

(2) NS indicates no cleanup objective or background level is available.

(3) ND indicates compound was not detected.

(4) J indicates an estimated concentration.

(5) Shaded values exceed 6NYCRR Part 375 Unrestricted Use Soil Cleanup Objectives.

(6) NA indicates compound was not analyzed.

(7) R indicates rejected value

**Table 3**  
**Summary of Soil Analytical Data**  
**Former Hunts Point Gas Works**  
**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-1 SB-1(10-12)-20131101 E4277-25 10 - 12 ft CTECH E4277 SOIL 11/1/2013 13:45	SB-1 SB-1(15-17)-20131101 E4277-26 15 - 17 ft CTECH E4277 SOIL 11/1/2013 13:55	SB-2 SB-2(5-10)-20131101 E4277-21 5 - 10 ft CTECH E4277 SOIL 11/1/2013 9:45
CAS NO.	COMPOUND		UNITS:	2/4/2014	2/4/2014
	<b>VOLATILES</b>				
67-64-1	ACETONE	0.05	mg/kg	ND	ND
75-15-0	CARBON DISULFIDE	NS	mg/kg	ND	ND
98-82-8	ISOPROPYLBENZENE (CUMENE)	NS	mg/kg	ND	ND
108-10-1	METHYL ISOBUTYL KETONE	NS	mg/kg	ND	ND
108-87-2	METHYLCYCLOHEXANE	NS	mg/kg	ND	ND
75-09-2	METHYLENE CHLORIDE	0.05	mg/kg	0.0037 J	0.0052 J
100-42-5	STYRENE	NS	mg/kg	ND	ND
1634-04-4	TERT-BUTYL METHYL ETHER	0.93	mg/kg	ND	ND
87-61-6	1,2,3-TRICHLOROBENZENE	NS	mg/kg	ND	ND
120-82-1	1,2,4-TRICHLOROBENZENE	NS	mg/kg	ND	ND
106-46-7	1,4-DICHLOROBENZENE	1.8	mg/kg	ND	ND
	<b>BTEX</b>				
71-43-2	BENZENE	0.06	mg/kg	ND	ND
100-41-4	ETHYLBENZENE	1	mg/kg	ND	ND
108-88-3	TOLUENE	0.7	mg/kg	ND	ND
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	0.26	mg/kg	ND	ND
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	0.26	mg/kg	ND	ND
	<b>SEMIVOLATILES</b>				
100-52-7	BENZALDEHYDE	NS	mg/kg	R	R
85-68-7	BENZYL BUTYL PHTHALATE	NS	mg/kg	ND	0.31 J
92-52-4	BIPHENYL (DIPHENYL)	NS	mg/kg	ND	ND
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	NS	mg/kg	ND	ND
86-74-8	CARBAZOLE	NS	mg/kg	ND	0.0928 J
132-64-9	DIBENZOFURAN	NS	mg/kg	ND	ND
84-66-2	DIETHYL PHTHALATE	NS	mg/kg	ND	ND
131-11-3	DIMETHYL PHTHALATE	NS	mg/kg	0.48	0.44 J
105-67-9	2,4-DIMETHYLPHENOL	NS	mg/kg	ND	ND
95-48-7	2-METHYLPHENOL (O-CRESOL)	0.33	mg/kg	ND	ND
MEPH3MEPH4	3- AND 4- METHYLPHENOL (TOTAL)	0.33	mg/kg	ND	ND
108-95-2	PHENOL	0.33	mg/kg	ND	ND

**Table 3**  
**Summary of Soil Analytical Data**  
**Former Hunts Point Gas Works**  
**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-1 SB-1(10-12)-20131101 E4277-25 10 - 12 ft CTECH E4277 SOIL 11/1/2013 13:45	SB-1 SB-1(15-17)-20131101 E4277-26 15 - 17 ft CTECH E4277 SOIL 11/1/2013 13:55	SB-2 SB-2(5-10)-20131101 E4277-21 5 - 10 ft CTECH E4277 SOIL 11/1/2013 9:45
CAS NO.	COMPOUND		UNITS:	2/4/2014	2/4/2014
	PAHs				
83-32-9	ACENAPHTHENE	20	mg/kg	ND	ND
208-96-8	ACENAPHTHYLENE	100	mg/kg	ND	0.19 J
120-12-7	ANTHRACENE	100	mg/kg	ND	0.38 J
56-55-3	BENZO(A)ANTHRACENE	1	mg/kg	ND	1.4
50-32-8	BENZO(A)PYRENE	1	mg/kg	ND	1.6
205-99-2	BENZO(B)FLUORANTHENE	1	mg/kg	ND	1.8
191-24-2	BENZO(G,H,I)PERYLENE	100	mg/kg	ND	1.2
207-08-9	BENZO(K)FLUORANTHENE	0.8	mg/kg	ND	0.84
218-01-9	CHRYSENE	1	mg/kg	ND	1.4
53-70-3	DIBENZ(A,H)ANTHRACENE	0.33	mg/kg	ND	0.25 J
206-44-0	FLUORANTHENE	100	mg/kg	ND	2
86-73-7	FLUORENE	30	mg/kg	ND	0.13 J
193-39-5	INDENO(1,2,3-C,D)PYRENE	0.5	mg/kg	ND	0.97
91-57-6	2-METHYLNAPHTHALENE	NS	mg/kg	ND	ND
91-20-3	NAPHTHALENE	12	mg/kg	ND	0.29 J
85-01-8	PHENANTHRENE	100	mg/kg	ND	1.2
129-00-0	PYRENE	100	mg/kg	ND	1.8
	INORGANICS				
7429-90-5	ALUMINUM	NS	mg/kg	12000	3100
7440-36-0	ANTIMONY	NS	mg/kg	ND	2.67 J
7440-38-2	ARSENIC	13	mg/kg	1.48	10.67
7440-39-3	BARIUM	350	mg/kg	71	416
7440-41-7	BERYLLIUM	7.2	mg/kg	0.365	0.087 J
7440-43-9	CADMIUM	2.5	mg/kg	ND	1.33
7440-70-2	CALCIUM	NS	mg/kg	8100	13500
7440-47-3	CHROMIUM, TOTAL	NS	mg/kg	25.1	30.7
7440-48-4	COBALT	NS	mg/kg	13.2	6.78
7440-50-8	COPPER	50	mg/kg	21.1	595
7439-89-6	IRON	NS	mg/kg	22600	31000
7439-92-1	LEAD	63	mg/kg	17.3 J	727 J
7439-95-4	MAGNESIUM	NS	mg/kg	5600	1800
7439-96-5	MANGANESE	1600	mg/kg	440	229
7439-97-6	MERCURY	0.18	mg/kg	0.039	0.67
7440-02-0	NICKEL	30	mg/kg	21.6 J	17.8 J
7440-09-7	POTASSIUM	NS	mg/kg	2000 J	906 J
7782-49-2	SELENIUM	3.9	mg/kg	1.36	2.96
7440-22-4	SILVER	2	mg/kg	1.3	2.21
7440-23-5	SODIUM	NS	mg/kg	116	649
7440-28-0	THALLIUM	NS	mg/kg	ND	ND
7440-62-2	VANADIUM	NS	mg/kg	40	23.2
7440-66-6	ZINC	109	mg/kg	58.5	17.7
57-12-5	CYANIDE	27	mg/kg	0.093 J	830
				ND	1.86

Notes:

(1) 6NYCRR Part 375 Environmental Remediation Programs (December 14, 2006)

(2) NS indicates no cleanup objective or background level is available.

(3) ND indicates compound was not detected.

(4) J indicates an estimated concentration.

(5) Shaded values exceed 6NYCRR Part 375 Unrestricted Use Soil Cleanup Objectives.

(6) NA indicates compound was not analyzed.

(7) R indicates rejected value

**Table 3**  
**Summary of Soil Analytical Data**  
**Former Hunts Point Gas Works**  
**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Unrestricted Use Soil Cleanup Objectives	Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-2 SB-2(10-15)-20131101 E4277-22 10 - 15 ft CTECH E4277 SOIL 11/1/2013 9:55	SB-3 SB-3(5-10)-20131101 E4277-23 5 - 10 ft CTECH E4277 SOIL 11/1/2013 11:15	SB-3 SB-3(12-14)-20131101 E4277-24 12 - 14 ft CTECH E4277 SOIL 11/1/2013 11:25
CAS NO.	COMPOUND		UNITS:	2/4/2014	2/4/2014	2/4/2014
	<b>VOLATILES</b>					
67-64-1	ACETONE	0.05	mg/kg	ND	ND	ND
75-15-0	CARBON DISULFIDE	NS	mg/kg	ND	ND	ND
98-82-8	ISOPROPYLBENZENE (CUMENE)	NS	mg/kg	ND	ND	ND
108-10-1	METHYL ISOBUTYL KETONE	NS	mg/kg	ND	ND	ND
108-87-2	METHYLCYCLOHEXANE	NS	mg/kg	ND	ND	ND
75-09-2	METHYLENE CHLORIDE	0.05	mg/kg	0.006 J	0.0063 J	0.0069 J
100-42-5	STYRENE	NS	mg/kg	ND	ND	ND
1634-04-4	TERT-BUTYL METHYL ETHER	0.93	mg/kg	ND	ND	ND
87-61-6	1,2,3-TRICHLOROBENZENE	NS	mg/kg	ND	ND	ND
120-82-1	1,2,4-TRICHLOROBENZENE	NS	mg/kg	ND	ND	ND
106-46-7	1,4-DICHLOROBENZENE	1.8	mg/kg	ND	ND	ND
	<b>BTEX</b>					
71-43-2	BENZENE	0.06	mg/kg	ND	ND	ND
100-41-4	ETHYLBENZENE	1	mg/kg	ND	ND	ND
108-88-3	TOLUENE	0.7	mg/kg	ND	ND	ND
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	0.26	mg/kg	ND	ND	ND
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	0.26	mg/kg	ND	ND	ND
	<b>SEMIVOLATILES</b>					
100-52-7	BENZALDEHYDE	NS	mg/kg	R	R	R
85-68-7	BENZYL BUTYL PHTHALATE	NS	mg/kg	ND	ND	ND
92-52-4	BIPHENYL (DIPHENYL)	NS	mg/kg	ND	ND	ND
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	NS	mg/kg	ND	ND	ND
86-74-8	CARBAZOLE	NS	mg/kg	0.29 J	ND	ND
132-64-9	DIBENZOFURAN	NS	mg/kg	0.35 J	ND	ND
84-66-2	DIETHYL PHTHALATE	NS	mg/kg	ND	ND	ND
131-11-3	DIMETHYL PHTHALATE	NS	mg/kg	0.89	ND	1
105-67-9	2,4-DIMETHYLPHENOL	NS	mg/kg	ND	ND	ND
95-48-7	2-METHYLPHENOL (O-CRESOL)	0.33	mg/kg	ND	ND	ND
MEPH3MEPH4	3- AND 4- METHYLPHENOL (TOTAL)	0.33	mg/kg	ND	ND	0.26 J
108-95-2	PHENOL	0.33	mg/kg	ND	ND	ND

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Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-2	SB-3	SB-3
CAS NO.	COMPOUND		SB-2(10-15)-20131101 E4277-22 10 - 15 ft CTECH E4277 SOIL 11/1/2013 9:55 2/4/2014	SB-3(5-10)-20131101 E4277-23 5 - 10 ft CTECH E4277 SOIL 11/1/2013 11:15 2/4/2014	SB-3(12-14)-20131101 E4277-24 12 - 14 ft CTECH E4277 SOIL 11/1/2013 11:25 2/4/2014
	PAHs				
83-32-9	ACENAPHTHENENE	20	mg/kg	0.36 J	ND
208-96-8	ACENAPHTHYLENE	100	mg/kg	0.23 J	ND
120-12-7	ANTHRACENE	100	mg/kg	1.5	0.97 J
56-55-3	BENZO(A)ANTHRACENE	1	mg/kg	3.6	6.3
50-32-8	BENZO(A)PYRENE	1	mg/kg	3.2	7.2
205-99-2	BENZO(B)FLUORANTHENENE	1	mg/kg	3.8	8.7
191-24-2	BENZO(G,H,I)PERYLENE	100	mg/kg	1.8	5.1
207-08-9	BENZO(K)FLUORANTHENENE	0.8	mg/kg	1.1	3.3 J
218-01-9	CHRYSENE	1	mg/kg	2.4	5.9
53-70-3	DIBENZ(A,H)ANTHRACENE	0.33	mg/kg	0.69	1.2 J
206-44-0	FLUORANTHENENE	100	mg/kg	7	9.9
86-73-7	FLUORENE	30	mg/kg	0.74	ND
193-39-5	INDENO(1,2,3-C,D)PYRENE	0.5	mg/kg	1.7	4.5
91-57-6	2-METHYLNAPHTHALENE	NS	mg/kg	ND	ND
91-20-3	NAPHTHALENE	12	mg/kg	0.2 J	ND
85-01-8	PHENANTHRENE	100	mg/kg	5	4.3
129-00-0	PYRENE	100	mg/kg	5.5	7
	INORGANICS				
7429-90-5	ALUMINUM	NS	mg/kg	11100	3900
7440-36-0	ANTIMONY	NS	mg/kg	ND	9.52 J
7440-38-2	ARSENIC	13	mg/kg	10.43	27.6
7440-39-3	BARIUM	350	mg/kg	168	354
7440-41-7	BERYLLIUM	7.2	mg/kg	0.426	ND
7440-43-9	CADMIUM	2.5	mg/kg	ND	2.73
7440-70-2	CALCIUM	NS	mg/kg	5900	38300
7440-47-3	CHROMIUM, TOTAL	NS	mg/kg	25.9	24.9
7440-48-4	COBALT	NS	mg/kg	12.3	9.71
7440-50-8	COPPER	50	mg/kg	35.3	361
7439-89-6	IRON	NS	mg/kg	32900	68200
7439-92-1	LEAD	63	mg/kg	168 J	1500 J
7439-95-4	MAGNESIUM	NS	mg/kg	6100	3800
7439-96-5	MANGANESE	1600	mg/kg	695	331
7439-97-6	MERCURY	0.18	mg/kg	0.765	3.85
7440-02-0	NICKEL	30	mg/kg	26.7 J	35.7 J
7440-09-7	POTASSIUM	NS	mg/kg	2700 J	476 J
7782-49-2	SELENIUM	3.9	mg/kg	2.22	3.95
7440-22-4	SILVER	2	mg/kg	2.29	4.94
7440-23-5	SODIUM	NS	mg/kg	1100	658
7440-28-0	THALLIUM	NS	mg/kg	ND	ND
7440-62-2	VANADIUM	NS	mg/kg	43.9	23.4
7440-66-6	ZINC	109	mg/kg	187	1100
57-12-5	CYANIDE	27	mg/kg	0.158 J	2.94

Notes:

(1) 6NYCRR Part 375 Environmental Remediation Programs (December 14, 2006)

(2) NS indicates no cleanup objective or background level is available.

(3) ND indicates compound was not detected.

(4) J indicates an estimated concentration.

(5) Shaded values exceed 6NYCRR Part 375 Unrestricted Use Soil Cleanup Objectives.

(6) NA indicates compound was not analyzed.

(7) R indicates rejected value

**Table 3**  
**Summary of Soil Analytical Data**  
**Former Hunts Point Gas Works**  
**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Unrestricted Use Soil Cleanup Objectives	Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-4 SB-4(15-17)-20131031 E4277-19 15 - 17 ft CTECH E4277 SOIL 10/31/2013 14:15 2/4/2014	SB-4 SB-4(19-20)-20131031 E4277-20 19 - 20 ft CTECH E4277 SOIL 10/31/2013 14:15 2/4/2014	SB-5 SB-5(5-10)-20131031 E4277-17 5 - 10 ft CTECH E4277 SOIL 10/31/2013 12:40 2/4/2014
CAS NO.	COMPOUND		UNITS:			
	VOLATILES					
67-64-1	ACETONE	0.05	mg/kg	ND	0.0408 J	ND
75-15-0	CARBON DISULFIDE	NS	mg/kg	ND	0.008 J	ND
98-82-8	ISOPROPYLBENZENE (CUMENE)	NS	mg/kg	29 J	0.0065 J	ND
108-10-1	METHYL ISOBUTYL KETONE	NS	mg/kg	ND	ND	ND
108-87-2	METHYLCYCLOHEXANE	NS	mg/kg	0.66 J	ND	ND
75-09-2	METHYLENE CHLORIDE	0.05	mg/kg	1.7 J	0.0175	ND
100-42-5	STYRENE	NS	mg/kg	ND	ND	ND
1634-04-4	TERT-BUTYL METHYL ETHER	0.93	mg/kg	ND	0.0106 J	ND
87-61-6	1,2,3-TRICHLOROBENZENE	NS	mg/kg	9.9	ND	ND
120-82-1	1,2,4-TRICHLOROBENZENE	NS	mg/kg	63.2 J	ND	ND
106-46-7	1,4-DICHLOROBENZENE	1.8	mg/kg	ND	0.0025 J	ND
	BTEX					
71-43-2	BENZENE	0.06	mg/kg	85.3	0.0384	ND
100-41-4	ETHYLBENZENE	1	mg/kg	518.3 J	0.0379	ND
108-88-3	TOLUENE	0.7	mg/kg	18.4	ND	ND
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	0.26	mg/kg	438.1	0.0146 J	ND
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	0.26	mg/kg	112.9	0.016	ND
	SEMICVOLATILES					
100-52-7	BENZALDEHYDE	NS	mg/kg	R	R	R
85-68-7	BENZYL BUTYL PHTHALATE	NS	mg/kg	ND	ND	ND
92-52-4	BIPHENYL (DIPHENYL)	NS	mg/kg	74.5	ND	ND
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	NS	mg/kg	ND	ND	ND
86-74-8	CARBAZOLE	NS	mg/kg	3.6 J	ND	1.4
132-64-9	DIBENZOFURAN	NS	mg/kg	23.2	ND	0.44
84-66-2	DIETHYL PHTHALATE	NS	mg/kg	ND	ND	ND
131-11-3	DIMETHYL PHTHALATE	NS	mg/kg	ND	0.6	0.66
105-67-9	2,4-DIMETHYLPHENOL	NS	mg/kg	ND	ND	ND
95-48-7	2-METHYLPHENOL (O-CRESOL)	0.33	mg/kg	ND	ND	ND
MEPH3MEPH4	3- AND 4- METHYLPHENOL (TOTAL)	0.33	mg/kg	ND	ND	ND
108-95-2	PHENOL	0.33	mg/kg	ND	ND	ND

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Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Unrestricted Use Soil Cleanup Objectives	SB-4 SB-4(15-17)-20131031 E4277-19 15 - 17 ft CTECH E4277 SOIL 10/31/2013 14:15 2/4/2014	SB-4 SB-4(19-20)-20131031 E4277-20 19 - 20 ft CTECH E4277 SOIL 10/31/2013 14:15 2/4/2014	SB-5 SB-5(5-10)-20131031 E4277-17 5 - 10 ft CTECH E4277 SOIL 10/31/2013 12:40 2/4/2014
CAS NO.	COMPOUND	Matrix: Sampled: Validated: UNITS:			
	PAHs				
83-32-9	ACENAPHTHENE	20	mg/kg 233.2	ND	ND
208-96-8	ACENAPHTHYLENE	100	mg/kg 33.3	ND	2.4
120-12-7	ANTHRACENE	100	mg/kg 133.6	0.13 J	5.4
56-55-3	BENZO(A)ANTHRACENE	1	mg/kg 80.2	0.23 J	8.2
50-32-8	BENZO(A)PYRENE	1	mg/kg 72.2	0.21 J	7.3
205-99-2	BENZO(B)FLUORANTHENE	1	mg/kg 59.1	0.24 J	7.7
191-24-2	BENZO(G,H,I)PERYLENE	100	mg/kg 29.2	ND	3.1
207-08-9	BENZO(K)FLUORANTHENE	0.8	mg/kg 14.1	ND	3.8
218-01-9	CHRYSENE	1	mg/kg 73.4	0.18 J	7.3
53-70-3	DIBENZA(H)ANTHRACENE	0.33	mg/kg 6.9	ND	0.87
206-44-0	FLUORANTHENE	100	mg/kg 145.4	0.41 J	14.8
86-73-7	FLUORENE	30	mg/kg 165	ND	1.1
193-39-5	INDENO(1,2,3-C,D)PYRENE	0.5	mg/kg 21.6	ND	3.2
91-57-6	2-METHYLNAPHTHALENE	NS	mg/kg 766.4	0.16 J	0.3 J
91-20-3	NAPHTHALENE	12	mg/kg 1899.2	0.48 J	1.2
85-01-8	PHENANTHRENE	100	mg/kg 547.3	0.38 J	12.3
129-00-0	PYRENE	100	mg/kg 253.6	0.37 J	13
	INORGANICS				
7429-90-5	ALUMINUM	NS	mg/kg 4000	15500	2700
7440-36-0	ANTIMONY	NS	mg/kg 1.74 J	ND	ND
7440-38-2	ARSENIC	13	mg/kg 12	10.52	5.42
7440-39-3	BARIUM	350	mg/kg 78.3	54.5	170
7440-41-7	BERYLLIUM	7.2	mg/kg 0.291 J	0.643	0.204 J
7440-43-9	CADMUM	2.5	mg/kg 0.3 J	ND	ND
7440-70-2	CALCIUM	NS	mg/kg 2000	3700	3200
7440-47-3	CHROMIUM, TOTAL	NS	mg/kg 16.7	35	24.8
7440-48-4	COBALT	NS	mg/kg 9.46	17.2	5.01
7440-50-8	COPPER	50	mg/kg 92.1	28.2	129
7439-89-6	IRON	NS	mg/kg 46400	40100	13000
7439-92-1	LEAD	63	mg/kg 292 J	63.9 J	737 J
7439-95-4	MAGNESIUM	NS	mg/kg 374	8700	1200
7439-96-5	MANGANESE	1600	mg/kg 113	808	49.9
7439-97-6	MERCURY	0.18	mg/kg 3.11	0.461	0.199
7440-02-0	NICKEL	30	mg/kg 21.6 J	37.2 J	11.2 J
7440-09-7	POTASSIUM	NS	mg/kg 597 J	4100 J	584 J
7782-49-2	SELENIUM	3.9	mg/kg 3.67	2.49	1.95
7440-22-4	SILVER	2	mg/kg 2.97	2.81	0.862
7440-23-5	SODIUM	NS	mg/kg 430	3400	299
7440-28-0	THALLIUM	NS	mg/kg ND	ND	ND
7440-62-2	VANADIUM	NS	mg/kg 19.5	58.1	24.5
7440-66-6	ZINC	109	mg/kg 495	219	72
57-12-5	CYANIDE	27	mg/kg 34.9	1.84	0.79

Notes:

(1) 6NYCRR Part 375 Environmental Remediation Programs (December 14, 2006)

(2) NS indicates no cleanup objective or background level is available.

(3) ND indicates compound was not detected.

(4) J indicates an estimated concentration.

(5) Shaded values exceed 6NYCRR Part 375 Unrestricted Use Soil Cleanup Objectives.

(6) NA indicates compound was not analyzed.

(7) R indicates rejected value

**Table 3**  
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**Former Hunts Point Gas Works**  
**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Unrestricted Use Soil Cleanup Objectives	Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-5 SB-5(10-15)-20131031 E4277-18 10 - 15 ft CTECH E4277 SOIL 10/31/2013 12:50 2/4/2014	SB-6 SB-6(8-10)-20131031 E4277-12 8 - 10 ft CTECH E4277 SOIL 10/31/2013 10:20 2/4/2014	SB-6 SB-6(10-15)-20131031 E4277-15 10 - 15 ft CTECH E4277 SOIL 10/31/2013 10:30 2/4/2014
CAS NO.	COMPOUND		UNITS:			
	<b>VOLATILES</b>					
67-64-1	ACETONE	0.05	mg/kg	ND	ND	0.0637
75-15-0	CARBON DISULFIDE	NS	mg/kg	ND	ND	ND
98-82-8	ISOPROPYLBENZENE (CUMENE)	NS	mg/kg	ND	ND	ND
108-10-1	METHYL ISOBUTYL KETONE	NS	mg/kg	ND	ND	ND
108-87-2	METHYLCYCLOHEXANE	NS	mg/kg	ND	ND	ND
75-09-2	METHYLENE CHLORIDE	0.05	mg/kg	0.004 J	0.0078 J	ND
100-42-5	STYRENE	NS	mg/kg	ND	ND	ND
1634-04-4	TERT-BUTYL METHYL ETHER	0.93	mg/kg	ND	ND	ND
87-61-6	1,2,3-TRICHLOROBENZENE	NS	mg/kg	ND	ND	ND
120-82-1	1,2,4-TRICHLOROBENZENE	NS	mg/kg	ND	ND	ND
106-46-7	1,4-DICHLOROBENZENE	1.8	mg/kg	ND	ND	ND
	<b>BTEX</b>					
71-43-2	BENZENE	0.06	mg/kg	ND	ND	ND
100-41-4	ETHYLBENZENE	1	mg/kg	ND	ND	ND
108-88-3	TOLUENE	0.7	mg/kg	ND	ND	ND
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	0.26	mg/kg	ND	ND	ND
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	0.26	mg/kg	ND	ND	ND
	<b>SEMIVOLATILES</b>					
100-52-7	BENZALDEHYDE	NS	mg/kg	R	R	R
85-68-7	BENZYL BUTYL PHTHALATE	NS	mg/kg	ND	ND	ND
92-52-4	BIPHENYL (DIPHENYL)	NS	mg/kg	ND	ND	ND
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	NS	mg/kg	ND	ND	ND
86-74-8	CARBAZOLE	NS	mg/kg	ND	ND	0.12 J
132-64-9	DIBENZOFURAN	NS	mg/kg	ND	ND	0.15 J
84-66-2	DIETHYL PHTHALATE	NS	mg/kg	ND	ND	0.15 J
131-11-3	DIMETHYL PHTHALATE	NS	mg/kg	0.31 J	0.45	0.68
105-67-9	2,4-DIMETHYLPHENOL	NS	mg/kg	ND	ND	ND
95-48-7	2-METHYLPHENOL (O-CRESOL)	0.33	mg/kg	ND	ND	ND
MEPH3MEPH4	3- AND 4- METHYLPHENOL (TOTAL)	0.33	mg/kg	ND	ND	ND
108-95-2	PHENOL	0.33	mg/kg	ND	ND	ND

**Table 3**  
**Summary of Soil Analytical Data**  
**Former Hunts Point Gas Works**  
**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-5 SB-5(10-15)-20131031 E4277-18 10 - 15 ft CTECH E4277 SOIL 10/31/2013 12:50	SB-6 SB-6(8-10)-20131031 E4277-12 8 - 10 ft CTECH E4277 SOIL 10/31/2013 10:20	SB-6 SB-6(10-15)-20131031 E4277-15 10 - 15 ft CTECH E4277 SOIL 10/31/2013 10:30
CAS NO.	COMPOUND		UNITS:	2/4/2014	2/4/2014
	PAHs				
83-32-9	ACENAPHTHENE	20	mg/kg	ND	ND
208-96-8	ACENAPHTHYLENE	100	mg/kg	ND	0.34 J
120-12-7	ANTHRACENE	100	mg/kg	ND	0.32 J
56-55-3	BENZO(A)ANTHRACENE	1	mg/kg	0.13 J	0.75
50-32-8	BENZO(A)PYRENE	1	mg/kg	0.12 J	0.73
205-99-2	BENZO(B)FLUORANTHENE	1	mg/kg	0.15 J	0.79
191-24-2	BENZO(G,H,I)PERYLENE	100	mg/kg	ND	0.55
207-08-9	BENZO(K)FLUORANTHENE	0.8	mg/kg	ND	0.38 J
218-01-9	CHRYSENE	1	mg/kg	0.11 J	0.71
53-70-3	DIBENZ(A,H)ANTHRACENE	0.33	mg/kg	ND	0.12 J
206-44-0	FLUORANTHENE	100	mg/kg	0.21 J	1.4
86-73-7	FLUORENE	30	mg/kg	ND	0.25 J
193-39-5	INDENO(1,2,3-C,D)PYRENE	0.5	mg/kg	ND	0.45
91-57-6	2-METHYLNAPHTHALENE	NS	mg/kg	ND	0.17 J
91-20-3	NAPHTHALENE	12	mg/kg	0.19 J	0.18 J
85-01-8	PHENANTHRENE	100	mg/kg	0.17 J	1.2
129-00-0	PYRENE	100	mg/kg	0.2 J	1.5
	INORGANICS				
7429-90-5	ALUMINUM	NS	mg/kg	6300	7300
7440-36-0	ANTIMONY	NS	mg/kg	ND	1.36 J
7440-38-2	ARSENIC	13	mg/kg	7.29	15.9
7440-39-3	BARIUM	350	mg/kg	377	203
7440-41-7	BERYLLIUM	7.2	mg/kg	0.474	0.34
7440-43-9	CADMIUM	2.5	mg/kg	0.619	0.404
7440-70-2	CALCIUM	NS	mg/kg	5000	15000
7440-47-3	CHROMIUM, TOTAL	NS	mg/kg	20	62.1
7440-48-4	COBALT	NS	mg/kg	7.22	7.31
7440-50-8	COPPER	50	mg/kg	120	79.6
7439-89-6	IRON	NS	mg/kg	18200	34300
7439-92-1	LEAD	63	mg/kg	494 J	227 J
7439-95-4	MAGNESIUM	NS	mg/kg	2300	3200
7439-96-5	MANGANESE	1600	mg/kg	128	165
7439-97-6	MERCURY	0.18	mg/kg	0.319	0.39
7440-02-0	NICKEL	30	mg/kg	18.5 J	55.1 J
7440-09-7	POTASSIUM	NS	mg/kg	1200 J	1100 J
7782-49-2	SELENIUM	3.9	mg/kg	1.94	2.5
7440-22-4	SILVER	2	mg/kg	1.31	2.41
7440-23-5	SODIUM	NS	mg/kg	1600	1300
7440-28-0	THALLIUM	NS	mg/kg	ND	ND
7440-62-2	VANADIUM	NS	mg/kg	29	41.2
7440-66-6	ZINC	109	mg/kg	259	336
57-12-5	CYANIDE	27	mg/kg	4.13	7.32

Notes:

(1) 6NYCRR Part 375 Environmental Remediation Programs (December 14, 2006)

(2) NS indicates no cleanup objective or background level is available.

(3) ND indicates compound was not detected.

(4) J indicates an estimated concentration.

(5) Shaded values exceed 6NYCRR Part 375 Unrestricted Use Soil Cleanup Objectives.

(6) NA indicates compound was not analyzed.

(7) R indicates rejected value

**Table 3**  
**Summary of Soil Analytical Data**  
**Former Hunts Point Gas Works**  
**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-7	SB-7	SB-7
CAS NO.	COMPOUND		SB-7(8-10)-20131031 E4277-10 8 - 10 ft CTECH E4277 SOIL 10/31/2013 8:35 2/4/2014	SB-17(8-10)-20131031 E4277-16 8 - 10 ft CTECH E4277 SOIL 10/31/2013 8:45 2/4/2014	SB-17(8-10)-20131031 E4277-11 16 - 18 ft CTECH E4277 SOIL 10/31/2013 10:20 2/4/2014
	<b>VOLATILES</b>				
67-64-1	ACETONE	0.05	mg/kg	0.0337	0.0616
75-15-0	CARBON DISULFIDE	NS	mg/kg	ND	ND
98-82-8	ISOPROPYLBENZENE (CUMENE)	NS	mg/kg	ND	ND
108-10-1	METHYL ISOBUTYL KETONE	NS	mg/kg	ND	ND
108-87-2	METHYLCYCLOHEXANE	NS	mg/kg	ND	ND
75-09-2	METHYLENE CHLORIDE	0.05	mg/kg	0.0044 J	ND
100-42-5	STYRENE	NS	mg/kg	ND	ND
1634-04-4	TERT-BUTYL METHYL ETHER	0.93	mg/kg	ND	ND
87-61-6	1,2,3-TRICHLOROBENZENE	NS	mg/kg	ND	ND
120-82-1	1,2,4-TRICHLOROBENZENE	NS	mg/kg	ND	ND
106-46-7	1,4-DICHLOROBENZENE	1.8	mg/kg	ND	ND
	<b>BTEX</b>				
71-43-2	BENZENE	0.06	mg/kg	ND	ND
100-41-4	ETHYLBENZENE	1	mg/kg	ND	ND
108-88-3	TOLUENE	0.7	mg/kg	ND	ND
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	0.26	mg/kg	ND	ND
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	0.26	mg/kg	ND	ND
	<b>SEMIVOLATILES</b>				
100-52-7	BENZALDEHYDE	NS	mg/kg	R	R
85-68-7	BENZYL BUTYL PHTHALATE	NS	mg/kg	ND	ND
92-52-4	BIPHENYL (DIPHENYL)	NS	mg/kg	ND	ND
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	NS	mg/kg	ND	0.11 J
86-74-8	CARBAZOLE	NS	mg/kg	0.0801 J	0.13 J
132-64-9	DIBENZOFURAN	NS	mg/kg	ND	0.0847 J
84-66-2	DIETHYL PHTHALATE	NS	mg/kg	ND	ND
131-11-3	DIMETHYL PHTHALATE	NS	mg/kg	0.24 J	0.61 J
105-67-9	2,4-DIMETHYLPHENOL	NS	mg/kg	ND	ND
95-48-7	2-METHYLPHENOL (O-CRESOL)	0.33	mg/kg	ND	ND
MEPH3MEPH4	3- AND 4- METHYLPHENOL (TOTAL)	0.33	mg/kg	ND	ND
108-95-2	PHENOL	0.33	mg/kg	ND	ND

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**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

						Dup of SB-7(8-10)-20131031
Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-7	SB-7	SB-7	SB-7
CAS NO.	COMPOUND		SB-7(8-10)-20131031	E4277-10	E4277-16	SB-7(16-18)-20131031
	PAHs	Unrestricted Use Soil Cleanup Objectives	UNITS:			
83-32-9	ACENAPHTHENE	20	mg/kg	ND	ND	ND
208-96-8	ACENAPHTHYLENE	100	mg/kg	0.23 J	0.32 J	ND
120-12-7	ANTHRACENE	100	mg/kg	0.33 J	0.58 J	0.2 J
56-55-3	BENZO(A)ANTHRACENE	1	mg/kg	1.2	1.7	0.99
50-32-8	BENZO(A)PYRENE	1	mg/kg	1.2	1.9	0.89
205-99-2	BENZO(B)FLUORANTHENE	1	mg/kg	1.5	2	1.1
191-24-2	BENZO(G,H,I)PERYLENE	100	mg/kg	0.89 J	1.6 J	0.53
207-08-9	BENZO(K)FLUORANTHENE	0.8	mg/kg	0.36 J	0.9 J	0.39 J
218-01-9	CHRYSENE	1	mg/kg	1.2	1.5	0.89
53-70-3	DIBENZ(A,H)ANTHRACENE	0.33	mg/kg	0.21 J	0.25 J	0.14 J
206-44-0	FLUORANTHENE	100	mg/kg	1.7	2.7	1.3
86-73-7	FLUORENE	30	mg/kg	0.15 J	0.25 J	ND
193-39-5	INDENO(1,2,3-C,D)PYRENE	0.5	mg/kg	0.83	1.3	0.49
91-57-6	2-METHYLNAPHTHALENE	NS	mg/kg	0.0869 J	0.15 J	ND
91-20-3	NAPHTHALENE	12	mg/kg	0.13 J	0.27 J	ND
85-01-8	PHENANTHRENE	100	mg/kg	1	1.7	0.51
129-00-0	PYRENE	100	mg/kg	1.8	2.8	1.4
	INORGANICS					
7429-90-5	ALUMINUM	NS	mg/kg	6800	7500	2400
7440-36-0	ANTIMONY	NS	mg/kg	0.76 J	1.11 J	3.91 J
7440-38-2	ARSENIC	13	mg/kg	4.36	5.98	12.8
7440-39-3	BARIUM	350	mg/kg	195	201	240
7440-41-7	BERYLLIUM	7.2	mg/kg	0.317	0.36	0.097 J
7440-43-9	CADMIUM	2.5	mg/kg	2.36	2.58	5.08
7440-70-2	CALCIUM	NS	mg/kg	6400	9800	14200
7440-47-3	CHROMIUM, TOTAL	NS	mg/kg	20.6	21.9	17
7440-48-4	COBALT	NS	mg/kg	9.55	10.81	6.35
7440-50-8	COPPER	50	mg/kg	120	126	414
7439-89-6	IRON	NS	mg/kg	22400	28900	52900
7439-92-1	LEAD	63	mg/kg	301 J	283 J	719 J
7439-95-4	MAGNESIUM	NS	mg/kg	2800	5100	1400
7439-96-5	MANGANESE	1600	mg/kg	237	342	336
7439-97-6	MERCURY	0.18	mg/kg	0.361	0.454	1.12
7440-02-0	NICKEL	30	mg/kg	29.4 J	37.5 J	212 J
7440-09-7	POTASSIUM	NS	mg/kg	2300 J	2400 J	419 J
7782-49-2	SELENIUM	3.9	mg/kg	1.1	0.856 J	2.62
7440-22-4	SILVER	2	mg/kg	0.661	0.99	1.86
7440-23-5	SODIUM	NS	mg/kg	242	245	684
7440-28-0	THALLIUM	NS	mg/kg	ND	ND	ND
7440-62-2	VANADIUM	NS	mg/kg	41.8	50.9	20.8
7440-66-6	ZINC	109	mg/kg	315	372	719
57-12-5	CYANIDE	27	mg/kg	ND	1.3 J	1.78

Notes:

(1) 6NYCRR Part 375 Environmental Remediation Programs (December 14, 2006)

(2) NS indicates no cleanup objective or background level is available.

(3) ND indicates compound was not detected.

(4) J indicates an estimated concentration.

(5) Shaded values exceed 6NYCRR Part 375 Unrestricted Use Soil Cleanup Objectives.

(6) NA indicates compound was not analyzed.

(7) R indicates rejected value

**Table 3**  
**Summary of Soil Analytical Data**  
**Former Hunts Point Gas Works**  
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**Site Characterization - October and November 2013**

Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Unrestricted Use Soil Cleanup Objectives	Location ID:	SB-8	SB-8	SB-9
CAS NO.	COMPOUND		Sample ID:	SB-8(15-16)-20131030	E4277-08	SB-9(5-10)-20131030
			Lab Sample Id:	E4277-07	17 - 19 ft	E4277-05
		Depth:	15 - 16 ft		5 - 10 ft	
		Source:	CTECH		CTECH	
		SDG:	E4277		E4277	
		Matrix:	SOIL		SOIL	
		Sampled:	10/30/2013 14:55		10/30/2013 14:55	
		Validated:	2/4/2014		2/4/2014	
		UNITS:				
	VOLATILES					
67-64-1	ACETONE	0.05	mg/kg	0.0525	ND	0.0236 J
75-15-0	CARBON DISULFIDE	NS	mg/kg	0.0023 J	0.003 J	0.0035 J
98-82-8	ISOPROPYLBENZENE (CUMENE)	NS	mg/kg	0.0025 J	ND	ND
108-10-1	METHYL ISOBUTYL KETONE	NS	mg/kg	0.0061 J	ND	0.0054 J
108-87-2	METHYLCYCLOHEXANE	NS	mg/kg	ND	ND	0.0014 J
75-09-2	METHYLENE CHLORIDE	0.05	mg/kg	0.0025 J	0.0022 J	0.0035 J
100-42-5	STYRENE	NS	mg/kg	0.0037 J	ND	0.0094
1634-04-4	TERT-BUTYL METHYL ETHER	0.93	mg/kg	ND	ND	ND
87-61-6	1,2,3-TRICHLOROBENZENE	NS	mg/kg	ND	ND	ND
120-82-1	1,2,4-TRICHLOROBENZENE	NS	mg/kg	ND	ND	ND
106-46-7	1,4-DICHLOROBENZENE	1.8	mg/kg	ND	ND	ND
	BTEX					
71-43-2	BENZENE	0.06	mg/kg	0.0137	ND	0.0097
100-41-4	ETHYLBENZENE	1	mg/kg	0.0045 J	ND	0.004 J
108-88-3	TOLUENE	0.7	mg/kg	0.0093	ND	0.0213
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	0.26	mg/kg	0.0266	ND	0.0768
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	0.26	mg/kg	0.013	ND	0.0367
	SEMIVOLATILES					
100-52-7	BENZALDEHYDE	NS	mg/kg	R	R	R
85-68-7	BENZYL BUTYL PHTHALATE	NS	mg/kg	ND	ND	ND
92-52-4	BIPHENYL (DIPHENYL)	NS	mg/kg	0.63	ND	0.4
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	NS	mg/kg	ND	ND	ND
86-74-8	CARBAZOLE	NS	mg/kg	0.97 J	ND	4.1 J
132-64-9	DIBENZOFURAN	NS	mg/kg	1.4	ND	2.2
84-66-2	DIETHYL PHTHALATE	NS	mg/kg	ND	ND	ND
131-11-3	DIMETHYL PHTHALATE	NS	mg/kg	0.72	0.61	0.41
105-67-9	2,4-DIMETHYLPHENOL	NS	mg/kg	ND	ND	0.11 J
95-48-7	2-METHYLPHENOL (O-CRESOL)	0.33	mg/kg	ND	ND	0.12 J
MEPH3MEPH4	3- AND 4- METHYLPHENOL (TOTAL)	0.33	mg/kg	0.14 J	ND	0.33 J
108-95-2	PHENOL	0.33	mg/kg	0.14 J	ND	0.24 J

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**Site Characterization - October and November 2013**

Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-8	SB-8	SB-9
CAS NO.	COMPOUND		SB-8(15-16)-20131030 E4277-07 15 - 16 ft CTECH E4277 SOIL 10/30/2013 14:55 2/4/2014	SB-8(17-19)-20131030 E4277-08 17 - 19 ft CTECH E4277 SOIL 10/30/2013 14:55 2/4/2014	SB-8(17-19)-20131030 E4277-05 5 - 10 ft CTECH E4277 SOIL 10/30/2013 13:30 2/4/2014
	PAHs				
83-32-9	ACENAPHTHENE	20	mg/kg	0.7	ND
208-96-8	ACENAPHTHYLENE	100	mg/kg	2.5	ND
120-12-7	ANTHRACENE	100	mg/kg	3.9	ND
56-55-3	BENZO(A)ANTHRACENE	1	mg/kg	4.9	ND
50-32-8	BENZO(A)PYRENE	1	mg/kg	3.1	ND
205-99-2	BENZO(B)FLUORANTHENE	1	mg/kg	2.7 J	ND
191-24-2	BENZO(G,H,I)PERYLENE	100	mg/kg	1.9 J	ND
207-08-9	BENZO(K)FLUORANTHENE	0.8	mg/kg	1.9	ND
218-01-9	CHRYSENE	1	mg/kg	2.7	ND
53-70-3	DIBENZ(A,H)ANTHRACENE	0.33	mg/kg	0.76	ND
206-44-0	FLUORANTHENE	100	mg/kg	9.2	ND
86-73-7	FLUORENE	30	mg/kg	3.2	ND
193-39-5	INDENO(1,2,3-C,D)PYRENE	0.5	mg/kg	1.6	ND
91-57-6	2-METHYLNAPHTHALENE	NS	mg/kg	1.9	ND
91-20-3	NAPHTHALENE	12	mg/kg	6.5	ND
85-01-8	PHENANTHRENE	100	mg/kg	15	ND
129-00-0	PYRENE	100	mg/kg	9.7	ND
	INORGANICS				
7429-90-5	ALUMINUM	NS	mg/kg	4100	2000
7440-36-0	ANTIMONY	NS	mg/kg	4.33 J	3.23 J
7440-38-2	ARSENIC	13	mg/kg	13.4	5.75
7440-39-3	BARIUM	350	mg/kg	275	157
7440-41-7	BERYLLIUM	7.2	mg/kg	0.226 J	0.183 J
7440-43-9	CADMIUM	2.5	mg/kg	1.42	ND
7440-70-2	CALCIUM	NS	mg/kg	6500	18900
7440-47-3	CHROMIUM, TOTAL	NS	mg/kg	20.5	19.1
7440-48-4	COBALT	NS	mg/kg	7.93	7.17
7440-50-8	COPPER	50	mg/kg	390	63.6
7439-89-6	IRON	NS	mg/kg	24000	21000
7439-92-1	LEAD	63	mg/kg	884 J	261 J
7439-95-4	MAGNESIUM	NS	mg/kg	2500	3900
7439-96-5	MANGANESE	1600	mg/kg	200	185
7439-97-6	MERCURY	0.18	mg/kg	2.18	0.131
7440-02-0	NICKEL	30	mg/kg	96.2 J	20.1 J
7440-09-7	POTASSIUM	NS	mg/kg	799 J	410 J
7782-49-2	SELENIUM	3.9	mg/kg	2.35	1.66
7440-22-4	SILVER	2	mg/kg	2.08	1.58
7440-23-5	SODIUM	NS	mg/kg	387	460
7440-28-0	THALLIUM	NS	mg/kg	ND	ND
7440-62-2	VANADIUM	NS	mg/kg	24.1	23.1
7440-66-6	ZINC	109	mg/kg	3200	197
57-12-5	CYANIDE	27	mg/kg	2	0.109 J

Notes:

(1) 6NYCRR Part 375 Environmental Remediation Programs (December 14, 2006)

(2) NS indicates no cleanup objective or background level is available.

(3) ND indicates compound was not detected.

(4) J indicates an estimated concentration.

(5) Shaded values exceed 6NYCRR Part 375 Unrestricted Use Soil Cleanup Objectives.

(6) NA indicates compound was not analyzed.

(7) R indicates rejected value

**Table 3**  
**Summary of Soil Analytical Data**  
**Former Hunts Point Gas Works**  
**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Unrestricted Use Soil Cleanup Objectives	Location ID:	SB-9	SB-10	SB-10
CAS NO.	COMPOUND		Sample ID:	SB-9(10-15)-20131030	SB-10(5-10)-20131030	SB-10(10-12)-20131030
			Lab Sample Id:	E4277-06	E4277-03	E4277-04
		Depth:	10 - 15 ft	5 - 10 ft	10 - 12 ft	
		Source:	CTECH	CTECH	CTECH	
		SDG:	E4277	E4277	E4277	
		Matrix:	SOIL	SOIL	SOIL	
		Sampled:	10/30/2013 13:40	10/30/2013 11:20	10/30/2013 11:30	
		Validated:	2/4/2014	2/4/2014	2/4/2014	
		UNITS:				
	<b>VOLATILES</b>					
67-64-1	ACETONE	0.05	mg/kg	0.0814	0.0476	0.018 J
75-15-0	CARBON DISULFIDE	NS	mg/kg	0.0017 J	0.002 J	0.0026 J
98-82-8	ISOPROPYLBENZENE (CUMENE)	NS	mg/kg	ND	ND	ND
108-10-1	METHYL ISOBUTYL KETONE	NS	mg/kg	ND	ND	ND
108-87-2	METHYLCYCLOHEXANE	NS	mg/kg	ND	ND	ND
75-09-2	METHYLENE CHLORIDE	0.05	mg/kg	0.0028 J	0.0048 J	0.005 J
100-42-5	STYRENE	NS	mg/kg	ND	ND	ND
1634-04-4	TERT-BUTYL METHYL ETHER	0.93	mg/kg	ND	ND	ND
87-61-6	1,2,3-TRICHLOROBENZENE	NS	mg/kg	ND	ND	ND
120-82-1	1,2,4-TRICHLOROBENZENE	NS	mg/kg	ND	ND	ND
106-46-7	1,4-DICHLOROBENZENE	1.8	mg/kg	ND	ND	ND
	<b>BTEX</b>					
71-43-2	BENZENE	0.06	mg/kg	ND	ND	ND
100-41-4	ETHYLBENZENE	1	mg/kg	ND	ND	ND
108-88-3	TOLUENE	0.7	mg/kg	ND	ND	ND
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	0.26	mg/kg	ND	ND	ND
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	0.26	mg/kg	ND	ND	ND
	<b>SEMIVOLATILES</b>					
100-52-7	BENZALDEHYDE	NS	mg/kg	R	0.0993 J	R
85-68-7	BENZYL BUTYL PHTHALATE	NS	mg/kg	ND	ND	ND
92-52-4	BIPHENYL (DIPHENYL)	NS	mg/kg	ND	ND	ND
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	NS	mg/kg	ND	ND	ND
86-74-8	CARBAZOLE	NS	mg/kg	0.11 J	ND	0.13 J
132-64-9	DIBENZOFURAN	NS	mg/kg	ND	ND	0.13 J
84-66-2	DIETHYL PHTHALATE	NS	mg/kg	ND	ND	ND
131-11-3	DIMETHYL PHTHALATE	NS	mg/kg	0.59	0.59	0.54
105-67-9	2,4-DIMETHYLPHENOL	NS	mg/kg	ND	ND	ND
95-48-7	2-METHYLPHENOL (O-CRESOL)	0.33	mg/kg	ND	ND	ND
MEPH3MEPH4	3- AND 4- METHYLPHENOL (TOTAL)	0.33	mg/kg	ND	ND	ND
108-95-2	PHENOL	0.33	mg/kg	ND	ND	ND

**Table 3**  
**Summary of Soil Analytical Data**  
**Former Hunts Point Gas Works**  
**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-9 SB-9(10-15)-20131030 E4277-06 10 - 15 ft CTECH E4277 SOIL 10/30/2013 13:40	SB-10 SB-10(5-10)-20131030 E4277-03 5 - 10 ft CTECH E4277 SOIL 10/30/2013 11:20	SB-10 SB-10(10-12)-20131030 E4277-04 10 - 12 ft CTECH E4277 SOIL 10/30/2013 11:30
CAS NO.	COMPOUND		UNITS:	2/4/2014	2/4/2014
	PAHs				
83-32-9	ACENAPHTHENENE	20	mg/kg	ND	0.21 J
208-96-8	ACENAPHTHYLENE	100	mg/kg	0.24 J	0.0942 J
120-12-7	ANTHRACENE	100	mg/kg	0.48	0.92
56-55-3	BENZO(A)ANTHRACENE	1	mg/kg	0.95	1.4
50-32-8	BENZO(A)PYRENE	1	mg/kg	0.77	0.92
205-99-2	BENZO(B)FLUORANTHENENE	1	mg/kg	0.86 J	1.1 J
191-24-2	BENZO(G,H,I)PERYLENE	100	mg/kg	0.76 J	0.5 J
207-08-9	BENZO(K)FLUORANTHENENE	0.8	mg/kg	0.29 J	0.5
218-01-9	CHRYSENE	1	mg/kg	0.75	1.1
53-70-3	DIBENZ(A,H)ANTHRACENE	0.33	mg/kg	0.19 J	0.24 J
206-44-0	FLUORANTHENENE	100	mg/kg	1.7	2.6
86-73-7	FLUORENE	30	mg/kg	0.23 J	0.31 J
193-39-5	INDENO(1,2,3-C,D)PYRENE	0.5	mg/kg	0.55	0.49
91-57-6	2-METHYLNAPHTHALENE	NS	mg/kg	ND	ND
91-20-3	NAPHTHALENE	12	mg/kg	0.2 J	0.15 J
85-01-8	PHENANTHRENE	100	mg/kg	1.2	2.1
129-00-0	PYRENE	100	mg/kg	1.7	2.2
	INORGANICS				
7429-90-5	ALUMINUM	NS	mg/kg	5200	7800
7440-36-0	ANTIMONY	NS	mg/kg	1.15 J	0.732 J
7440-38-2	ARSENIC	13	mg/kg	6.71	4.85
7440-39-3	BARIUM	350	mg/kg	99.1	135
7440-41-7	BERYLLIUM	7.2	mg/kg	0.323	0.25 J
7440-43-9	CADMIUM	2.5	mg/kg	0.862	0.14 J
7440-70-2	CALCIUM	NS	mg/kg	49900	22900
7440-47-3	CHROMIUM, TOTAL	NS	mg/kg	17.2	19.3
7440-48-4	COBALT	NS	mg/kg	9.24	9.87
7440-50-8	COPPER	50	mg/kg	70.3	56.3
7439-89-6	IRON	NS	mg/kg	17000	22200
7439-92-1	LEAD	63	mg/kg	161 J	173 J
7439-95-4	MAGNESIUM	NS	mg/kg	18200	7600
7439-96-5	MANGANESE	1600	mg/kg	176	229
7439-97-6	MERCURY	0.18	mg/kg	0.573	0.303
7440-02-0	NICKEL	30	mg/kg	32.3 J	26.5 J
7440-09-7	POTASSIUM	NS	mg/kg	912 J	1800 J
7782-49-2	SELENIUM	3.9	mg/kg	1.27	2.09
7440-22-4	SILVER	2	mg/kg	1.41	1.3
7440-23-5	SODIUM	NS	mg/kg	258	623
7440-28-0	THALLIUM	NS	mg/kg	ND	ND
7440-62-2	VANADIUM	NS	mg/kg	25.6	34.8
7440-66-6	ZINC	109	mg/kg	246	151
57-12-5	CYANIDE	27	mg/kg	0.104 J	0.522

Notes:

(1) 6NYCRR Part 375 Environmental Remediation Programs (December 14, 2006)

(2) NS indicates no cleanup objective or background level is available.

(3) ND indicates compound was not detected.

(4) J indicates an estimated concentration.

(5) Shaded values exceed 6NYCRR Part 375 Unrestricted Use Soil Cleanup Objectives.

(6) NA indicates compound was not analyzed.

(7) R indicates rejected value

**Table 3**  
**Summary of Soil Analytical Data**  
**Former Hunts Point Gas Works**  
**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Unrestricted Use Soil Cleanup Objectives	Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-11 SB-11(5-10)-20131030 E4277-01 5 - 10 ft CTECH E4277 SOIL 10/30/2013 9:45 2/4/2014	SB-11 SB-11(15-20)-20131030 E4277-02 15 - 20 ft CTECH E4277 SOIL 10/30/2013 10:05 2/4/2014
CAS NO.	COMPOUND		UNITS:		
	VOLATILES				
67-64-1	ACETONE	0.05	mg/kg	0.0295 J	0.0362 J
75-15-0	CARBON DISULFIDE	NS	mg/kg	ND	0.0036 J
98-82-8	ISOPROPYLBENZENE (CUMENE)	NS	mg/kg	ND	ND
108-10-1	METHYL ISOBUTYL KETONE	NS	mg/kg	ND	ND
108-87-2	METHYLCYCLOHEXANE	NS	mg/kg	ND	ND
75-09-2	METHYLENE CHLORIDE	0.05	mg/kg	0.0042 J	0.0056 J
100-42-5	STYRENE	NS	mg/kg	ND	ND
1634-04-4	TERT-BUTYL METHYL ETHER	0.93	mg/kg	ND	ND
87-61-6	1,2,3-TRICHLOROBENZENE	NS	mg/kg	ND	ND
120-82-1	1,2,4-TRICHLOROBENZENE	NS	mg/kg	ND	ND
106-46-7	1,4-DICHLOROBENZENE	1.8	mg/kg	ND	ND
	BTEX				
71-43-2	BENZENE	0.06	mg/kg	ND	ND
100-41-4	ETHYLBENZENE	1	mg/kg	ND	ND
108-88-3	TOLUENE	0.7	mg/kg	ND	ND
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	0.26	mg/kg	ND	ND
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	0.26	mg/kg	ND	ND
	SEMOVATILES				
100-52-7	BENZALDEHYDE	NS	mg/kg	R	R
85-68-7	BENZYL BUTYL PHTHALATE	NS	mg/kg	ND	ND
92-52-4	BIPHENYL (DIPHENYL)	NS	mg/kg	ND	ND
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	NS	mg/kg	0.0886 J	ND
86-74-8	CARBAZOLE	NS	mg/kg	ND	ND
132-64-9	DIBENZOFURAN	NS	mg/kg	ND	ND
84-66-2	DIETHYL PHTHALATE	NS	mg/kg	ND	ND
131-11-3	DIMETHYL PHTHALATE	NS	mg/kg	0.65	0.72
105-67-9	2,4-DIMETHYLPHENOL	NS	mg/kg	ND	ND
95-48-7	2-METHYLPHENOL (O-CRESOL)	0.33	mg/kg	ND	ND
MEPH3MEPH4	3- AND 4- METHYLPHENOL (TOTAL)	0.33	mg/kg	ND	ND
108-95-2	PHENOL	0.33	mg/kg	ND	ND

**Table 3**  
**Summary of Soil Analytical Data**  
**Former Hunts Point Gas Works**  
**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

Con Ed - Halleck Street Validated Soil Analytical Data Detected Compound Summary		Unrestricted Use Soil Cleanup Objectives	Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-11 SB-11(5-10)-20131030 E4277-01 5 - 10 ft CTECH E4277 SOIL 10/30/2013 9:45 2/4/2014	SB-11 SB-11(15-20)-20131030 E4277-02 15 - 20 ft CTECH E4277 SOIL 10/30/2013 10:05 2/4/2014
CAS NO.	COMPOUND		UNITS:		
	PAHs				
83-32-9	ACENAPHTHENE	20	mg/kg	ND	ND
208-96-8	ACENAPHTHYLENE	100	mg/kg	ND	ND
120-12-7	ANTHRACENE	100	mg/kg	0.14 J	0.12 J
56-55-3	BENZO(A)ANTHRACENE	1	mg/kg	0.72	0.24 J
50-32-8	BENZO(A)PYRENE	1	mg/kg	0.66	0.19 J
205-99-2	BENZO(B)FLUORANTHENE	1	mg/kg	0.76 J	0.18 J
191-24-2	BENZO(G,H,I)PERYLENE	100	mg/kg	0.48 J	0.12 J
207-08-9	BENZO(K)FLUORANTHENE	0.8	mg/kg	0.3 J	ND
218-01-9	CHRYSENE	1	mg/kg	0.5	0.17 J
53-70-3	DIBENZ(A,H)ANTHRACENE	0.33	mg/kg	0.19 J	ND
206-44-0	FLUORANTHENE	100	mg/kg	0.82	0.37 J
86-73-7	FLUORENE	30	mg/kg	ND	ND
193-39-5	INDENO(1,2,3-C,D)PYRENE	0.5	mg/kg	0.42	0.0999 J
91-57-6	2-METHYLNAPHTHALENE	NS	mg/kg	ND	ND
91-20-3	NAPHTHALENE	12	mg/kg	ND	ND
85-01-8	PHENANTHRENE	100	mg/kg	0.41	0.38 J
129-00-0	PYRENE	100	mg/kg	0.83	0.42 J
	INORGANICS				
7429-90-5	ALUMINUM	NS	mg/kg	4300	9600
7440-36-0	ANTIMONY	NS	mg/kg	1.5 J	2.26 J
7440-38-2	ARSENIC	13	mg/kg	5.77	20.8
7440-39-3	BARIUM	350	mg/kg	237	207
7440-41-7	BERYLLIUM	7.2	mg/kg	0.348	0.271 J
7440-43-9	CADMUM	2.5	mg/kg	ND	0.839
7440-70-2	CALCIUM	NS	mg/kg	4300	7300
7440-47-3	CHROMIUM, TOTAL	NS	mg/kg	12.6	34
7440-48-4	COBALT	NS	mg/kg	11.7	13.9
7440-50-8	COPPER	50	mg/kg	346	130
7439-89-6	IRON	NS	mg/kg	22300	77700
7439-92-1	LEAD	63	mg/kg	884 J	338 J
7439-95-4	MAGNESIUM	NS	mg/kg	1300	4700
7439-96-5	MANGANESE	1600	mg/kg	128	564
7439-97-6	MERCURY	0.18	mg/kg	0.709	2.67
7440-02-0	NICKEL	30	mg/kg	14.5 J	33.2 J
7440-09-7	POTASSIUM	NS	mg/kg	871 J	2300 J
7782-49-2	SELENIUM	3.9	mg/kg	1.97	3.85
7440-22-4	SILVER	2	mg/kg	1.64	5.38
7440-23-5	SODIUM	NS	mg/kg	508	2300
7440-28-0	THALLIUM	NS	mg/kg	ND	ND
7440-62-2	VANADIUM	NS	mg/kg	19.4	41
7440-66-6	ZINC	109	mg/kg	112	288
57-12-5	CYANIDE	27	mg/kg	8.79	9.13

Notes:

- (1) 6NYCRR Part 375 Environmental Remediation Programs (December 14, 2006)
- (2) NS indicates no cleanup objective or background level is available.
- (3) ND indicates compound was not detected.
- (4) J indicates an estimated concentration.
- (5) Shaded values exceed 6NYCRR Part 375 Unrestricted Use Soil Cleanup Objectives.
- (6) NA indicates compound was not analyzed.
- (7) R indicates rejected value

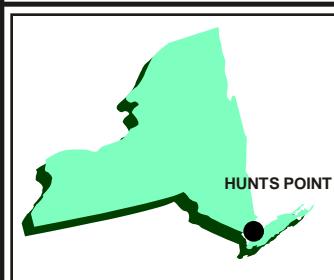
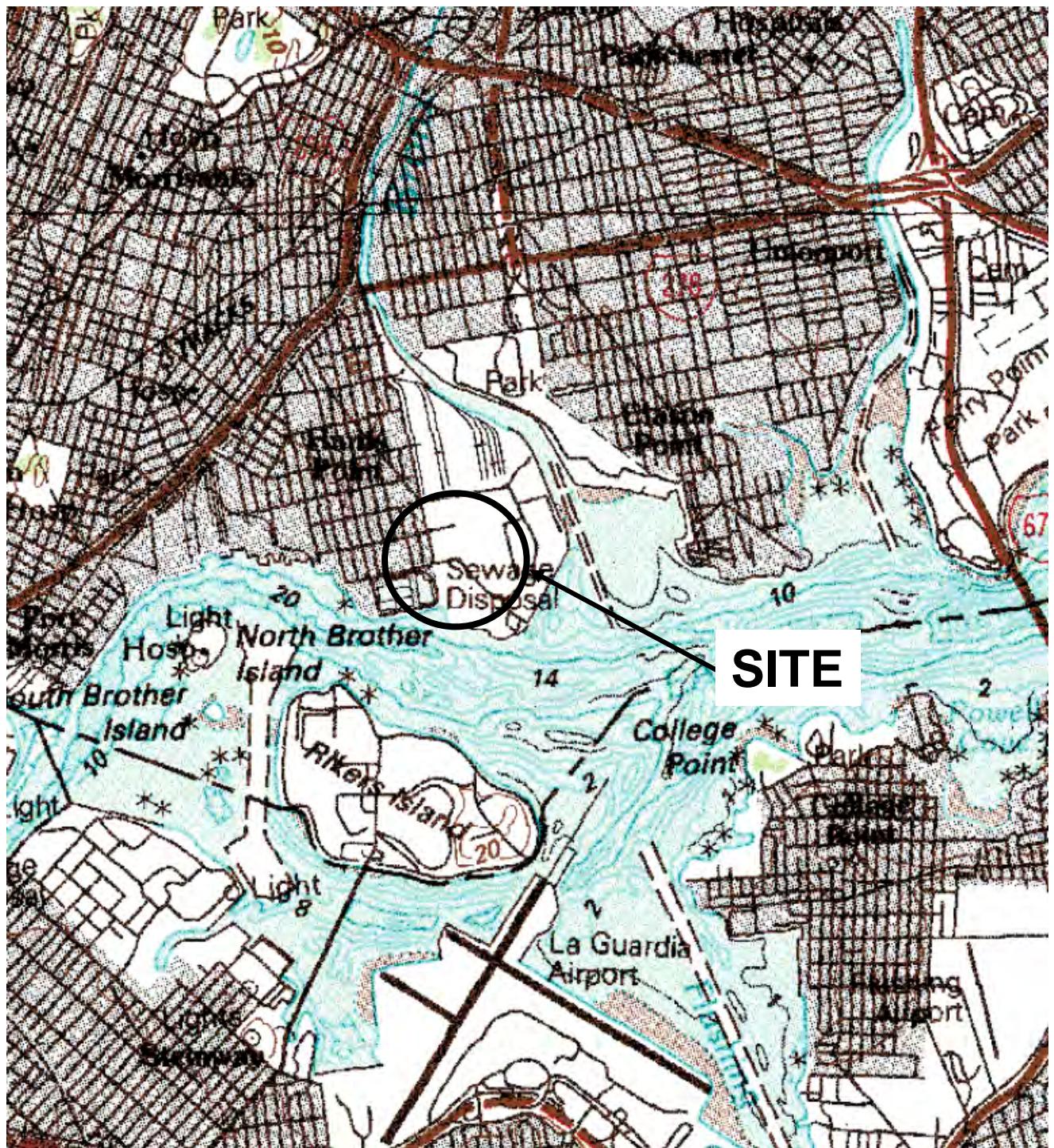
**Table 4**  
**Summary of Groundwater Analytical Data**  
**Former Hunts Point Gas Works**  
**Consolidated Edison Company of New York**  
**Site Characterization - October and November 2013**

							Dup of MW-3-20131127
Con Ed - Halleck Street Validated Groundwater Analytical Data Detected Compound Summary		NYSDEC Ambient Water Quality Standards/Guidance Criteria	Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	MW-1 MW-1-20131127 E4638-06	MW-2 MW-2-20131127 E4638-05	MW-3 MW-3-20131127 E4638-01	MW-3 MW-3A-20131127 E4638-04
CAS NO.	COMPOUND		UNITS:				
1634-04-4	VOLATILES TERT-BUTYL METHYL ETHER	10 (G)	ug/l	ND	6.8	ND	0.56 J
131-11-3	SEMIVOLATILES DIMETHYL PHTHALATE	50 (G)	ug/l	ND	ND	ND	8.6 J
	INORGANICS						
7429-90-5	ALUMINUM	NS	ug/l	920	252	48.2 J	53.6
7440-38-2	ARSENIC	25	ug/l	ND	8.81 J	ND	ND
7440-39-3	BARIUM	1000	ug/l	319	84.3	1400	1400
7440-43-9	CADMIUM	5	ug/l	0.833 J	ND	ND	ND
7440-70-2	CALCIUM	NS	ug/l	148200	142000	175900	170000
7440-47-3	CHROMIUM, TOTAL	50	ug/l	34.8	1.97 J	ND	5.67 J
7439-89-6	IRON	300	ug/l	50500	28100	33500	31300
7439-92-1	LEAD	25	ug/l	3.31 J	25.3	6.64	6.74
7439-95-4	MAGNESIUM	35000 (G)	ug/l	34600	18400	58900	59400
7439-96-5	MANGANESE	300	ug/l	2900	472	719	698
7440-02-0	NICKEL	100	ug/l	13.4 J	ND	ND	ND
7440-09-7	POTASSIUM	NS	ug/l	24100	18500	41000	40000
7782-49-2	SELENIUM	10	ug/l	5.78 J	ND	ND	ND
7440-23-5	SODIUM	20000	ug/l	610600	186000	635100	628900
7440-28-0	THALLIUM	0.5 (G)	ug/l	3.56 J	ND	ND	ND
7440-66-6	ZINC	2000 (G)	ug/l	11.2 J	12.3 J	ND	ND
57-12-5	CYANIDE	200	ug/l	12	28	687	681

Notes:

- Indicates concentration exceeds standard or guidance value.
- (G) Indicates guidance value.
- NS No standard or guidance value available.
- ND Indicates compound was not detected.
- J Indicates an estimated concentration.
- ug/L Micrograms per liter

## **FIGURES**



LATITUDE: N40° 48' 33"  
LONGITUDE: W73° 52' 48"



SOURCE: DeLORME 3-D  
TOPOQUAD PROGRAM

New York  
Quadrangle

## FIGURE 1

CONSOLIDATED EDISON COMPANY OF NEW YORK  
FORMER HUNTS POINT GAS WORKS  
HALLECK STREET SIDEWALK AREA  
BRONX, NEW YORK

## SITE VICINITY MAP

**PARSONS**

200 COTTONTAIL ROAD, SOMERSET NJ 08873 PHONE: (732) 537-3500



#### Legend

- Approximate Site Boundary (Hunts Point MGP OU-3)
- Approximate Boundary of Former Gas Works Facility
- Approximate Boundary of Parcels

**FIGURE 2**

CONSOLIDATED EDISON COMPANY OF NEW YORK  
FORMER HUNTS POINT GAS WORKS  
HALLECK STREET SIDEWALK AREA  
BRONX, NEW YORK

#### **Site Location Map**

**PARSONS**

200 COTTONTAIL LANE SOUTH, SOMERSET, NJ 08873 PHONE: (732) 537-3500



FILE NAME: P:\CONED\448468 HALLECK STREET\CAD\FIGURES\448468-SK005.DWG  
PLOT DATE: 3/5/2014 9:35 AM PLOTTED BY: RUSSO, JILL

## LEGEND:

- SITE BOUNDARY  
MGP FACILITIES  
SOIL BORING LOCATION  
MONITORING WELL  
LOCATION

A horizontal number line starting at -400 and ending at 400. The line is marked with tick marks at intervals of 100, labeled as follows: 200, 100, 0, 200, 400. The labels 200, 100, and 0 are positioned above the line, while the labels 200 and 400 are positioned below the line.

SCALE: 1"=200'

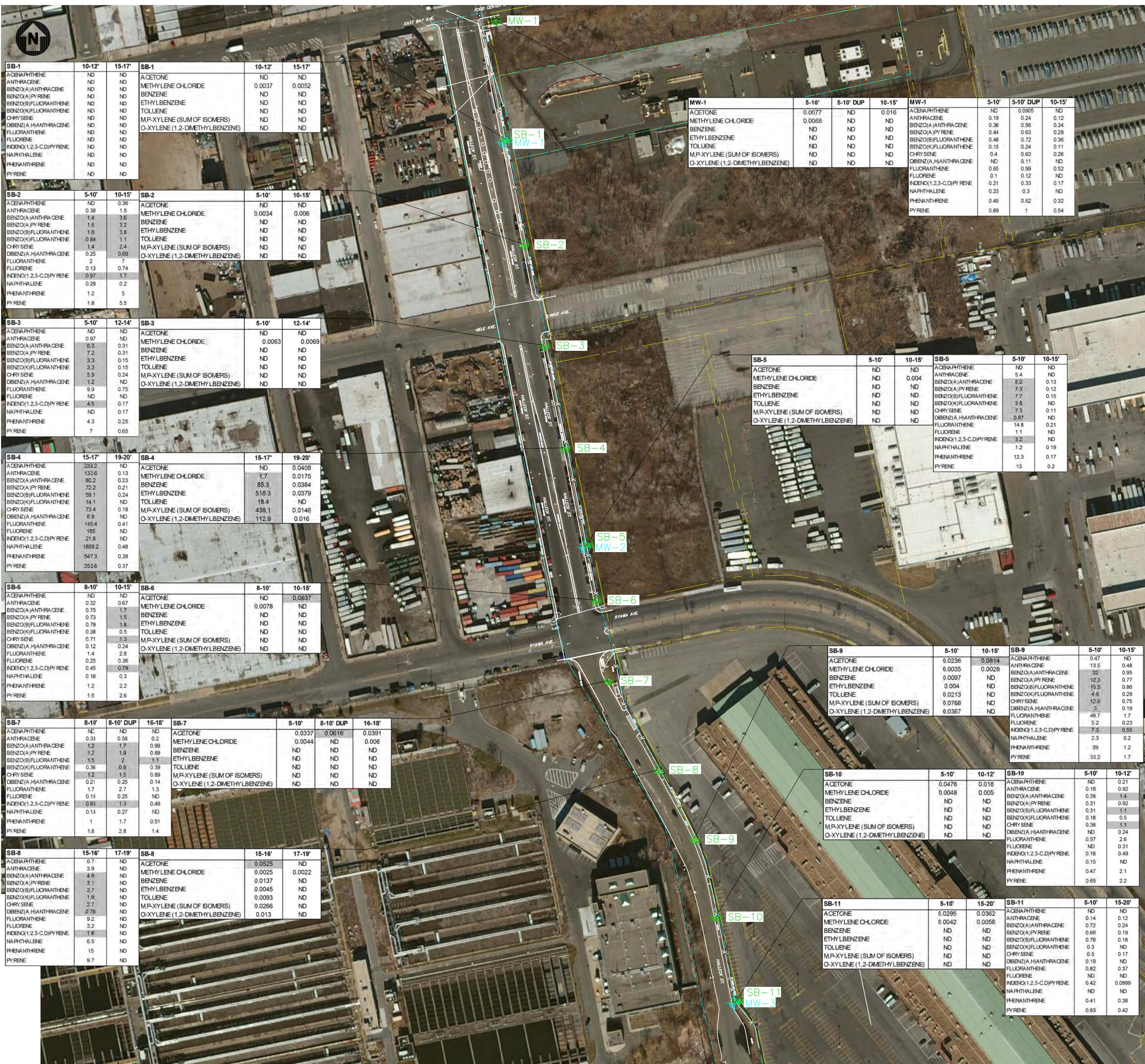
### FIGURE 3

CONSOLIDATED EDISON  
HALLECK STREET SCR  
NEW YORK, NEW YORK

## SAMPLE LOCATION MAP

**PARSONS**

200 COTTONTAIL LANE SOUTH, SOMERSET, N.J. 08873, PHONE: 732-537-3500



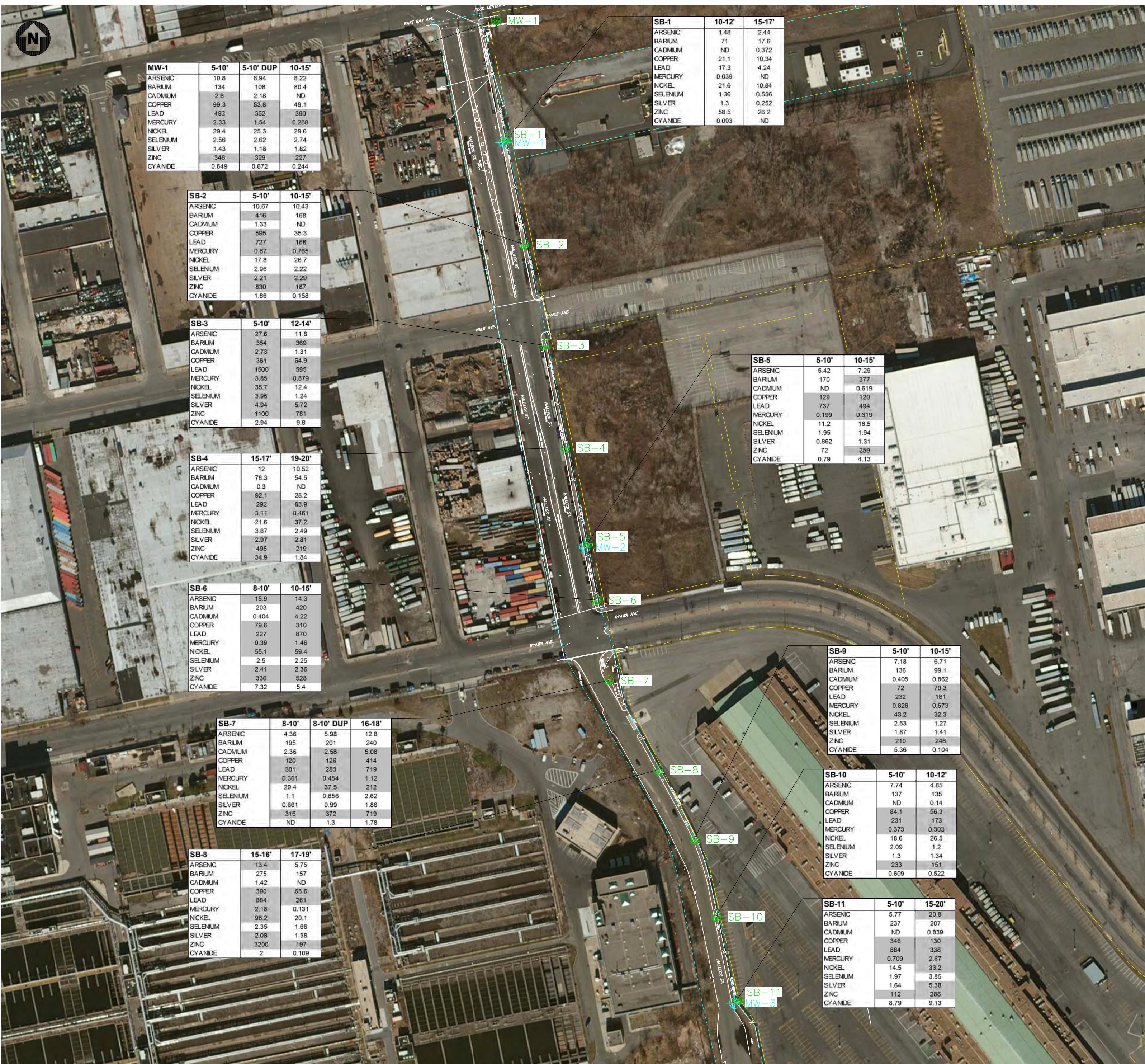
Unrestricted Use Soil Cleanup Objectives Exceedence Criteria		
COMPOUND	VALUE	UNITS
ACETONE	0.05	mg/kg
METHYLENE CHLORIDE	0.05	mg/kg
BENZENE	0.06	mg/kg
ETHYLBENZENE	1.00	mg/kg
TOLUENE	0.70	mg/kg
M.P-XYLENE (SUM OF ISOMERS)	0.26	mg/kg
O-XYLENE (1,2-DIMETHYLBENZENE)	0.26	mg/kg

Unrestricted Use Soil Cleanup Objectives Exceedence Criteria		
COMPOUND	VALUE	UNITS
ACENAPHTHENE	20	mg/kg
ANTHRACENE	100	mg/kg
BENZO(A)ANTHRACENE	1	mg/kg
BENZO(A)PYRENE	1	mg/kg
BENZO(K)FLUORANTHENE	0.8	mg/kg
DIBENZ(A,H)ANTHRACENE	0.33	mg/kg
FLUORANTHENE	100	mg/kg
FLUORENE	30	mg/kg
INDENO(1,2,3-C,D)PYRENE	0.5	mg/kg
NAPHTHALENE	12	mg/kg
PHENANTHRENE	100	mg/kg
PYRENE	100	mg/kg

150 0 150 300

SCALE: 1"=150'

**FIGURE 4**  
CONSOLIDATED EDISON  
HALLECK STREET SCR  
NEW YORK, NY  
SUMMARY OF VOCs AND SVOCs  
IN SOIL



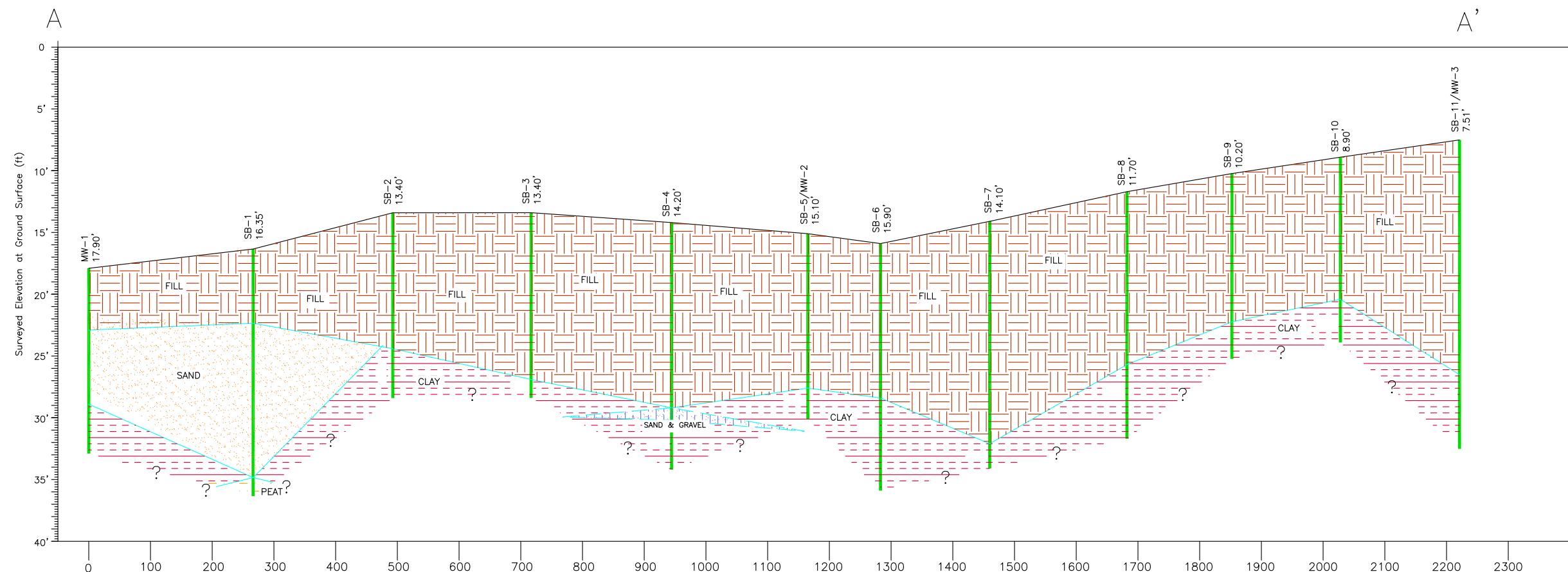


FILE NAME: P:\CONED\448468 HALLECK STREET\CAD\FIGURES\448468-SK003.DWG  
PLOT DATE: 7/3/2014 1:51 PM PLOTTED BY: RUSSO, JILL

FIGURE 6	
CONSOLIDATED EDISON HALLECK STREET SCR NEW YORK, NEW YORK	
SUMMARY OF GROUNDWATER SAMPLE RESULTS	

**PARSONS**

200 COTTONTAIL LANE SOUTH, SOMERSET, NJ. 08873, PHONE: 732-537-3500



10 5 0 10 20  
SCALE: 1"=10'

200 100 0 200 400  
SCALE: 1"=200'

LEGEND:

- FILL
- SAND
- SAND AND GRAVEL
- CLAY
- PEAT
- SOIL BORING

FIGURE 7

CONSOLIDATED EDISON OF NEW YORK  
HALLECK STREET SCR

CROSS SECTION A-A'

**PARSONS**

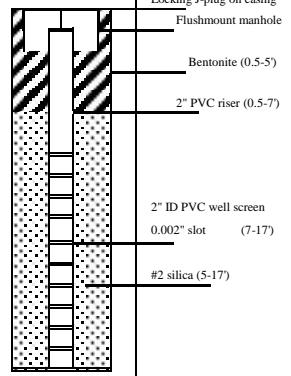
200 COTTONTAIL LANE SOUTH, SOMERSET, N.J. 08873, PHONE: 732-537-3500

## **APPENDIX A**

### **SOIL BORING AND MONITORING WELL LOGS**

				PARSONS DRILLING RECORD			BORING/WELL ID: MW-1		
							Sheet 1 of 1		
Contractor:	Aquifer Drilling & Testing			PROJECT NAME: Consolidated Edison - Former Hunts Point Gas Works			Location Description:		
Driller:	Chris Iodici			PROJECT NUMBER: 446110-04000			Eastern corner of Halleck Street/Food Center Drive Intersection		
Inspector:	Zohar Levy								
Rig Type:	Track Mounted Geoprobe Rig								
GROUNDWATER OBSERVATIONS									
Water Level	DTW	DTW			Weather: Partly cloudy, high 40s			Location Plan	
		12.15						See Site Plan	
Date		11-8-2013			Date/Time Start: 11-4-2013/1050				
Time		0830			Date/Time Finish: 11-4-2013/1110				
Meas. From		Top of Casing							
Sample Depth	Location/ Sample I.D.	SPT	Rec. (%)	PID (ppm)	FIELD IDENTIFICATION OF MATERIAL			SCHEMATIC	COMMENTS
+2									
+1									
0					0-6" Grass, roots, organics 6"-5' Dry, dark brown, fine to medium SAND, some cobble, weathered rock boulders				
1									
2									
3									
4									
5									
6									
7									
8									
9									
10	MW-1 (5-10)	Geoprobe	55	0.0	0-20" Dry, dark brown/grey, medium SAND, trace medium angular to sub-angular gravel 20-30" Moist, grey/dark orange, fine SAND and Silt, some clay, trace sub-angular gravel				
11									
12									
13									
14									
15	MW-1 (10-15)	Geoprobe	65	0.0	0-10" Moist, grey/dark orange, fine SAND and Silt, some clay, trace sub-angular gravel 10-40" Wet, grey, CLAY				
					End of Boring at 15 ft bgs				
SAMPLING METHOD					Hand cleared to 5' bgs, Geoprobe drilled from 5' to 15' bgs. SS = SPLIT SPOON A = AUGER CUTTINGS C = CORED WH = WEIGHT OF HAMMER (RODS)				
					Duplicate sample collected from MW-1 (5-10) interval.				

				PARSONS DRILLING RECORD		BORING/WELL ID: SB-1/MW-1				
						Sheet 1 of 1				
Contractor: Aquifer Drilling & Testing				PROJECT NAME: Consolidated Edison - Former Hunts Point Gas Works		Location Description:				
Driller: Andrea Babel				PROJECT NUMBER: 448468.01000		Halleck St on west side of property boundary				
Inspector: Zohar Levy										
Rig Type: Track Mounted Geoprobe Rig										
GROUNDWATER OBSERVATIONS										
Water Level	DTW	DTW		Weather: Rain, high 60's, windy		Location Plan				
~12' bgs										
Date	11-1-2013			Date/Time Start: 11-1-2013/1310						
Time	1310			Date/Time Finish: 11-1-2013/1400		See Site Plan				
Meas. From	ft bgs - Soil cuttings	Top of Casing								
Sample Depth	Location/ Sample I.D.	SPT	Rec. (%)	PID (ppm)	FIELD IDENTIFICATION OF MATERIAL					
+2										
+1										
0	Vac-tron	NA	0.0	0-6" Grass, roots, organics 6"-5' Dry, dark brown, fine to medium SAND, some cobble, weathered rock boulders						
1										
2										
3										
4										
5	Geoprobe	80	0.0	0-8" Dry, brown, fine to medium SAND, little fine to medium sub-angular to sub-round gravel, trace brick 8-48" Moist, grey/olive, fine to medium SAND, some silt, trace sub-angular gravel						
6										
7										
8										
9										
10	SB-1 (10-12)	Geoprobe	80	0.0	0-20" Moist, grey/olive, fine to medium SAND, some silt, trace sub-angular gravel 20-44" Wet, grey/olive, fine to medium SAND, some silt, trace sub-angular gravel 44-48" Wet, dark grey, fine to medium SAND, some silt, trace sub-angular gravel					
11										
12										
13										
14	SB-1 (15-17)	Geoprobe	100	0.0	44-48" Wet, dark grey, fine to medium SAND, some silt, trace sub-angular gravel 0-16" Wet, grey/olive, fine to medium SAND, some silt, trace sub-angular gravel 16-30" Wet, grey, fine to medium SAND, trace silt 30-60" Peat					
15										
16										
17										
18										
19					End of Boring at 20 ft bgs					
20										
SAMPLING METHOD				Hand cleared to 5 ft bgs; Geoprobe drilled from 5 to 20 ft bgs. Soil boring (SB-1) was converted to a monitoring well (MW-1) on 11/4/13.						
SS = SPLIT SPOON										
A = AUGER CUTTINGS										
C = CORED										
WH = WEIGHT OF HAMMER (RODS)										



				PARSONS DRILLING RECORD				BORING/WELL ID: SB-2 Sheet 1 of 1		
Contractor: Aquifer Drilling & Testing				PROJECT NAME: Consolidated Edison - Former Hunts Point Gas Works				Location Description:		
Driller: Andrea Babel				PROJECT NUMBER: 448468.01000				Halleck St on west side of property boundary		
Inspector: Zohar Levy										
Rig Type: Track Mounted Geoprobe Rig										
GROUNDWATER OBSERVATIONS								Location Plan		
Water Level	DTW	DTW		Weather: Rain, high 60's, windy						
~10' bgs										
Date	11-1-2013			Date/Time Start: 11-1-2013/0940						
Time	940									
Meas. From	ft bgs - Soil cuttings	Top of Casing		Date/Time Finish: 11-1-2013/0955						
Sample Depth	Location/ Sample I.D.	SPT	Rec. (%)	PID (ppm)	FIELD IDENTIFICATION OF MATERIAL				SCHEMATIC	COMMENTS
+2										
+1										
0				NA	0-3" Concrete 3-7" GRAVEL					
1					7"-5' Dry, dark brown, dense, fine to coarse SAND, some Cobble, little fine to coarse sub-angular and sub-round gravel, trace brick, trace boulder, trace silt					
2										
3										
4										
5					0-8" Dry, brown, fine to medium SAND, some sub-angular Gravel, trace brick					
6					8-20" Weathered schist, cobble					
7					20-40" Dry/moist, brown/orange, fine angular GRAVEL and Ash, little fine sand					
8										
9										
10					0-12" Wet, brown/black, fine to medium SAND and fine to medium angular Gravel, some ash, little glass, trace wood					
11					12-42" Wet, grey, CLAY					
12										
13										
14										
15					End of Boring at 15 ft bgs					
SAMPLING METHOD				Hand cleared to 5 ft bgs; Geoprobe drilled from 5 to 15 ft bgs.						
SS = SPLIT SPOON										
A = AUGER CUTTINGS										
C = CORED										
WH = WEIGHT OF HAMMER (RODS)										

				PARSONS DRILLING RECORD			BORING/WELL ID: SB-3 Sheet 1 of 1		
Contractor:	Aquifer Drilling & Testing			PROJECT NAME: Consolidated Edison - Former Hunts Point Gas Works			Location Description:		
Driller:	Andrea Babel			PROJECT NUMBER: 448468.01000			Halleck St on west side of property boundary		
Inspector:	Zohar Levy								
Rig Type:	Track Mounted Geoprobe Rig								
GROUNDWATER OBSERVATIONS							Location Plan		
Water Level	DTW	DTW		Weather: Rain, high 60's, windy					
~10' bgs									
Date	11-1-2013			Date/Time Start: 11-1-2013/1110					
Time	1110								
Meas. From	ft bgs - Soil cuttings	Top of Casing		Date/Time Finish: 11-1-2013/1125					
Sample Depth	Location/ Sample I.D.	SPT	Rec. (%)	PID (ppm)	FIELD IDENTIFICATION OF MATERIAL			SCHEMATIC	COMMENTS
+2									
+1									
0					NA 0-3" Concrete 3-8" GRAVEL				
1					0.0 8"-5' Dry, brown, dense, fine to coarse SAND, some fine to coarse sub-angular and sub-round gravel, some cobble, trace brick, trace concrete, trace wood, trace silt				
2									
3									
4									
5					0-40" Dry, brown/black, fine to medium SAND, some brick, little cobble, little f-c angular-subangular Gravel				
6					40-48" Moist, dark brown/orange, fine to medium SAND and fine to medium angular Gravel, some ash				
7									
8									
9									
10					0-24" Wet, dark brown/orange, fine to medium SAND and fine to medium angular Gravel, some ash				
11					24-44" Moist/wet, black, fine SAND and fine angular Gravel, some wood, trace ash, trace glass				
12					44-55" Wet, grey, CLAY				
13									
14									
15					End of Boring at 15 ft bgs				
SAMPLING METHOD				Hand cleared to 5 ft bgs; Geoprobe drilled from 5 to 15 ft bgs.					
SS = SPLIT SPOON									
A = AUGER CUTTINGS									
C = CORED									
WH = WEIGHT OF HAMMER (RODS)									

				PARSONS DRILLING RECORD				BORING/WELL ID: SB-4 Sheet 1 of 1			
Contractor: Aquifer Drilling & Testing				PROJECT NAME: Consolidated Edison - Former Hunts Point Gas Works				Location Description:			
Driller: Andrea Babel				PROJECT NUMBER: 448468.01000				Halleck St on west side of property boundary			
Inspector: Zohar Levy											
Rig Type: Track Mounted Geoprobe Rig											
GROUNDWATER OBSERVATIONS								Location Plan			
Water Level	DTW	DTW		Weather: Clear, low 60s				See Site Plan			
	~11' bgs										
Date	10-31-2013			Date/Time Start: 10-31-2013/1345							
Time	1345			Date/Time Finish: 10-31-2013/1415							
Meas. From	ft bgs - Soil cuttings	Top of Casing									
Sample Depth	Location/ Sample I.D.	SPT	Rec. (%)	PID (ppm)	FIELD IDENTIFICATION OF MATERIAL				SCHEMATIC	COMMENTS	
+2											
+1											
0	Vac-tron	NA	NA	0.0	0-6" Concrete 6-10" GRAVEL 10"-5' Dry, brown, fine to medium SAND, some fine to coarse sub-angular gravel, little cobble, little brick, trace concrete, trace wood						
1											
2			Geoprobe	75	0.5	0-30" Dry, brown/black, fine to medium SAND, little fine to medium sub-angular gravel, trace brick 30-40" Dry/moist, brown/grey, fine to medium SAND, trace fine sub-angular gravel, hydrocarbon odor					
3											
4			52.0								
5											
6	Geoprobe	40	137.0	0-10" Dry/moist, brown/grey, fine to medium SAND, trace fine sub-angular gravel, hydrocarbon odor 10-24" Moist/wet, black, fine SAND and Wood, NAPL saturated, strong hydrocarbon odor							
7											
8					698.0						
9											
10	Geoprobe	100	1260.0	0-24" Wet, black, fine to medium SAND and fine angular to sub-angular Gravel, trace wood, NAPL saturated, strong hydrocarbon odor							
11											
12					350.0	24-30" Wet, grey/black, CLAY, NAPL staining					
13											
14			3.5	30-60" Wet, grey, CLAY							
15											
16	SB-4 (15-17)										
17											
18	Geoprobe										
19											
20					End of Boring at 20 ft bgs						
SAMPLING METHOD				Hand cleared to 5 ft bgs; Geoprobe drilled from 5 to 20 ft bgs.							
SS = SPLIT SPOON				NAPL observed from ~11-18 ft bgs.							
A = AUGER CUTTINGS				NAPL sample collected from 15-17 ft bgs and submitted for Fingerprint Analysis.							
C = CORED											
WH = WEIGHT OF HAMMER (RODS)											

				PARSONS DRILLING RECORD		BORING/WELL ID: SB-5/MW-2
						Sheet 1 of 1
				PROJECT NAME: Consolidated Edison - Former Hunts Point Gas Works		Location Description:
				PROJECT NUMBER: 448468.01000		Halleck St on west side of property boundary
GROUNDWATER OBSERVATIONS						Location Plan
Water Level	DTW	DTW		Weather: Clear, low 60s		See Site Plan
	~10' bgs	9.75				
Date	10-31-2013	11-8-2013		Date/Time Start: 10-31-2013/1240		
Time	1240	0830		Date/Time Finish: 10-31-2013/1300		
Meas. From	ft bgs - Soil cuttings	Top of Casing		FIELD IDENTIFICATION OF MATERIAL		SCHEMATIC
Sample Depth	Location/ Sample I.D.	SPT	Rec. (%)	PID (ppm)		COMMENTS
+2						
+1						
0	Vac-tron	NA	0.0	0-3" Concrete		
1				3-7" GRAVEL		
2				7"-4.5" Dry, dark grey, dense, fine to coarse SAND, some cobble, little fine to coarse sub-angular to sub-round gravel, trace brick, trace boulder		
3				4.5-5" Dry, brown with red-orange mottling, dense, fine to coarse SAND, some cobble, little fine to coarse sub-angular gravel, trace silt		
4						
5	SB-5 (5-10)	Geoprobe	40	0-24" Dry, grey/black, fine to medium SAND and Ash, some fine angular to sub-angular gravel		
6						
7						
8						
9						
10	SB-5 (10-15)	Geoprobe	80	0-30" Wet, grey/black, fine to medium SAND and Ash, some fine angular to sub-angular gravel		
11				30-48" Wet, grey, CLAY		
12						
13						
14						
15						
16	End of Boring at 16 ft bgs					
SAMPLING METHOD				Hand cleared to 5 ft bgs; Geoprobe drilled from 5 to 16 ft bgs		
SS = SPLIT SPOON				Soil boring (SB-5) was converted to a monitoring well (MW-2) on 11/5/13.		
A = AUGER CUTTINGS						
C = CORED						
WH = WEIGHT OF HAMMER (RODS)						

				PARSONS DRILLING RECORD				BORING/WELL ID: SB-6			
								Sheet 1 of 1			
Contractor: Aquifer Drilling & Testing				PROJECT NAME: Consolidated Edison - Former Hunts Point Gas Works				Location Description:			
Driller: Andrea Babel				PROJECT NUMBER: 448468.01000				Halleck St on west side of property boundary			
Inspector: Zohar Levy											
Rig Type: Track Mounted Geoprobe Rig											
GROUNDWATER OBSERVATIONS								Location Plan			
Water Level	DTW	DTW		Weather: Clear, low 60s							
~10' bgs											
Date	10-31-2013			Date/Time Start: 10-31-2013/1015				See Site Plan			
Time	1015			Date/Time Finish: 10-31-2013/1040							
Meas. From	ft bgs - Soil cuttings	Top of Casing									
Sample Depth	Location/ Sample I.D.	SPT	Rec. (%)	PID (ppm)	FIELD IDENTIFICATION OF MATERIAL				SCHEMATIC	COMMENTS	
+2											
+1											
0				NA	0-3" Concrete 3-8" GRAVEL						
1					8"-5' Dry, dark grey, dense, fine to coarse SAND, some cobble, little fine to coarse sub-angular to sub-round gravel, trace brick						
2											
3											
4											
5					0-30" Dry, grey/brown, fine to medium SAND, some fine to medium angular to sub-angular gravel, little ash, trace brick						
6					30-40" Moist, grey/brown, fine to medium SAND, some fine to medium angular to sub-angular gravel, little ash, trace brick						
7											
8	SB-6 (8-10)				0-10" Wet, grey/brown, fine to medium SAND, some fine to medium angular to sub-angular gravel, little ash, trace brick						
9					10-26" Dry/moist, black, fine to medium SAND and medium to coarse sub-angular Gravel						
10					26-30" Wet, grey, CLAY						
11											
12											
13											
14											
15					0-50" Wet, grey, CLAY						
16											
17											
18											
19											
20					End of Boring at 20 ft bgs						
SAMPLING METHOD				Hand cleared to 5 ft bgs; Geoprobe drilled from 5 to 20 ft bgs.							
SS = SPLIT SPOON											
A = AUGER CUTTINGS											
C = CORED											
WH = WEIGHT OF HAMMER (RODS)											

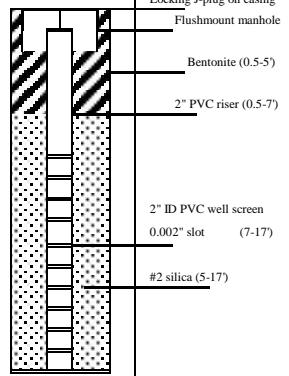
				PARSONS DRILLING RECORD				BORING/WELL ID: SB-7 Sheet 1 of 1		
Contractor: Aquifer Drilling & Testing				PROJECT NAME: Consolidated Edison - Former Hunts Point Gas Works				Location Description:		
Driller: Andrea Babel				PROJECT NUMBER: 448468.01000				Halleck Street on west side of property boundary		
Inspector: Zohar Levy										
Rig Type: Track Mounted Geoprobe Rig										
GROUNDWATER OBSERVATIONS								Location Plan		
Water Level	DTW	DTW		Weather: Clear, low 60s				See Site Plan		
	~10' bgs									
Date	10-31-2013			Date/Time Start: 10-31-2013/0835						
Time	835			Date/Time Finish: 10-31-2013/0850						
Meas. From	ft bgs - Soil cuttings	Top of Casing								
Sample Depth	Location/ Sample I.D.	SPT	Rec. (%)	PID (ppm)	FIELD IDENTIFICATION OF MATERIAL				SCHEMATIC	COMMENTS
+2										
+1										
0					0-6" Grass, roots, organics					
1					6"-5" Dry, brown, fine to coarse SAND, some cobble, little fine to coarse sub-angular gravel, little brick, trace concrete debris					
2										
3										
4										
5					0-40" Dry, brown/black, fine to medium SAND, some fine to coarse angular to sub-round gravel, trace wood, trace fabric, trace brick					
6					40-60" Moist, brown/black, fine to medium SAND, some fine angular to sub-round gravel, trace wood, trace fabric, trace brick, trace ash					
7										
8	SB-7 (8-10)									
9										
10					0-36" Wet, grey/black, medium SAND and fine angular Gravel, some shells, little ash					
11										
12										
13										
14										
15					0-40" Wet, grey/black, medium SAND and fine angular Gravel, some shells, little ash, trace glass, trace wood					
16	SB-7 (16-18)				40-60" Wet, grey, CLAY					
17										
18										
19										
20					End of Boring at 20 ft bgs					
SAMPLING METHOD				Hand cleared to 5 ft bgs; Geoprobe drilled from 5 to 20 ft bgs						
SS = SPLIT SPOON										
A = AUGER CUTTINGS										
C = CORED										
WH = WEIGHT OF HAMMER (RODS)										

				PARSONS DRILLING RECORD				BORING/WELL ID: SB-8 Sheet 1 of 1		
Contractor: Aquifer Drilling & Testing				PROJECT NAME: Consolidated Edison - Former Hunts Point Gas Works				Location Description:		
Driller: Andrea Babel				PROJECT NUMBER: 448468.01000				Halleck Street on west side of property boundary		
Inspector: Zohar Levy										
Rig Type: Track Mounted Geoprobe Rig										
GROUNDWATER OBSERVATIONS								Location Plan		
Water Level	DTW	DTW		Weather: Clear, low 60s				See Site Plan		
	~8' bgs									
Date	10-30-2013			Date/Time Start: 10-30-2013/1430						
Time	1430			Date/Time Finish: 10-30-2013/1455						
Meas. From	ft bgs - Soil cuttings	Top of Casing								
Sample Depth	Location/ Sample I.D.	SPT	Rec. (%)	PID (ppm)	FIELD IDENTIFICATION OF MATERIAL				SCHEMATIC	COMMENTS
+2										
+1										
0					0-6" Grass, roots, organics					
1					6"-2' Dry, brown, fine to coarse SAND, some cobble, little brick, trace concrete debris					
2					2-2.5' Concrete debris					
3					2.5-5' Dry, brown, fine to coarse SAND, some cobble, little brick, little concrete debris					
4										
5					0-12" Dry, brown, fine to medium SAND, some fine sub-round gravel					
6					12-30" Moist, grey/brown, fine to medium SAND, some fine angular to sub-round gravel, little brick, trace wood					
7					30-42" Moist/wet, black, fine to medium SAND and angular to sub-angular gravel, little wood, trace brick					
8										
9										
10					0-36" Wet, black, fine to medium SAND and fine to medium angular to sub-angular Gravel, trace wire					
11										
12										
13										
14										
15	SB-8 (15-16)			10.7	0-12" Wet, black, fine to medium SAND and fine to medium angular to sub-angular Gravel, little wood, slight hydrocarbon odor					
16					12-60" Wet, grey, CLAY, trace shells					
17	SB-8 (17-19)			100						
18										
19										
20					End of Boring at 20 ft bgs					
SAMPLING METHOD				Hand cleared to 5 ft bgs; Geoprobe drilled from 5 to 20 ft bgs						
SS = SPLIT SPOON										
A = AUGER CUTTINGS										
C = CORED										
WH = WEIGHT OF HAMMER (RODS)										

				PARSONS DRILLING RECORD			BORING/WELL ID: SB-9		
Contractor: Aquifer Drilling & Testing				PROJECT NAME: Consolidated Edison - Former Hunts Point Gas Works			Sheet 1 of 1		
Driller: Andrea Babel				PROJECT NUMBER: 448468.01000			Location Description:		
Inspector: Zohar Levy							Halleck Street on west side of property boundary		
Rig Type: Track Mounted Geoprobe Rig									
GROUNDWATER OBSERVATIONS							Location Plan		
Water Level	DTW	DTW		Weather: Clear, low 60s			See Site Plan		
	~9' bgs								
Date	10-30-2013			Date/Time Start: 10-30-2013/1320					
Time	1320			Date/Time Finish: 10-30-2013/1340					
Meas. From	ft bgs - Soil cuttings	Top of Casing		FIELD IDENTIFICATION OF MATERIAL			SCHEMATIC	COMMENTS	
Sample Depth	Location/ Sample I.D.	SPT	Rec. (%)	PID (ppm)					
+2									
+1									
0					0-6" Grass, roots, organics 6"-5' Dry, brown, fine to coarse SAND, some cobble, little fine to coarse sub-angular gravel, little brick, trace concrete debris				
1									
2									
3									
4									
5					0-20" Moist, brown, fine to medium SAND, some fine to medium sub-angular to sub-round gravel, trace silt, trace brick				
6					20-36" Moist, black, fine to medium SAND, some fine angular to sub-angular gravel, little silt				
7					36-42" COBBLE				
8					42-48" Moist/wet, black, fine to medium SAND and fine angular Gravel, little coal				
9									
10					0-18" Moist/wet, black, fine to medium SAND and fine to medium angular to sub-round Gravel, trace coal, trace wood				
11					18-48" Wet, grey, CLAY				
12					30-60" Wet, grey, CLAY, trace shells				
13									
14									
15					End of Boring at 15 ft bgs				
SAMPLING METHOD				Hand cleared to 5 ft bgs; Geoprobe drilled from 5 to 15 ft bgs.					
SS = SPLIT SPOON									
A = AUGER CUTTINGS									
C = CORED									
WH = WEIGHT OF HAMMER (RODS)									

				PARSONS DRILLING RECORD				BORING/WELL ID: SB-10		
								Sheet 1 of 1		
Contractor: Aquifer Drilling & Testing				PROJECT NAME: Consolidated Edison - Former Hunts Point Gas Works				Location Description:		
Driller: Andrea Babel				PROJECT NUMBER: 448468.01000				Halleck Street on west side of property boundary		
Inspector: Zohar Levy										
Rig Type: Track Mounted Geoprobe Rig										
GROUNDWATER OBSERVATIONS								Location Plan		
Water Level	DTW	DTW		Weather: Clear, low 60s				See Site Plan		
	~10' bgs									
Date	10-30-2013			Date/Time Start: 10-30-2013/1115						
Time	1115			Date/Time Finish: 10-30-2013/1130						
Meas. From	ft bgs - Soil cuttings	Top of Casing		FIELD IDENTIFICATION OF MATERIAL				SCHEMATIC	COMMENTS	
Sample Depth	Location/ Sample I.D.	SPT	Rec. (%)	PID (ppm)						
+2										
+1										
0					0-6" Grass, roots					
1					6"-4.5" Dry, brown, fine to coarse SAND, some fine to coarse sub-angular to sub-round gravel, little brick, little cobble, trace concrete					
2					4.5"- Dry, brown, fine to coarse SAND, some fine to coarse sub-angular to sub-round gravel, some brick, little cobble, trace concrete					
3										
4										
5					0-12" Dry, brown, fine to medium SAND, some fine to coarse sub-round gravel					
6					12-32" Moist, grey, medium SAND, some fine to medium angular to sub-angular gravel, trace brick					
7					32-48" Moist, black/grey, medium SAND and fine to medium angular to sub-angular Gravel					
8										
9										
10					0-16" Wet, black/grey, medium SAND and fine to medium angular to sub-angular Gravel, trace brick					
11	SB-10 (5-10)	Geoprobe	80	0.0	16-30" Wet, grey, CLAY, some Shells					
12					30-60" Wet, grey, CLAY, trace shells					
13										
14										
15					End of Boring at 15 ft bgs					
SAMPLING METHOD					Hand cleared to 5 ft bgs; Geoprobe drilled from 5 to 15 ft bgs					
SS = SPLIT SPOON										
A = AUGER CUTTINGS										
C = CORED										
WH = WEIGHT OF HAMMER (RODS)										

				PARSONS DRILLING RECORD		BORING/WELL ID: SB-11/MW-3		
				PROJECT NAME: Consolidated Edison - Former Hunts Point Gas Works		Sheet 1 of 1		
				PROJECT NUMBER: 448468.01000		Location Description:		
						Halleck Street on west side of property boundary		
						<b>Location Plan</b>		
						See Site Plan		
GROUNDWATER OBSERVATIONS								
Water Level	DTW	DTW		Weather: Clear, low 60s				
-7' bgs		5.95						
Date	10-30-2013	11-8-2013		Date/Time Start: 10-30-2013/0930				
Time	930	0830		Date/Time Finish: 10-30-2013/1015				
Meas. From	ft bgs - Soil cuttings	Top of Casing						
Sample Depth	Location/ Sample I.D.	SPT	Rec. (%)	PID (ppm)	FIELD IDENTIFICATION OF MATERIAL		SCHEMATIC	COMMENTS
+2								
+1								
0					0-6" Grass, roots 6-4" Moist/dry, brown, fine to coarse SAND, some fine to coarse sub-angular to sub-round gravel, little concrete debris, trace brick 4-5" Moist/dry, brown, fine to coarse SAND, some fine to coarse sub-angular to sub-round gravel and concrete debris, trace brick			Locking J-plug on casing
1								Flushmount manhole
2								Bentonite (0.5')
3								2" PVC riser (0.5-7')
4								
5					0-18" Dry, grey/brown, medium SAND and fine to medium angular to sub-angular Gravel, little ash 18-24" Moist, black, medium SAND and fine to medium angular to sub-angular Gravel, little ash 24-30" Wet, black, medium SAND and fine to medium angular to sub-angular Gravel, little ash			2" ID PVC well screen
6								0.002" slot (7-17')
7								#2 silica (5-17)
8								
9								
10					0-24" Wet, black, medium SAND and fine to medium angular to sub-angular Gravel, trace concrete			
11								
12								
13								
14								
15					0-20" Wet, black, medium SAND and fine to medium angular to sub-angular Gravel, trace concrete, trace porcelain 20-24" Moist, grey, CLAY			
16								
17								
18								
19								
20					0-50" Moist, grey, CLAY, trace shells			
21								
22								
23								
24								
25					End of Boring at 25 ft bgs			
SAMPLING METHOD				Hand cleared to 5 ft bgs; Geoprobe drilled from 5 to 25 ft bgs				
SS = SPLIT SPOON				Soil boring (SB-11) was converted to a monitoring well (MW-3) on 11/5/13.				
A = AUGER CUTTINGS								
C = CORED								
WH = WEIGHT OF HAMMER (RODS)								



**APPENDIX B**

**GROUNDWATER SAMPLING LOGS**

**PARSONS**  
**GROUNDWATER SAMPLING RECORD**

SITE NAME: Con Edison (Farmington Hill Water) *Halleck*  
 PROJECT NUMBER: 446407-00000  
 Purge Date: 11/27/13  
 Sampling Date: 11/27/13  
 Samplers: Zohar Levy  
 SAMPLE ID: MW-1  
 Sampling Method: Low flow purge (Monsoon Pump)

**WELL PURGING**

Static Water Level (TOC):  
Depth to Well Bottom (TOC):

**CALCULATIONS:**  
2-Inch Casing:  
3-Inch Casing:  
4-Inch Casing:  
Method:

Ft. of Water in Well      X (GAL / FT) = \_\_\_\_\_ Gallons  
 Ft. of Water in Well      x 0.16 = \_\_\_\_\_ Gallons  
 Ft. of Water in Well      x 0.32 = \_\_\_\_\_ Gallons  
 Ft. of Water in Well      x 0.64 = \_\_\_\_\_ Gallons  
 Low Flow Pump

**SAMPLE DESCRIPTION**

Odor:  
Other:

**FIELD TESTS**

Time  
Depth To Water (TOC) (ft)  
Depth To Pump (TOC) (ft)  
Flow Rate (ml/min)  
Volume of Water Purged  
pH (s.u.)  
Conductivity (mS/cm)  
Turbidity (NTUs)  
Dissolved Oxygen (mg/L)  
Temperature (Degrees C)  
ORP (mV)  
Salinity (%)  
TDS (g/L)

PURGE	SAMPLE							
1105	1150	1155	1200	1205	1210	1215	1220	
12.4								
16.15								
~200	~250	~250	~250	~250	~250	~250	~250	~250
~0.25	~0.25	~0.25	~0.25	~0.25	~0.25	~0.25	~0.25	~0.25
6.77	6.77	6.77	6.77	6.77	6.77	6.77	6.77	6.77
3.16	3.46	3.65	4.26	4.36	4.4	4.43	4.45	
1.9	6.38	9.15	23.4	11.0	50.7	30.2	10.3	
17.03	17.67	17.77	17.79	17.88	17.95	17.94	17.92	
-48	-68	-77	-90	-93	-98	-100	-101	
1.7	1.9	1.9	2.3	2.3	2.3	2.3	2.3	
2.04	2.37	2.34	2.72	2.75	2.78	2.78	2.8	

**SAMPLE ANALYSIS / LABORATORY**

Analyze For:

TCL VOC's , TCL SVOCs , TAL Metals, CN

Shipped Via:  
Laboratory

Chemtech

Other Notes:

**PARSONS**  
**GROUNDWATER SAMPLING RECORD**

**SITE NAME:** Con Edison (Ewington Ct Holder) Heller St.  
**PROJECT NUMBER:** 44607-00000  
**Purge Date:** 11/27/13  
**Sampling Date:** 11/27/13  
**Samplers:** Zohar Levy of Parsons / Somerset, NJ  
**SAMPLE ID:** MW-2  
**Sampling Method:** Low flow purge (Monsoon Pump)

**WELL PURGING**

Static Water Level (TOC):

Depth to Well Bottom (TOC):

**CALCULATIONS:**

2-inch Casing:

3-inch Casing:

4-inch Casing:

Method:

9.47

16.1

Ft. of Water in Well \_\_\_\_\_ X (GAL / FT) = \_\_\_\_\_ Gallons  
 Ft. of Water in Well \_\_\_\_\_ x 0.16 = \_\_\_\_\_ Gallons  
 Ft. of Water in Well \_\_\_\_\_ x 0.32 = \_\_\_\_\_ Gallons  
 Ft. of Water in Well \_\_\_\_\_ x 0.64 = \_\_\_\_\_ Gallons  
 Low Flow Pump

**SAMPLE DESCRIPTION**

Odor :

Other :

**FIELD TESTS**

Time

Depth To Water (TOC) (ft)

Depth To Pump (TOC) (ft)

Flow Rate (ml/min)

Volume of Water Purged

pH (s.u.)

Conductivity (mS/cm)

Turbidity (NTUs)

Dissolved Oxygen (mg/L)

Temperature (Degrees C)

ORP (mV)

Salinity (%)

TDS (g/L)

PURGE	SAMPLE							
1015	1020	1025	1030	1035	1040	1045	1050	1050
15.1								3
~250	~250	~200	~250	~250	~250	~250	~250	~250
~0.5	~1	-1.25	+1.5	2	-2.5	-3	3.5	
7.21	7.21	7.21	7.20	7.19	7.18	7.16	7.18	
2.25	2.22	2.18	2.15	2.13	2.12	2.11	2.10	
615	367	203	131	69.3	48.2	20.9	0	
22.23	7.87	2.01	1.67	1.48	1.91	1.36	1.36	
17.24	17.29	17.03	17.66	17.73	17.85	17.83	17.81	
-124	-127	-131	-103	-144	-146	-147	-147	
1.2	1.2	1.2	1.1	1.1	1.1	1.1	1.1	
1.44	1.41	1.38	1.37	1.36	1.35	1.35	1.35	

**SAMPLE ANALYSIS / LABORATORY**

Analyze For:

TCL VOC's , TCL SVOCs , TAC Metals, CN

Shipped Via:

Chemtech

Laboratory

Other Notes:

**PARSONS**  
**GROUNDWATER SAMPLING RECORD**

**SITE NAME:** Con Edison (Parson St. Holder) *Halkirk St.*  
**PROJECT NUMBER:** 446197-00000  
**Purge Date:** 4/27/13  
**Sampling Date:** 4/27/13  
**Samplers:** Zohar Levy of Parsons / Somerset, NJ  
  
**SAMPLE ID:** MW-3  
**Sampling Method:** Low flow purge (Monsoon Pump)

**WELL PURGING**

Static Water Level (TOC):  
Depth to Well Bottom (TOC):

**CALCULATIONS:**  
2-inch Casing:  
3-inch Casing:  
4-inch Casing:  
Method:

*5.6*  
*17.25*  
 Ft. of Water in Well \_\_\_\_\_ X (GAL / FT) = \_\_\_\_\_ Gallons  
 Ft. of Water in Well \_\_\_\_\_ x 0.16 = \_\_\_\_\_ Gallons  
 Ft. of Water in Well \_\_\_\_\_ x 0.32 = \_\_\_\_\_ Gallons  
 Ft. of Water in Well \_\_\_\_\_ x 0.64 = \_\_\_\_\_ Gallons  
 Low Flow Pump

**SAMPLE DESCRIPTION**

Odor:  
Other:

**FIELD TESTS**

Time  
Depth To Water (TOC) (ft)  
Depth To Pump (TOC) (ft)  
Flow Rate (ml/min)  
Volume of Water Purged  
pH (s.u.)  
Conductivity (mS/cm)  
Turbidity (NTUs)  
Dissolved Oxygen (mg/L)  
Temperature (Degrees C)  
ORP (mV)  
Salinity (%)  
TDS (g/L)

PURGE	SAMPLE							
0810	0815	0820	0825	0830	0835	0840	0845	
5.91	6.0	5.93	5.85	5.97	5.96	5.81	5.73	
16.25								
~350	~350	~350	~350	~350	~350	~350	~350	
-0.75	-1.5	-2.25	-3	-3.5	-4	-4.5	-5	
7.12	7.16	7.17	7.17	7.17	7.18	7.18	7.18	
5.56	5.48	6.24	6.28	6.3	6.3	6.31	6.3	
3.41	1.32	49.7	5.5	0	0	0	0	
3.77	1.75	1.43	1.19	1.05	1.0	.98	.97	
17.71	17.35	17.65	17.78	17.73	17.5	17.82	17.79	
-119	-152	-190	-199	-156	-161	-165	-168	
3.0	3.3	3.4	3.6	3.4	3.4	3.5	3.5	
3.52	3.78	3.95	3.95	3.95	3.95	3.95	3.95	

**SAMPLE ANALYSIS / LABORATORY**

Analyze For: TCL VOC's , TCL SVOCs , TAL Met-S, CN

Shipped Via: Chemtech

Laboratory

Other Notes:

**APPENDIX C**

**DATA USABILITY SUMMARY REPORT**

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# **DATA USABILITY SUMMARY REPORT**

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## **HALLECK STREET SIDEWALK SITE**

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*Prepared For:*



**CONSOLIDATED EDISON COMPANY OF NEW YORK, INC.**

**31-01 20<sup>th</sup> Avenue  
Long Island City, NY 11105**

*Prepared By:*

**PARSONS**

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**FEBRUARY 2014**

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## LIST OF ATTACHMENTS

### ATTACHMENT A VALIDATED LABORATORY DAT

#### ATTACHMENT A-1 VALIDATED LABORATORY DATA FOR SOIL SAMPLES

#### ATTACHMENT A-2 VALIDATED LABORATORY DATA FOR GROUNDWATER SAMPLES

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PARSONS

# **SECTION 1**

## **DATA USABILITY SUMMARY**

Soil and groundwater samples were collected from the Consolidated Edison Halleck Street site in Bronx, New York from October 30, 2013 through November 27, 2013. Analytical results from these samples were validated and reviewed by Parsons for usability with respect to the following requirements:

- Work Plan,
- NYSDEC Analytical Services Protocol (ASP), and
- USEPA Region II Standard Operating Procedures (SOPs) for organic and inorganic data review.

The analytical laboratory for this project was Chemtech. This laboratory is certified to perform project analyses through the New York State Department of Health (NYSDOH) Environmental Laboratory Approval Program (ELAP).

### **1.1 LABORATORY DATA PACKAGES**

The laboratory data package turnaround time, defined as the time from sample receipt by the laboratory to receipt of the analytical data packages by Parsons, was 28 days on average for the project samples.

The data packages received from Chemtech were paginated, complete, and overall were of good quality. Comments on specific quality control (QC) and other requirements are discussed in detail in the attached data validation report which is summarized in Section 2.

### **1.2 SAMPLING AND CHAIN-OF-CUSTODY**

The samples were collected, properly preserved, shipped under a chain-of-custody (COC) record, and received at Chemtech within one day of sampling. All samples were received intact and in good condition at Chemtech.

### **1.3 LABORATORY ANALYTICAL METHODS**

The soil and groundwater samples that were collected from the site were analyzed for volatile organic compounds (VOCs), semivolatile organic compounds (SVOCs), metals, and cyanide. Summaries of issues concerning these laboratory analyses are presented in Subsections 1.3.1 through 1.3.3. The data qualifications resulting from the data validation review and statements on the laboratory analytical precision, accuracy, representativeness, completeness, and comparability (PARCC) are discussed for each analytical method in Section 2. The laboratory data were reviewed and may be qualified with the following validation flags:

- "U" - not detected at the value given,
- "UJ" - estimated and not detected at the value given,

"J" - estimated at the value given,

"N" - presumptive evidence at the value given, and

"R" - unusable value.

The validated laboratory data were tabulated and are presented in Attachment A.

### **1.3.1 Volatile Organic Analysis**

Soil and groundwater samples were analyzed for VOCs using the USEPA SW-846 8260C analytical method. Certain reported results for the VOC samples were qualified as estimated based upon laboratory control sample (LCS) recoveries, instrument calibrations, and internal standard responses. Certain reported VOC analytical results were considered unusable and qualified "R" based upon poor instrument calibrations. The reported VOC analytical results were 99.5% to 100% complete (i.e., usable) for the soil and groundwater data, respectively. PARCC requirements were met overall.

### **1.3.2 Semivolatile Organic Analysis**

Soil and groundwater samples were analyzed for SVOCs using the USEPA SW-846 8270D analytical method. Certain reported results for the SVOC samples were qualified as estimated based upon laboratory control sample (LCS) recoveries, matrix spike/matrix spike duplicate (MS/MSD) precision and accuracy, instrument calibrations, and field duplicate precision. Certain reported SVOC analytical results were considered unusable and qualified "R" based upon poor LCS recoveries and MS/MSD precision and accuracy. The reported SVOC analytical results were 98.6% to 100% complete (i.e., usable) for the soil and groundwater data, respectively. PARCC requirements were met overall.

### **1.3.3 Inorganics Analysis**

Soil and groundwater samples were analyzed for metals and cyanide using the USEPA SW-846 6010C/7470A/7471A/9012B analytical methods. Certain reported results for the inorganics samples were qualified as estimated based upon matrix spike recoveries, interference check sample recoveries, serial dilutions, and field duplicate precision. The reported inorganic analytical results were considered 100% complete (i.e., usable) for the soil and groundwater data. PARCC requirements were met.

## **SECTION 2**

### **DATA VALIDATION REPORT**

#### **2.1 SOIL**

Data review has been completed for data packages generated by Chemtech containing soil samples collected from the site. All of these samples were properly preserved, shipped under a COC record, and received intact by the analytical laboratory. The analytical results were presented by the laboratory in two sample delivery groups (SDGs): E4277 and E4340. Data validation was performed for all samples in accordance with the most current editions of the USEPA Region II SOPs for organic and inorganic data review. This data validation and usability report is presented by analysis type and the validated laboratory data are presented in Attachment A-1.

##### **2.1.1 Volatiles**

The following items were reviewed for compliancy in the volatile analysis:

- Custody documentation
- Holding times
- Surrogate recoveries
- Matrix spike/matrix spike duplicate (MS/MSD) precision and accuracy
- Laboratory control sample (LCS) recoveries
- Laboratory method blank and field equipment blank contamination
- GC/MS instrument performance
- Initial and continuing calibrations
- Internal standard area counts and retention times
- Field duplicate precision
- Sample result verification and identification
- Quantitation limits
- Data completeness

These items were considered compliant and acceptable in accordance with the validation protocols with the exception of MS/MSD precision and accuracy, LCS recoveries, blank contamination, initial and continuing calibrations, and internal standard responses as discussed below.

#### MS/MSD Precision and Accuracy

All MS/MSD precision (relative percent difference; RPD) and accuracy (percent recovery; %R) measurements were considered acceptable and within QC limits for designated spiked

project samples with the exception of the high MS/MSD accuracy results for isopropylbenzene, 1,1,2,2-tetrachloroethane, and 1,2-dibromo-3-chloropropane, the low MSD accuracy result for 1,2,4-trichlorobenzene, and the high precision result for methylcyclohexane during the spiked analyses of sample SB-6 (8-10); and the high MS accuracy result for isopropylbenzene during the spiked analyses of sample MW-1 (10-15). Validation qualification of the parent samples for these compounds was not required.

### LCS Recoveries

All LCS recoveries associated with project samples were considered acceptable and within QC limits with the exception of the low LCS recovery for 1,2-dichlorobenzene (80%R; QC limit 82-118%R) associated with samples SB-7 (8-10), SB-7 (16-18), SB-6 (8-10), and SB-4 (19-20); the high LCS recoveries for methylene chloride (135%R; QC limit 73-134%R) and 1,1,2,2-tetrachloroethane (125%R; QC limit 79-124%R) associated with samples SB-6 (10-15), SB-17 (8-10), and MW-1 (10-15); and the high LCS recovery for cis-1,2-dichloroethene (142%R; QC limit 78-122%R) associated with samples SB-2 (5-10), SB-2 (10-15), SB-3 (5-10), SB-3 (12-14), SB-1 (10-12), SB-1 (15-17), SB-5 (5-10), and SB-5 (10-15). Positive results for those compounds where LCS recoveries exceeded the QC limits were considered estimated, possibly biased high, and qualified "J" for the affected samples. Results for those compounds where LCS recoveries fell below the QC limits were considered estimated, possibly biased low, with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

### Blank Contamination

The laboratory method blank associated with samples SB-6 (10-15), SB-17 (8-10), and MW-1 (10-15) contained methylene chloride at a concentration of 4.6 µg/kg; and the laboratory method blank associated with sample MW-11 (5-10) contained methylene chloride at a concentration of 2 µg/kg. Therefore, methylene chloride results less than the validation action concentrations were considered not detected and qualified "U" for the affected samples.

### Initial and Continuing Calibrations

All initial calibration compounds were compliant with a minimum average relative response factor (RRF) of 0.05 and a maximum percent relative standard deviation (%RSD) of 20% with the exception of 1,2-dibromo-3-chloropropane (21.2%RSD) in the initial calibration associated with samples FB110113, TB110113, FB110413, and TB110413; carbon tetrachloride (63.5%RSD), 4-methyl-2-pentanone (24.9%RSD), ethylbenzene (29%RSD), bromoform (22.6%RSD), isopropylbenzene (27.9%RSD), and 1,2,4-trichlorobenzene (21.4%RSD) in the initial calibration associated with sample SB-4 (15-17); and chloroethane (20.6%RSD) and trichlorofluoromethane (23.3%RSD) in the initial calibration associated with samples SB-2 (5-10), SB-2 (10-15), SB-3 (5-10), SB-3 (12-14), SB-1 (10-12), SB-1 (15-17), SB-5 (5-10), SB-5 (10-15), MW-11 (5-10), and MW-1 (5-10). The results for these compounds were considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

All continuing calibration compounds were compliant with a minimum RRF of 0.05 and a maximum percent difference (%D) within  $\pm 20\%$  with the exception of methyl acetate (20.49%D), 1,2-dichloropropane (21.76%D), bromoform (25.91%D), 1,1,2,2-tetrachloroethane

(24.33%D), 1,2,4-trichlorobenzene (20.44%D), 1,2,3-trichlorobenzene (20.45%D), and 1,4-dioxane (100%D) in the continuing calibration associated with samples SB-11 (5-10), SB-11 (15-20), SB-10 (5-10), SB-10 (10-12), SB-9 (5-10), SB-9 (10-15), SB-8 (15-16), and SB-8 (17-19); bromochloromethane (21.65%D, -21.65%D) in the continuing calibrations associated with samples SB-7 (8-10), SB-7 (16-18), SB-6 (8-10), SB-6 (10-15), SB-17 (8-10), SB-4(19-20), and MW-1 (10-15); trichlorofluoromethane (20.01%D) and cis-1,2-dichloroethene (37.56%D) in the continuing calibration associated with samples SB-2 (5-10), SB-2 (10-15), SB-3 (5-10), SB-3 (12-14), SB-1 (10-12), SB-1 (15-17), SB-5 (5-10), and SB-5 (10-15); acetone (-35.16%D), 2-butanone (22.81%D), carbon tetrachloride (-27.52%D), 2-hexanone (23.91%D), 1,2,4-trichlorobenzene (28.6%D), and 1,4-dioxane (50%D) in the continuing calibration associated with sample SB-4 (15-17); vinyl chloride (-20.15%D), chloroethane (-20.55%D), 1,1-dichloroethane (33.9%D), 2-hexanone (33.49%D), dibromochloromethane (25.96%D), 1,2-dibromoethane (23.02%D), and bromoform (37.04%D) in the continuing calibration associated with samples FB110113, TB110113, FB110413, and TB110413; chloroethane (30.36%D) and trichlorofluoromethane (37.78%D) in the continuing calibration associated with sample MW-11 (5-10); and chloroethane (28.19%D) and trichlorofluoromethane (38.49%D) in the continuing calibration associated with sample MW-1 (5-10). Therefore, the sample results for these compounds were considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples. However, nondetected sample results for those compounds where the %D was greater than 90% were considered unusable and qualified "R" for the affected samples.

#### Internal Standard Responses

All internal standard (IS) responses and retention times were within specified QC ranges based on associated calibration standards (i.e., sample's area count within -50% to +100% and retention times within  $\pm 0.5$  minutes of the standard) with the exception of the low IS response for 1,4-dichlorobenzene-d4 in samples SB-11 (5-10), SB-10 (5-10), SB-7 (8-10), SB-7 (16-18), SB-6 (8-10), SB-2 (5-10), SB-2 (10-15), and SB-5 (10-15). These samples were reanalyzed yielding similar IS responses. Therefore, sample results associated with this IS were considered estimated, possibly biased low, with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

#### Usability

All volatile soil sample results were considered usable following data validation with the exception of certain nondetected results based upon poor instrument calibration linearity for 1,4-dioxane.

#### Summary

The quality assurance objectives for measurement data included considerations for precision, accuracy, representativeness, completeness, and comparability. The volatile soil data presented by Chemtech were 99.5% complete (i.e., usable). The validated volatile laboratory data are tabulated and presented in Attachment A-1.

## 2.1.2 Semivolatiles

The following items were reviewed for compliancy in the semivolatile analysis:

- Custody documentation
- Holding times
- Surrogate recoveries
- MS/MSD precision and accuracy
- LCS recoveries
- Laboratory method blank and field equipment blank contamination
- GC/MS instrument performance
- Initial and continuing calibrations
- Internal standard area counts and retention times
- Field duplicate precision
- Sample result verification and identification
- Quantitation limits
- Data completeness

These items were considered compliant and acceptable in accordance with the validation protocols with the exception of surrogate recoveries, MS/MSD precision and accuracy, LCS recoveries, initial and continuing calibrations, and field duplicate precision as discussed below.

### Surrogate Recoveries

All sample surrogate recoveries were considered acceptable and within QC limits with the exception of the low recovery for the surrogate phenol-d6 (QC limit 34-127%R) in sample SB-7 (8-10) (31%R). Validation qualification of this sample was not required since only one acid surrogate fell below the QC limit.

### MS/MSD Precision and Accuracy

All MS/MSD precision (relative percent difference; RPD) and accuracy (percent recovery; %R) measurements were considered acceptable and within QC limits for designated spiked project samples with the exception of the high MSD accuracy results for fluoranthene and benzo(b)fluoranthene, the low MSD accuracy result for hexachlorocyclopentadiene, the less than 10% MS/MSD accuracy results for benzaldehyde, and the many high precision results during the spiked analyses of sample SB-6(8-10); and the 0% MS/MSD accuracy results for benzaldehyde and the high precision results for 2,4-dinitrophenol and 4,6-dinitro-2-methylphenol during the spiked analyses of sample MW-1 (5-10). Validation qualification of these compounds was not required for the parent samples with the exception of benzaldehyde. The nondetected benzaldehyde results were considered unusable and qualified “R” for the parent samples.

## LCS Recoveries

All LCS recoveries were considered acceptable and within QC limits with the exception of the high LCS recoveries for hexachlorocyclopentadiene (136%R; QC limit 38-122%R), carbazole (106%R; QC limit 57-102%R), di-n-butylphthalate (106%R; QC limit 57-103%R), benzo(b)fluoranthene (118%R; QC limit 56-103%R), benzo(g,h,i)perylene (106%R; QC limit 56-105%R) associated with samples SB-11 (5-10), SB-11 (15-20), SB-10 (5-10), SB-10 (10-12), SB-9 (5-10), SB-9 (10-15), SB-8 (15-16), and SB-8 (17-19); the less than 10% LCS recovery for benzaldehyde associated with all samples; and the high LCS recoveries for hexachlorocyclopentadiene (127%R, 124%R; QC limit 38-122%R) associated with all soil samples except SB-11 (5-10), SB-11 (15-20), SB-10 (5-10), SB-10 (10-12), SB-9 (5-10), SB-9 (10-15), SB-8 (15-16), and SB-8 (17-19). Therefore, results for those compounds where LCS recoveries fell below the QC limit were considered estimated, possibly biased low, with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples. However, the benzaldehyde results which were nondetects were considered unusable and qualified "R" for the affected samples. Positive results for those samples where LCS recoveries exceeded the QC limit were considered estimated, possibly biased high, and qualified "J" for the affected samples.

## Initial and Continuing Calibrations

All initial calibration compounds were compliant with a minimum relative response factor (RRF) of 0.05 and a maximum percent relative standard deviation (%RSD) of 20% with the exception of 2,4-dinitrophenol (30.6%RSD) in the initial calibration associated with samples SB-11 (5-10), SB-11 (15-20), SB-10 (5-10), SB-10 (10-12), SB-9 (5-10), SB-9 (10-15), SB-8 (15-16), SB-8 (17-19), SB-1 (15-17), FB110113, and FB110413; and 4,6-dinitro-2-methylphenol (26%RSD) in the initial calibration associated with soil samples in SDGs E4277 and E4340 except SB-11 (5-10), SB-11 (15-20), SB-10 (5-10), SB-10 (10-12), SB-9 (5-10), SB-9 (10-15), SB-8 (15-16), SB-8 (17-19), SB-1 (15-17), and FB110113. The results for these compounds were considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

All continuing calibration compounds were compliant with a minimum RRF of 0.05 and a maximum percent difference (%D) within  $\pm 20\%$  with the exception of 2,4-dinitrophenol (54.1%D), 4,6-dinitro-2-methylphenol (21.5%D), atrazine (22.4%D), benzo(b)fluoranthene (20.7%D), and benzo(k)fluoranthene (-21.2%D) in the continuing calibration associated with sample SB-8 (17-19); nitrobenzene (-20.3%D) and pentachlorophenol (-23.4%D) in the continuing calibration associated with samples FB110113 and SB-1 (15-17); 2,4-dinitrophenol (36.1%D) and 4,6-dinitro-2-methylphenol (35.4%D) in the continuing calibration associated with samples SB-7 (8-10), SB-7 (16-18), SB-6 (8-10), SB-4 (15-17), and MW-1 (10-15); phenol (-27.8%D), hexachlorocyclopentadiene (22.6%D), 2,4-dinitrophenol (57.4%D), 4,6-dinitro-2-methylphenol (58.2%D), bis(2-ethylhexyl)phthalate (22.4%D), and di-n-octylphthalate (29.2%D) in the continuing calibration associated with samples SB-6 (10-15), SB-17 (8-10), SB-5 (5-10), SB-5 (10-15), SB-4 (19-20), SB-2 (5-10), SB-2 (10-15), SB-3 (5-10), SB-3 (12-14), and SB-1 (10-12); phenol (-20.7%D), 2-nitroaniline (20.3%D), 2,4-dinitrophenol (80.3%D), 4-nitrophenol (26.1%D), 2,4-dinitrotoluene (23.4%D), 4-nitroaniline (20.7%D), 4,6-dinitro-2-methylphenol (68.4%D), and benzo(k)fluoranthene (20.7%D) in the continuing calibration associated with samples MW-1 (5-10) and MW-11 (5-10); and benzaldehyde (-20.2%D),

hexachlorocyclopentadiene (-26.2%D), 2,4-dinitrophenol (-22.4%D), and benzo(k)fluoranthene (23.5%D) in the continuing calibration associated with sample FB110413 . Therefore, results for these compounds were considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

#### Field Duplicate Precision

All field duplicate precision results were considered acceptable with the exception of the precision results for dimethylphthalate, anthracene, benzo(k)fluoranthene, and benzo(g,h,i)perylene for the field duplicate pair SB-7 (8-10) and SB-17 (8-10). Therefore, the results for these compounds were considered estimated and qualified "J" for these samples.

#### Usability

All semivolatile soil sample results were considered usable following data validation with the exception of certain nondetected results based upon poor MS/MSD and LCS recoveries.

#### Summary

The quality assurance objectives for measurement data included considerations for precision, accuracy, representativeness, completeness, and comparability. The semivolatile soil data presented by Chemtech were 98.6% complete (i.e., usable). The validated semivolatile laboratory data are tabulated and presented in Attachment A-1.

### **2.1.3 Inorganics**

The following items were reviewed for compliancy in the inorganics analysis:

- Custody documentation
- Holding times
- Initial and continuing calibration verifications
- Initial and continuing calibration blank, and laboratory preparation blank, and field equipment blank contamination
- Inductively coupled plasma (ICP) interference check sample (ICS)
- Matrix spike (MS) recoveries
- Laboratory duplicate precision
- Field duplicate precision
- Laboratory control sample (LCS) recoveries
- ICP serial dilutions
- Sample result verification and identification
- Quantitation limits
- Data completeness

These items were considered compliant and acceptable in accordance with the validation protocols with the exception of blank contamination, ICS recoveries, matrix spike recoveries, serial dilutions, and field duplicate precision as discussed below.

#### Blank Contamination

The field equipment blank FB110113 associated with samples in SDG E4277 contained calcium and sodium at concentrations of 86.5 and 41.3 µg/L, respectively. Since sample results were not affected by the contamination in this blank, validation qualification was not required.

#### ICS Recoveries

All ICS recoveries were considered acceptable and within the 80-120%R QC limit for all analytes with the exception of the low ICS recovery for barium (72.5%R) and the high ICS recoveries for cobalt (125.5%R, 130.3%R) associated with samples in SDG E4340. Therefore, barium results were considered estimated, possibly biased low, with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples. Positive cobalt results were considered estimated, possibly biased high, and qualified "J" for the affected samples.

#### Matrix Spike Recoveries

All the MS recoveries for designated spiked project samples were within the 75-125%R QC limit with sample concentrations less than four times the spiking concentration with the exception of the low MS recoveries for antimony (26%R) and lead (24%R, 22%R) and the high MS recoveries for nickel (146%R, 144%R) and potassium (186%R) associated with all soil samples in SDG E4277; and the low MS recoveries for antimony (21.1%R, 21.8%R) associated with soil samples in SDG E4340. Therefore, positive results for those analytes where MS recoveries exceeded the QC limit were considered estimated, possibly biased high, and qualified "J" for the affected samples. Results for those analytes where MS recoveries fell below the QC limit were considered estimated, possibly biased low, with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

#### ICP Serial Dilutions

All serial dilution results for designated project samples were considered acceptable with a percent difference (%D) less than 10% for all ICP analytes with the exception of aluminum (12%D), barium (19%D), calcium (23%D), chromium (19%D), copper (30%D), magnesium (19%D), manganese (29%D), potassium (18%D), sodium (18%D), and vanadium (21%D) associated with soil samples in SDG E4340. Therefore, positive results for these analytes were considered estimated and qualified "J" for the affected samples.

#### Field Duplicate Precision

All field duplicate precision results were considered acceptable with the exception of the precision results for cyanide for the field duplicate pair SB-7 (8-10) and SB-17 (8-10). Therefore, the cyanide results for these samples were considered estimated with the positive result qualified "J" and the nondetected result qualified "UJ".

## Usability

All inorganics soil sample results were considered usable following data validation

## Summary

The quality assurance objectives for measurement data included considerations for precision, accuracy, representativeness, completeness, and comparability. The inorganics soil data presented by Chemtech were 100% complete (i.e., usable). The validated soil inorganics laboratory data are tabulated and presented in Attachment A-1.

## **2.2 GROUNDWATER**

Data review has been completed for data packages generated by Chemtech containing groundwater samples collected from the site. All of these samples were properly preserved, shipped under a COC record, and received intact by the analytical laboratory. The analytical results were presented by the laboratory in one sample delivery group (SDG): E4638. Data validation was performed for all samples in accordance with the most current editions of the USEPA Region II SOPs for organic and inorganic data review. This data validation and usability report is presented by analysis type and the validated laboratory data are presented in Attachment A-2.

### **2.2.1 Volatiles**

The following items were reviewed for compliancy in the volatile analysis:

- Custody documentation
- Holding times
- Surrogate recoveries
- Matrix spike/matrix spike duplicate (MS/MSD) precision and accuracy
- Laboratory control sample (LCS) recoveries
- Laboratory method blank and field equipment/trip blank contamination
- GC/MS instrument performance
- Initial and continuing calibrations
- Internal standard area counts and retention times
- Field duplicate precision
- Sample result verification and identification
- Quantitation limits
- Data completeness

These items were considered compliant and acceptable in accordance with the validation protocols with the exception of MS/MSD precision and accuracy, LCS recoveries, blank contamination, and initial and continuing calibrations as discussed below.

## MS/MSD Precision and Accuracy

All MS/MSD precision (relative percent difference; RPD) and accuracy (percent recovery; %R) measurements were considered acceptable and within QC limits for designated project spiked samples with the exception of the high precision results for carbon tetrachloride, 1,1,2-trichloroethane, 2-hexanone, tetrachloroethene, isopropylbenzene, 1,1,2,2-tetrachloroethane, and 1,3-dichlorobenzene during the spiked analyses of sample MW-3. Validation qualification of the parent sample was not required.

## LCS Recoveries

All LCS recoveries were considered acceptable and within QC limits for all compounds with the exception of the high LCS recovery for bromochloromethane (150%R; 65-130%R) associated with samples MW-1, -2, and -3. Validation qualification was not required since this compound was not detected in these samples.

## Blank Contamination

The field equipment blank FB112713 associated with groundwater samples contained methylene chloride at a concentration of 1.1 µg/L. Since this compound was not detected in the project samples, validation qualification was not required.

## Initial and Continuing Calibrations

All initial calibration compounds were compliant with a minimum relative response factor (RRF) of 0.05 and a maximum percent relative standard deviation (%RSD) of 20% with the exception of cyclohexane (40.7%RSD) in the initial calibration associated with all groundwater samples. The results for this compound were considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

All continuing calibration compounds were compliant with a minimum RRF of 0.05 and a maximum percent difference (%D) within  $\pm 20\%$  with the exception of cyclohexane (40.7%D) in the continuing calibration associated with all groundwater samples. Therefore, the sample results for this compound were considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

## Usability

All volatile groundwater sample results were considered usable following data validation.

## Summary

The quality assurance objectives for measurement data included considerations for precision, accuracy, representativeness, completeness, and comparability. The volatile groundwater data presented by Chemtech were 100% complete (i.e., usable). The validated volatile laboratory data are tabulated and presented in Attachment A-2.

## **2.2.2 Semivolatiles**

The following items were reviewed for compliancy in the semivolatile analysis:

- Custody documentation
- Holding times
- Surrogate recoveries
- MS/MSD precision and accuracy
- LCS recoveries
- Laboratory method blank and field equipment blank contamination
- GC/MS instrument performance
- Initial and continuing calibrations
- Internal standard area counts and retention times
- Field duplicate precision
- Sample result verification and identification
- Quantitation limits
- Data completeness

These items were considered compliant and acceptable in accordance with the validation protocols with the exception of MS/MSD precision and accuracy, LCS recoveries, and initial and continuing calibrations as discussed below.

### MS/MSD Precision and Accuracy

All MS/MSD precision (relative percent difference; RPD) and accuracy (percent recovery; %R) measurements were considered acceptable and within QC limits for designated spiked project samples with the exception of the low MSD accuracy result for 2,3,4,6-tetrachlorophenol during the spiked analyses of sample MW-3. Validation qualification of the parent sample was not required.

### LCS Recoveries

All LCS recoveries were considered acceptable and within QC limits with the exception of the high LCS recoveries for 2-methylphenol (99%R; QC limit 32-94%R), 3+4-methylphenols (95%R; QC limit 24-91%R), and hexachlorocyclopentadiene (150%R; QC limit 42-121%R) associated with all groundwater samples. Validation qualification of the groundwater samples was not required since these compounds were not detected.

### Initial and Continuing Calibrations

All initial calibration compounds were compliant with a minimum relative response factor (RRF) of 0.05 and a maximum percent relative standard deviation (%RSD) of 20% with the exception of 2,4-dinitrophenol (29.7%RSD) in the initial calibration associated with all

groundwater samples. The results for this compound were considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

All continuing calibration compounds were compliant with a minimum RRF of 0.05 and a maximum percent difference (%D) within  $\pm 20\%$  with hexachlorocyclopentadiene (39.2%D), 2-nitroaniline (24.3%D), 2,6-dinitrotoluene (20.6%D), 4-nitroaniline (24.9%D), bis(2-ethylhexyl)phthalate (22%D), and di-n-octylphthalate (27.8%D) in the continuing calibration associated with all groundwater samples. Therefore, the sample results for these compounds were considered estimated with positive results qualified "J" and nondetected results qualified "UJ" for the affected samples.

### Usability

All semivolatile groundwater sample results were considered usable following data validation.

### Summary

The quality assurance objectives for measurement data included considerations for precision, accuracy, representativeness, completeness, and comparability. The semivolatile groundwater data presented by Chemtech were 100% complete (i.e., usable). The validated semivolatile laboratory data are tabulated and presented in Attachment A-2.

## **2.2.3 Inorganics**

The following items were reviewed for compliancy in the inorganics analysis:

- Custody documentation
- Holding times
- Initial and continuing calibration verifications
- Initial and continuing calibration blank, and laboratory preparation blank, and field equipment blank contamination
- Inductively coupled plasma (ICP) interference check sample (ICS)
- Matrix spike (MS) recoveries
- Laboratory duplicate precision
- Field duplicate precision
- Laboratory control sample (LCS) recoveries
- ICP serial dilutions
- Sample result verification and identification
- Quantitation limits
- Data completeness

These items were considered compliant and acceptable in accordance with the validation protocols with the exception of blank contamination and field duplicate precision as discussed below.

#### Blank Contamination

The field equipment blank FB112713 associated with the groundwater samples contained calcium at a concentration of 1200 µg/L. Validation qualification was not required since samples were not affected by the contamination in this blank.

#### Field Duplicate Precision

All field duplicate precision results were considered acceptable for the field duplicate pair MW-3 and MW-3A with the exception of the precision for chromium. Therefore, the chromium results were considered estimated with the positive result qualified "J" and the nondetected result qualified "UJ" for the affected parent sample and its field duplicate.

#### Usability

All inorganics groundwater sample results were considered usable following data validation.

#### Summary

The quality assurance objectives for measurement data included considerations for precision, accuracy, representativeness, completeness, and comparability. The inorganics groundwater data presented by Chemtech were 100% complete (i.e., usable). The validated groundwater inorganics laboratory data are tabulated and presented in Attachment A-2.

**ATTACHMENT A**

**VALIDATED LABORATORY DATA**

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**PARSONS**

**ATTACHMENT A-1**

**VALIDATED LABORATORY DATA FOR SOIL SAMPLES**

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**PARSONS**

Dup of MW-1(5-10)-20131104						
Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	MW-1 MW-1(5-10)-20131104 E4340-01 5 - 10 ft CTECH E4340 SOIL 11/4/2013 10:50 2/4/2014	MW-1 MW-1(5-10)-20131104 E4340-05 5 - 10 ft CTECH E4340 SOIL 11/4/2013 10:40 2/4/2014	MW-1 MW-1(10-15)-20131104 E4340-02 10 - 15 ft CTECH E4340 SOIL 11/4/2013 11:00 2/4/2014	SB-1 SB-1(10-12)-20131101 E4277-25 10 - 12 ft CTECH E4277 SOIL 11/1/2013 13:45 2/4/2014
CAS NO.	COMPOUND	UNITS:				
	VOLATILES					
71-55-6	1,1,1-TRICHLOROETHANE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/kg	0.82 U	0.72 U	0.71 U	0.43 U
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/kg	1.6 U	1.4 U	1.4 U	0.84 U
75-34-3	1,1-DICHLOROETHANE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
75-35-4	1,1-DICHLOROETHENE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
120-82-1	1,2,4-TRICHLOROBENZENE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/kg	1.6 U	1.4 U	1.4 U	0.81 U
106-93-4	1,2-DIBromoETHANE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
95-50-1	1,2-DICHLOROBENZENE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
107-06-2	1,2-DICHLOROETHANE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
78-87-5	1,2-DICHLOROPROPANE	ug/kg	0.46 U	0.4 U	0.4 U	0.24 U
541-73-1	1,3-DICHLOROBENZENE	ug/kg	0.66 U	0.58 U	0.57 U	0.35 U
106-46-7	1,4-DICHLOROBENZENE	ug/kg	0.73 U	0.64 U	0.64 U	0.38 U
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/kg	180 U	160 U	160 U	93.3 U
591-78-6	2-HEXANONE	ug/kg	4.5 U	3.9 U	3.9 U	2.3 U
67-64-1	ACETONE	ug/kg	7.7 J	3.9 U	16 J	2.3 U
71-43-2	BENZENE	ug/kg	0.68 U	0.59 U	0.59 U	0.35 U
74-97-5	BROMOCHLOROMETHANE	ug/kg	0.89 U	0.78 U	0.78 UJ	0.47 U
75-27-4	BROMODICHLOROMETHANE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
75-25-2	BROMOFORM	ug/kg	1.3 U	1.2 U	1.1 U	0.69 U
74-83-9	BROMOMETHANE	ug/kg	1.8 U	1.6 U	1.6 U	0.93 U
75-15-0	CARBON DISULFIDE	ug/kg	0.89 U	0.78 U	2.5 J	0.47 U
56-23-5	CARBON TETRACHLORIDE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
108-90-7	CHLOROBENZENE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
75-00-3	CHLOROETHANE	ug/kg	0.89 UJ	0.78 UJ	0.78 U	0.47 UJ
67-66-3	CHLOROFORM	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
74-87-3	CHLOROMETHANE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 UJ
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
110-82-7	CYCLOHEXANE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
124-48-1	DIBROMOCHLOROMETHANE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
75-71-8	DICHLORODIFLUOROMETHANE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
100-41-4	ETHYLBENZENE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/kg	0.86 U	0.75 U	0.75 U	0.45 U
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/kg	1.3 U	1.1 U	1.1 U	0.67 U
79-20-9	METHYL ACETATE	ug/kg	1.8 U	1.6 U	1.6 U	0.93 U
78-93-3	METHYL ETHYL KETONE	ug/kg	5.5 U	4.8 U	4.8 U	2.9 U
108-10-1	METHYL ISOBUTYL KETONE	ug/kg	4.5 U	3.9 U	3.9 U	2.3 U
108-87-2	METHYLCYCLOHEXANE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
75-09-2	METHYLENE CHLORIDE	ug/kg	6.8 J	7.8 U	9.4 UJ	3.7 J
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
100-42-5	STYRENE	ug/kg	0.8 U	0.7 U	0.7 U	0.42 U
1634-04-4	TERT-BUTYL METHYL ETHER	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
108-88-3	TOLUENE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U
75-69-4	TRICHLOROFUOROMETHANE	ug/kg	0.89 UJ	0.78 UJ	0.78 U	0.47 UJ
75-01-4	VINYL CHLORIDE	ug/kg	0.89 U	0.78 U	0.78 U	0.47 U

Dup of MW-1(5-10)-20131104						
Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	MW-1 MW-1(5-10)-20131104 E4340-01 5 - 10 ft CTECH E4340 SOIL 11/4/2013 10:50 2/4/2014	MW-1 MW-1(5-10)-20131104 E4340-05 5 - 10 ft CTECH E4340 SOIL 11/4/2013 10:40 2/4/2014	MW-1 MW-1(10-15)-20131104 E4340-02 10 - 15 ft CTECH E4340 SOIL 11/4/2013 11:00 2/4/2014	SB-1 SB-1(10-12)-20131101 E4277-25 10 - 12 ft CTECH E4277 SOIL 11/1/2013 13:45 2/4/2014
CAS NO.	COMPOUND	UNITS:				
	SEMIVOLATILES					
95-94-3	1,2,4,5-TETRACHLOROBENZENE	ug/kg	20.2 U	17.7 U	19.5 U	15.2 U
58-90-2	2,3,4,6-TETRACHLOROPHENOL	ug/kg	20.2 U	17.7 U	19.5 U	15.2 U
95-95-4	2,4,5-TRICHLOROPHENOL	ug/kg	36.1 U	31.6 U	34.8 U	27.2 U
88-06-2	2,4,6-TRICHLOROPHENOL	ug/kg	15.7 U	13.8 U	15.2 U	11.9 U
120-83-2	2,4-DICHLOROPHENOL	ug/kg	19.6 U	17.2 U	18.9 U	14.8 U
105-67-9	2,4-DIMETHYLPHENOL	ug/kg	29.2 U	25.5 U	28.1 U	22 U
51-28-5	2,4-DINITROPHENOL	ug/kg	52.3 UJ	45.8 UJ	50.5 UJ	39.4 UJ
121-14-2	2,4-DINITROTOLUENE	ug/kg	15.4 UJ	13.5 UJ	14.9 U	11.6 U
606-20-2	2,6-DINITROTOLUENE	ug/kg	21 U	18.4 U	20.2 U	15.8 U
91-58-7	2-CHLORONAPHTHALENE	ug/kg	11.7 U	10.3 U	11.3 U	8.8 U
95-57-8	2-CHLOROPHENOL	ug/kg	27.2 U	23.8 U	26.2 U	20.5 U
91-57-6	2-METHYLNAPHTHALENE	ug/kg	13 U	140 J	12.5 U	9.8 U
95-48-7	2-METHYLPHENOL (O-CRESOL)	ug/kg	27.9 U	24.4 U	26.9 U	21 U
88-74-4	2-NITROANILINE	ug/kg	22.8 UJ	20 UJ	22 U	17.2 U
88-75-5	2-NITROPHENOL	ug/kg	24.8 U	21.7 U	24 U	18.7 U
MEPH3ME	3- AND 4- METHYLPHENOL (TOTAL)	ug/kg	26.7 U	23.4 U	25.7 U	20.1 U
91-94-1	3,3'-DICHLOROBENZIDINE	ug/kg	33 U	28.9 U	31.9 U	24.9 U
99-09-2	3-NITROANILINE	ug/kg	33 U	28.9 U	31.9 U	24.9 U
534-52-1	4,6-DINITRO-2-METHYLPHENOL	ug/kg	29.5 UJ	25.8 UJ	28.4 UJ	22.2 UJ
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/kg	10 U	8.8 U	9.7 U	7.6 U
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/kg	22.8 U	20 U	22 U	17.2 U
106-47-8	4-CHLOROANILINE	ug/kg	36.3 U	31.7 U	35 U	27.3 U
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/kg	27.9 U	24.4 U	26.9 U	21 U
100-01-6	4-NITROANILINE	ug/kg	67 UJ	58.6 UJ	64.6 U	50.4 U
100-02-7	4-NITROPHENOL	ug/kg	95.5 UJ	83.6 UJ	92.1 U	72 U
83-32-9	ACENAPHTHENE	ug/kg	14.5 U	90.5 J	14 U	10.9 U
208-96-8	ACENAPHTHYLENE	ug/kg	240 J	260 J	12.5 U	9.8 U
98-86-2	ACETOPHENONE	ug/kg	15.7 U	13.8 U	15.2 U	11.9 U
120-12-7	ANTHRACENE	ug/kg	190 J	240 J	120 J	7.9 U
1912-24-9	ATRAZINE	ug/kg	27.2 U	23.8 U	26.2 U	20.5 U
100-52-7	BENZALDEHYDE	ug/kg	R	R	R	R
56-55-3	BENZO(A)ANTHRACENE	ug/kg	360 J	560	340 J	18.5 U
50-32-8	BENZO(A)PYRENE	ug/kg	440 J	630	290 J	8.4 U
205-99-2	BENZO(B)FLUORANTHENE	ug/kg	480 J	720	360 J	12.7 U
191-24-2	BENZO(G,H,I)PERYLENE	ug/kg	300 J	420 J	190 J	15.7 U
207-08-9	BENZO(K)FLUORANTHENE	ug/kg	150 J	240 J	110 J	18.2 U
85-68-7	BENZYL BUTYL PHTHALATE	ug/kg	24.7 U	21.6 U	23.8 U	18.6 U
92-52-4	BIPHENYL (DIPHENYL)	ug/kg	19.4 U	17 U	18.8 U	14.6 U
111-91-1	BIS(2-CHLOROETHOXY) METHANE	ug/kg	29.6 U	25.9 U	28.6 U	22.3 U
111-44-4	BIS(2-CHLOROETHYL) ETHER	ug/kg	24.7 U	21.6 U	23.8 U	18.6 U
108-60-1	BIS(2-CHLOROISOPROPYL) ETHER	ug/kg	21.3 U	18.6 U	20.5 U	16 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/kg	18.2 U	15.9 U	17.6 U	13.7 UJ
105-60-2	CAPROLACTAM	ug/kg	23.9 U	20.9 U	23.1 U	18 U
86-74-8	CARBAZOLE	ug/kg	11.3 U	9.9 U	10.9 U	8.5 U
218-01-9	CHRYSENE	ug/kg	400 J	630	260 J	17.6 U
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/kg	14.8 U	110 J	14.3 U	11.2 U
132-64-9	DIBENZOFURAN	ug/kg	20.1 U	17.6 U	19.3 U	15.1 U
84-66-2	DIETHYL PHTHALATE	ug/kg	8 U	7 U	7.7 U	6 U
131-11-3	DIMETHYL PHTHALATE	ug/kg	610	800	530	480
84-74-2	DI-N-BUTYL PHTHALATE	ug/kg	40.4 U	35.4 U	39 U	30.5 U
117-84-0	DI-N-OCTYLPHTHALATE	ug/kg	5.9 U	5.1 U	5.7 U	4.4 UJ
206-44-0	FLUORANTHENE	ug/kg	650	990	520	7.8 U
86-73-7	FLUORENE	ug/kg	100 J	120 J	18.8 U	14.6 U
118-74-1	HEXACHLOROBENZENE	ug/kg	21 U	18.4 U	20.2 U	15.8 U
87-68-3	HEXACHLOROBUTADIENE	ug/kg	18.7 U	16.3 U	18 U	14.1 U
77-47-4	HEXACHLOROCYCLOPENTADIENE	ug/kg	12.5 U	10.9 U	12.1 U	9.4 UJ
67-72-1	HEXACHLOROETHANE	ug/kg	23 U	20.1 U	22.2 U	17.3 U
193-39-5	INDENO(1,2,3-C,D)PYRENE	ug/kg	210 J	330 J	170 J	12.9 U
78-59-1	ISOPHORONE	ug/kg	17 U	14.9 U	16.4 U	12.8 U

Dup of MW-1(5-10)-20131104						
Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	MW-1 MW-1(5-10)-20131104 E4340-01 5 - 10 ft CTECH E4340 SOIL 11/4/2013 10:50 2/4/2014	MW-1 MW-1(5-10)-20131104 E4340-05 5 - 10 ft CTECH E4340 SOIL 11/4/2013 10:40 2/4/2014	MW-1 MW-1(10-15)-20131104 E4340-02 10 - 15 ft CTECH E4340 SOIL 11/4/2013 11:00 2/4/2014	SB-1 SB-1(10-12)-20131101 E4277-25 10 - 12 ft CTECH E4277 SOIL 11/1/2013 13:45 2/4/2014
CAS NO.	COMPOUND	UNITS:				
	SEMIVOLATILES					
91-20-3	NAPHTHALENE	ug/kg	230 J	300 J	17.1 U	13.4 U
98-95-3	NITROBENZENE	ug/kg	19.4 U	17 U	18.8 U	14.6 U
621-64-7	N-NITROSODI-N-PROPYLAMINE	ug/kg	25.9 U	22.7 U	25 U	19.5 U
86-30-6	N-NITROSODIPHENYLAMINE	ug/kg	12.3 U	10.8 U	11.9 U	9.3 U
87-86-5	PENTACHLOROPHENOL	ug/kg	35.2 U	30.8 U	33.9 U	26.5 U
85-01-8	PHENANTHRENE	ug/kg	450 J	620	320 J	10.5 U
108-95-2	PHENOL	ug/kg	11.9 UJ	10.4 UJ	11.5 U	9 UJ
129-00-0	PYRENE	ug/kg	690	1000	540	9.3 U
	INORGANICS					
7429-90-5	ALUMINUM	mg/kg	14900 J	13800 J	13000 J	12000
7440-36-0	ANTIMONY	mg/kg	0.751 UJ	0.626 UJ	0.707 UJ	0.543 UJ
7440-38-2	ARSENIC	mg/kg	10.8	6.94	8.22	1.48
7440-39-3	BARIUM	mg/kg	134 J	108 J	80.4 J	71
7440-41-7	BERYLLIUM	mg/kg	0.435	0.361	0.52	0.365
7440-43-9	CADMIUM	mg/kg	2.6	2.18	0.076 U	0.058 U
7440-70-2	CALCIUM	mg/kg	4200 J	4200 J	4300 J	8100
7440-47-3	CHROMIUM, TOTAL	mg/kg	36.1 J	30.1 J	28.3 J	25.1
7440-48-4	COBALT	mg/kg	12.4 J	10.47 J	14.1 J	13.2
7440-50-8	COPPER	mg/kg	99.3 J	53.8 J	49.1 J	21.1
7439-89-6	IRON	mg/kg	36900	33900	34700	22600
7439-92-1	LEAD	mg/kg	493	352	390	17.3 J
7439-95-4	MAGNESIUM	mg/kg	5500 J	5000 J	6800 J	5600
7439-96-5	MANGANESE	mg/kg	323 J	270 J	641 J	440
7439-97-6	MERCURY	mg/kg	2.33	1.54	0.268	0.039
7440-02-0	NICKEL	mg/kg	29.4	25.3	29.6	21.6 J
7440-09-7	POTASSIUM	mg/kg	2900 J	2700 J	3300 J	2000 J
7782-49-2	SELENIUM	mg/kg	2.56	2.62	2.74	1.36
7440-22-4	SILVER	mg/kg	1.43	1.18	1.82	1.3
7440-23-5	SODIUM	mg/kg	271 J	245 J	607 J	116
7440-28-0	THALLIUM	mg/kg	0.362 U	0.302 U	0.341 U	0.262 U
7440-62-2	VANADIUM	mg/kg	50.5 J	44.9 J	50.5 J	40
7440-66-6	ZINC	mg/kg	346	329	227	58.5
57-12-5	CYANIDE	mg/kg	0.649	0.672	0.244 J	0.093 J

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-1 SB-1(15-17)-20131101 E4277-26	SB-2 SB-2(5-10)-20131101 E4277-21	SB-2 SB-2(10-15)-20131101 E4277-22	SB-3 SB-3(5-10)-20131101 E4277-23
CAS NO.	COMPOUND	UNITS:				
	VOLATILES					
71-55-6	1,1,1-TRICHLOROETHANE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/kg	0.56 U	0.71 UJ	0.94 UJ	0.66 U
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/kg	1.1 U	1.4 U	1.8 U	1.3 U
75-34-3	1,1-DICHLOROETHANE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
75-35-4	1,1-DICHLOROETHENE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/kg	0.61 U	0.77 UJ	1 UJ	0.72 U
120-82-1	1,2,4-TRICHLOROBENZENE	ug/kg	0.61 U	0.77 UJ	1 UJ	0.72 U
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/kg	1.1 U	1.3 UJ	1.8 UJ	1.3 U
106-93-4	1,2-DIBromoETHANE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
95-50-1	1,2-DICHLOROBENZENE	ug/kg	0.61 U	0.77 UJ	1 UJ	0.72 U
107-06-2	1,2-DICHLOROETHANE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
78-87-5	1,2-DICHLOROPROPANE	ug/kg	0.32 U	0.4 U	0.53 U	0.37 U
541-73-1	1,3-DICHLOROBENZENE	ug/kg	0.45 U	0.57 UJ	0.75 UJ	0.53 U
106-46-7	1,4-DICHLOROBENZENE	ug/kg	0.5 U	0.64 UJ	0.83 UJ	0.59 U
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/kg	120 U	150 U	200 U	140 U
591-78-6	2-HEXANONE	ug/kg	3 U	3.9 U	5.1 U	3.6 U
67-64-1	ACETONE	ug/kg	3 U	3.9 U	5.1 U	3.6 U
71-43-2	BENZENE	ug/kg	0.46 U	0.59 U	0.77 U	0.55 U
74-97-5	BROMOCHLOROMETHANE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
75-27-4	BROMODICHLOROMETHANE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
75-25-2	BROMOFORM	ug/kg	0.9 U	1.1 U	1.5 U	1.1 U
74-83-9	BROMOMETHANE	ug/kg	1.2 U	1.5 U	2 U	1.4 U
75-15-0	CARBON DISULFIDE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
56-23-5	CARBON TETRACHLORIDE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
108-90-7	CHLOROBENZENE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
75-00-3	CHLOROETHANE	ug/kg	0.61 UJ	0.77 UJ	1 UJ	0.72 UJ
67-66-3	CHLOROFORM	ug/kg	0.61 U	0.77 U	1 U	0.72 U
74-87-3	CHLOROMETHANE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/kg	0.61 UJ	0.77 UJ	1 UJ	0.72 UJ
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
110-82-7	CYCLOHEXANE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
124-48-1	DIBROMOCHLOROMETHANE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
75-71-8	DICHLORODIFLUOROMETHANE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
100-41-4	ETHYLBENZENE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/kg	0.58 U	0.74 UJ	0.98 UJ	0.69 U
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/kg	0.88 U	1.1 U	1.5 U	1 U
79-20-9	METHYL ACETATE	ug/kg	1.2 U	1.5 U	2 U	1.4 U
78-93-3	METHYL ETHYL KETONE	ug/kg	3.8 U	4.8 U	6.3 U	4.5 U
108-10-1	METHYL ISOBUTYL KETONE	ug/kg	3 U	3.9 U	5.1 U	3.6 U
108-87-2	METHYLCYCLOHEXANE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
75-09-2	METHYLENE CHLORIDE	ug/kg	5.2 J	3.4 J	6 J	6.3 J
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/kg	0.61 U	0.77 U	1 U	0.72 U
100-42-5	STYRENE	ug/kg	0.55 U	0.7 U	0.92 U	0.65 U
1634-04-4	TERT-BUTYL METHYL ETHER	ug/kg	0.61 U	0.77 U	1 U	0.72 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/kg	0.61 U	0.77 U	1 U	0.72 U
108-88-3	TOLUENE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/kg	0.61 U	0.77 U	1 U	0.72 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/kg	0.61 U	0.77 U	1 U	0.72 U
75-69-4	TRICHLOROFLUOROMETHANE	ug/kg	0.61 UJ	0.77 UJ	1 UJ	0.72 UJ
75-01-4	VINYL CHLORIDE	ug/kg	0.61 U	0.77 U	1 U	0.72 U

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-1 SB-1(15-17)-20131101 E4277-26	SB-2 SB-2(5-10)-20131101 E4277-21	SB-2 SB-2(10-15)-20131101 E4277-22	SB-3 SB-3(5-10)-20131101 E4277-23
CAS NO.	COMPOUND	UNITS:				
95-94-3	SEMIVOLATILES					
58-90-2	1,2,4,5-TETRACHLOROBENZENE	ug/kg	17.9 U	17.4 U	21.2 U	170 U
95-95-4	2,3,4,6-TETRACHLOROPHENOL	ug/kg	17.9 U	17.4 U	21.2 U	170 U
88-06-2	2,4,5-TRICHLOROPHENOL	ug/kg	31.9 U	31.2 U	37.8 U	310 U
120-83-2	2,4,6-TRICHLOROPHENOL	ug/kg	13.9 U	13.6 U	16.5 U	130 U
105-67-9	2,4-DICHLOROPHENOL	ug/kg	17.3 U	16.9 U	20.5 U	170 U
51-28-5	2,4-DIMETHYLPHENOL	ug/kg	25.8 U	25.2 U	30.6 U	250 U
121-14-2	2,4-DINITROTOLUENE	ug/kg	46.2 UJ	45.1 UJ	54.8 UJ	440 UJ
606-20-2	2,6-DINITROTOLUENE	ug/kg	13.6 U	13.3 U	16.2 U	130 U
91-58-7	2-CHLORONAPHTHALENE	ug/kg	18.5 U	18.1 U	22 U	180 U
95-57-8	2-CHLOROPHENOL	ug/kg	10.4 U	10.1 U	12.3 U	99.7 U
91-57-6	2-METHYLNAPHTHALENE	ug/kg	24 U	23.4 U	28.5 U	230 U
95-48-7	2-METHYLPHENOL (O-CRESOL)	ug/kg	11.4 U	11.2 U	13.6 U	110 U
88-74-4	2-NITROANILINE	ug/kg	24.7 U	24.1 U	29.3 U	240 U
88-75-5	2-NITROPHENOL	ug/kg	20.2 U	19.7 U	23.9 U	190 U
MEPH3ME	3- AND 4- METHYLPHENOL (TOTAL)	ug/kg	21.9 U	21.4 U	26 U	210 U
91-94-1	3,3'-DICHLOROBENZIDINE	ug/kg	23.6 U	23 U	28 U	230 U
99-09-2	3-NITROANILINE	ug/kg	29.2 U	28.5 U	34.6 U	280 U
534-52-1	4,6-DINITRO-2-METHYLPHENOL	ug/kg	29.2 U	28.5 U	34.6 U	280 U
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/kg	26 U	25.4 UJ	30.9 UJ	250 UJ
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/kg	8.9 U	8.7 U	10.5 U	85.3 U
106-47-8	4-CHLOROANILINE	ug/kg	20.2 U	19.7 U	23.9 U	190 U
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/kg	32 U	31.3 U	38 U	310 U
100-01-6	4-NITROANILINE	ug/kg	24.7 U	24.1 U	29.3 U	240 U
100-02-7	4-NITROPHENOL	ug/kg	59.1 U	57.8 U	70.2 U	570 U
83-32-9	ACENAPHTHENE	ug/kg	84.4 U	82.4 U	100 U	810 U
208-96-8	ACENAPHTHYLENE	ug/kg	12.8 U	12.5 U	360 J	120 U
98-86-2	ACETOPHENONE	ug/kg	11.4 U	190 J	230 J	110 U
120-12-7	ANTHRACENE	ug/kg	13.9 U	13.6 U	16.5 U	130 U
1912-24-9	ATRAZINE	ug/kg	9.3 U	380 J	1500	970 J
100-52-7	BENZALDEHYDE	ug/kg	24 U	23.4 U	28.5 U	230 U
56-55-3	BENZO(A)ANTHRACENE	ug/kg	R	R	R	R
50-32-8	BENZO(A)PYRENE	ug/kg	21.7 U	1400	3600	6300
205-99-2	BENZO(B)FLUORANTHENE	ug/kg	9.8 U	1600	3200	7200
191-24-2	BENZO(G,H,I)PERYLENE	ug/kg	14.9 U	1800	3800	8700
207-08-9	BENZO(K)FLUORANTHENE	ug/kg	18.4 U	1200	1800	5100
85-68-7	BENZYL BUTYL PHTHALATE	ug/kg	21.4 U	840	1100	3300 J
92-52-4	BIPHENYL (DIPHENYL)	ug/kg	21.8 U	310 J	25.9 U	210 U
111-91-1	BIS(2-CHLOROETHOXY) METHANE	ug/kg	17.2 U	16.8 U	20.4 U	170 U
111-44-4	BIS(2-CHLOROETHYL) ETHER	ug/kg	26.2 U	25.6 U	31.1 U	250 U
108-60-1	BIS(2-CHLOROISOPROPYL) ETHER	ug/kg	21.8 U	21.3 U	25.9 U	210 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/kg	18.8 U	18.4 U	22.3 U	180 U
105-60-2	BIS(2-ETHYLHEXYL) PHTHALATE	ug/kg	16.1 U	15.7 UJ	19.1 UJ	150 UJ
86-74-8	CAPROLACTAM	ug/kg	21.1 U	20.6 U	25.1 U	200 U
218-01-9	CARBAZOLE	ug/kg	9.9 U	92.8 J	290 J	95.8 U
53-70-3	CHRYSENE	ug/kg	9.9 U	1400	2400	5900
132-64-9	DIBENZ(A,H)ANTHRACENE	ug/kg	20.6 U	690	1200 J	170 U
84-66-2	DIBENZOFURAN	ug/kg	13.1 U	17.7 U	350 J	170 U
131-11-3	DIETHYL PHTHALATE	ug/kg	7.1 U	6.9 U	8.4 U	68.2 U
84-74-2	DI-N-BUTYL PHTHALATE	ug/kg	440 J	450	890	120 U
117-84-0	DI-N-OCTYL PHTHALATE	ug/kg	35.7 U	34.9 U	42.4 U	340 U
206-44-0	FLUORANTHENE	ug/kg	5.2 U	5.1 UJ	6.1 UJ	49.9 UJ
86-73-7	FLUORENE	ug/kg	9.1 U	2000	7000	9900
118-74-1	HEXACHLOROBENZENE	ug/kg	17.2 U	130 J	740	170 U
87-68-3	HEXACHLOROBUTADIENE	ug/kg	18.5 U	18.1 U	22 U	180 U
77-47-4	HEXACHLOROCYCLOPENTADIENE	ug/kg	16.5 U	16.1 U	19.6 U	160 U
67-72-1	HEXACHLOROETHANE	ug/kg	11 U	10.8 UJ	13.1 UJ	110 UJ
193-39-5	INDENO(1,2,3-C,D)PYRENE	ug/kg	20.3 U	19.8 U	24.1 U	200 U
78-59-1	ISOPHORONE	ug/kg	15.1 U	970	1700	4500
		ug/kg	15 U	14.6 U	17.8 U	140 U

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-1 SB-1(15-17)-20131101 E4277-26 15 - 17 ft CTECH E4277 SOIL 11/1/2013 13:55 2/4/2014	SB-2 SB-2(5-10)-20131101 E4277-21 5 - 10 ft CTECH E4277 SOIL 11/1/2013 9:45 2/4/2014	SB-2 SB-2(10-15)-20131101 E4277-22 10 - 15 ft CTECH E4277 SOIL 11/1/2013 9:55 2/4/2014	SB-3 SB-3(5-10)-20131101 E4277-23 5 - 10 ft CTECH E4277 SOIL 11/1/2013 11:15 2/4/2014
CAS NO.	COMPOUND	UNITS:				
	SEMIVOLATILES					
91-20-3	NAPHTHALENE	ug/kg	15.7 U	290 J	200 J	150 U
98-95-3	NITROBENZENE	ug/kg	17.2 UJ	16.8 U	20.4 U	170 U
621-64-7	N-NITROSODI-N-PROPYLAMINE	ug/kg	22.9 U	22.4 U	27.2 U	220 U
86-30-6	N-NITROSODIPHENYLAMINE	ug/kg	10.9 U	10.7 U	12.9 U	100 U
87-86-5	PENTACHLOROPHENOL	ug/kg	31.1 UJ	30.4 U	36.9 U	300 U
85-01-8	PHENANTHRENE	ug/kg	12.3 U	1200	5000	4300
108-95-2	PHENOL	ug/kg	10.5 U	10.3 UJ	12.5 UJ	100 UJ
129-00-0	PYRENE	ug/kg	10.9 U	1800	5500	7000
	INORGANICS					
7429-90-5	ALUMINUM	mg/kg	5900	3100	11100	3900
7440-36-0	ANTIMONY	mg/kg	0.639 UJ	2.67 J	0.744 UJ	9.52 J
7440-38-2	ARSENIC	mg/kg	2.44	10.67	10.43	27.6
7440-39-3	BARIUM	mg/kg	17.6	416	168	354
7440-41-7	BERYLLIUM	mg/kg	0.16 J	0.087 J	0.426	0.065 U
7440-43-9	CADMIUM	mg/kg	0.372	1.33	0.08 U	2.73
7440-70-2	CALCIUM	mg/kg	1300	13500	5900	38300
7440-47-3	CHROMIUM, TOTAL	mg/kg	14.9	30.7	25.9	24.9
7440-48-4	COBALT	mg/kg	5.15	6.78	12.3	9.71
7440-50-8	COPPER	mg/kg	10.34	595	35.3	361
7439-89-6	IRON	mg/kg	10500	31000	32900	68200
7439-92-1	LEAD	mg/kg	4.24 J	727 J	168 J	1500 J
7439-95-4	MAGNESIUM	mg/kg	2400	1800	6100	3800
7439-96-5	MANGANESE	mg/kg	129	229	695	331
7439-97-6	MERCURY	mg/kg	0.007 U	0.67	0.765	3.85
7440-02-0	NICKEL	mg/kg	10.84 J	17.8 J	26.7 J	35.7 J
7440-09-7	POTASSIUM	mg/kg	1300 J	906 J	2700 J	476 J
7782-49-2	SELENIUM	mg/kg	0.556 J	2.96	2.22	3.95
7440-22-4	SILVER	mg/kg	0.252 J	2.21	2.29	4.94
7440-23-5	SODIUM	mg/kg	762	649	1100	658
7440-28-0	THALLIUM	mg/kg	0.308 U	0.322 U	0.359 U	0.292 U
7440-62-2	VANADIUM	mg/kg	23.2	17.7	43.9	23.4
7440-66-6	ZINC	mg/kg	26.2	830	187	1100
57-12-5	CYANIDE	mg/kg	0.045 U	1.86	0.158 J	2.94

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-3 SB-3(12-14)-20131101 E4277-24 12 - 14 ft CTECH E4277 SOIL 11/1/2013 11:25 2/4/2014	SB-4 SB-4(15-17)-20131031 E4277-19 15 - 17 ft CTECH E4277 SOIL 10/31/2013 14:15 2/4/2014	SB-4 SB-4(19-20)-20131031 E4277-20 19 - 20 ft CTECH E4277 SOIL 10/31/2013 14:15 2/4/2014	SB-5 SB-5(5-10)-20131031 E4277-17 5 - 10 ft CTECH E4277 SOIL 10/31/2013 12:40 2/4/2014
CAS NO.	COMPOUND	UNITS:				
	VOLATILES					
71-55-6	1,1,1-TRICHLOROETHANE	ug/kg	0.96 U	400 U	1.3 U	0.71 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/kg	0.88 U	360 U	1.2 U	0.65 U
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/kg	0.96 U	400 U	1.3 U	0.71 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/kg	1.7 U	710 U	2.3 U	1.3 U
75-34-3	1,1-DICHLOROETHANE	ug/kg	0.96 U	400 U	1.3 U	0.71 U
75-35-4	1,1-DICHLOROETHENE	ug/kg	0.96 U	400 U	1.3 U	0.71 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/kg	0.96 U	9900	1.3 U	0.71 U
120-82-1	1,2,4-TRICHLOROBENZENE	ug/kg	0.96 U	63200 J	1.3 U	0.71 U
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/kg	1.7 U	690 U	2.2 U	1.2 U
106-93-4	1,2-DIBromoETHANE	ug/kg	0.96 U	400 U	1.3 U	0.71 U
95-50-1	1,2-DICHLOROBENZENE	ug/kg	0.96 U	400 U	1.3 UJ	0.71 U
107-06-2	1,2-DICHLOROETHANE	ug/kg	0.96 U	400 U	1.3 U	0.71 U
78-87-5	1,2-DICHLOROPROPANE	ug/kg	0.5 U	210 U	0.65 U	0.37 U
541-73-1	1,3-DICHLOROBENZENE	ug/kg	0.71 U	290 U	0.93 U	0.52 U
106-46-7	1,4-DICHLOROBENZENE	ug/kg	0.79 U	320 U	2.5 J	0.58 U
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/kg	190 U	79200 UJ	250 U	140 U
591-78-6	2-HEXANONE	ug/kg	4.8 U	2000 UJ	6.3 U	3.5 U
67-64-1	ACETONE	ug/kg	4.8 U	2000 UJ	40.8 J	3.5 U
71-43-2	BENZENE	ug/kg	0.73 U	85300	38.4	0.54 U
74-97-5	BROMOCHLOROMETHANE	ug/kg	0.96 U	400 U	1.3 U	0.71 U
75-27-4	BROMODICHLOROMETHANE	ug/kg	0.96 U	400 U	1.3 U	0.71 U
75-25-2	BROMOFORM	ug/kg	1.4 U	590 UJ	1.9 U	1 U
74-83-9	BROMOMETHANE	ug/kg	1.9 U	790 U	2.5 U	1.4 U
75-15-0	CARBON DISULFIDE	ug/kg	0.96 U	400 U	8 J	0.71 U
56-23-5	CARBON TETRACHLORIDE	ug/kg	0.96 U	400 UJ	1.3 U	0.71 U
108-90-7	CHLOROBENZENE	ug/kg	0.96 U	400 U	1.3 U	0.71 U
75-00-3	CHLOROETHANE	ug/kg	0.96 UJ	400 U	1.3 U	0.71 UJ
67-66-3	CHLOROFORM	ug/kg	0.96 U	400 U	1.3 U	0.71 U
74-87-3	CHLOROMETHANE	ug/kg	0.96 U	400 U	1.3 U	0.71 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/kg	0.96 UJ	400 U	1.3 U	0.71 UJ
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/kg	0.96 U	400 U	1.3 U	0.71 U
110-82-7	CYCLOHEXANE	ug/kg	0.96 U	400 U	1.3 U	0.71 U
124-48-1	DIBROMOCHLOROMETHANE	ug/kg	0.96 U	400 U	1.3 U	0.71 U
75-71-8	DICHLORODIFLUOROMETHANE	ug/kg	0.96 U	400 U	1.3 U	0.71 U
100-41-4	ETHYLBENZENE	ug/kg	0.96 U	518300 J	37.9	0.71 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/kg	0.92 U	29000 J	6.5 J	0.68 U
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/kg	1.4 U	438100	14.6 J	1 U
79-20-9	METHYL ACETATE	ug/kg	1.9 U	790 U	2.5 U	1.4 U
78-93-3	METHYL ETHYL KETONE	ug/kg	6 U	2500 UJ	7.8 U	4.4 U
108-10-1	METHYL ISOBUTYL KETONE	ug/kg	4.8 U	2000 UJ	6.3 U	3.5 U
108-87-2	METHYLCYCLOHEXANE	ug/kg	0.96 U	660 J	1.3 U	0.71 U
75-09-2	METHYLENE CHLORIDE	ug/kg	6.9 J	1700 J	17.5	0.71 U
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/kg	0.96 U	112900	16	0.71 U
100-42-5	STYRENE	ug/kg	0.87 U	360 U	1.1 U	0.64 U
1634-04-4	TERT-BUTYL METHYL ETHER	ug/kg	0.96 U	400 U	10.6 J	0.71 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/kg	0.96 U	400 U	1.3 U	0.71 U
108-88-3	TOLUENE	ug/kg	0.96 U	18400	1.3 U	0.71 U
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/kg	0.96 U	400 U	1.3 U	0.71 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/kg	0.96 U	400 U	1.3 U	0.71 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/kg	0.96 U	400 U	1.3 U	0.71 U
75-69-4	TRICHLOROFUOROMETHANE	ug/kg	0.96 UJ	400 U	1.3 U	0.71 UJ
75-01-4	VINYL CHLORIDE	ug/kg	0.96 U	400 U	1.3 U	0.71 U

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-3 SB-3(12-14)-20131101 E4277-24 12 - 14 ft CTECH E4277 SOIL 11/1/2013 11:25 2/4/2014	SB-4 SB-4(15-17)-20131031 E4277-19 15 - 17 ft CTECH E4277 SOIL 10/31/2013 14:15 2/4/2014	SB-4 SB-4(19-20)-20131031 E4277-20 19 - 20 ft CTECH E4277 SOIL 10/31/2013 14:15 2/4/2014	SB-5 SB-5(5-10)-20131031 E4277-17 5 - 10 ft CTECH E4277 SOIL 10/31/2013 12:40 2/4/2014
CAS NO.	COMPOUND	UNITS:				
	SEMIVOLATILES					
95-94-3	1,2,4,5-TETRACHLOROBENZENE	ug/kg	23.2 U	200 U	23 U	16.9 U
58-90-2	2,3,4,6-TETRACHLOROPHENOL	ug/kg	23.2 U	200 U	23 U	16.9 U
95-95-4	2,4,5-TRICHLOROPHENOL	ug/kg	41.5 U	350 U	41 U	30.2 U
88-06-2	2,4,6-TRICHLOROPHENOL	ug/kg	18.1 U	150 U	17.9 U	13.2 U
120-83-2	2,4-DICHLOROPHENOL	ug/kg	22.5 U	190 U	22.3 U	16.4 U
105-67-9	2,4-DIMETHYLPHENOL	ug/kg	33.5 U	280 U	33.1 U	24.4 U
51-28-5	2,4-DINITROPHENOL	ug/kg	60.2 UJ	510 UJ	59.4 UJ	43.8 UJ
121-14-2	2,4-DINITROTOLUENE	ug/kg	17.7 U	150 U	17.5 U	12.9 U
606-20-2	2,6-DINITROTOLUENE	ug/kg	24.1 U	200 U	23.8 U	17.6 U
91-58-7	2-CHLORONAPHTHALENE	ug/kg	13.5 U	110 U	13.3 U	9.8 U
95-57-8	2-CHLOROPHENOL	ug/kg	31.2 U	260 U	30.8 U	22.7 U
91-57-6	2-METHYLNAPHTHALENE	ug/kg	14.9 U	766400	160 J	300 J
95-48-7	2-METHYLPHENOL (O-CRESOL)	ug/kg	32.1 U	270 U	31.7 U	23.4 U
88-74-4	2-NITROANILINE	ug/kg	26.3 U	220 U	25.9 U	19.1 U
88-75-5	2-NITROPHENOL	ug/kg	28.6 U	240 U	28.2 U	20.8 U
MEPH3ME	3- AND 4- METHYLPHENOL (TOTAL)	ug/kg	260 J	260 U	30.3 U	22.4 U
91-94-1	3,3'-DICHLOROBENZIDINE	ug/kg	38 U	320 U	37.5 U	27.6 U
99-09-2	3-NITROANILINE	ug/kg	38 U	320 U	37.5 U	27.6 U
534-52-1	4,6-DINITRO-2-METHYLPHENOL	ug/kg	33.9 UJ	280 UJ	33.5 UJ	24.7 UJ
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/kg	11.5 U	96.9 U	11.4 U	8.4 U
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/kg	26.3 U	220 U	25.9 U	19.1 U
106-47-8	4-CHLOROANILINE	ug/kg	41.7 U	350 U	41.2 U	30.4 U
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/kg	32.1 U	270 U	31.7 U	23.4 U
100-01-6	4-NITROANILINE	ug/kg	77 U	650 U	76 U	56.1 U
100-02-7	4-NITROPHENOL	ug/kg	110 U	920 U	110 U	80 U
83-32-9	ACENAPHTHENE	ug/kg	16.7 U	233200	16.5 U	12.1 U
208-96-8	ACENAPHTHYLENE	ug/kg	14.9 U	33300	14.7 U	2400
98-86-2	ACETOPHENONE	ug/kg	18.1 U	150 U	17.9 U	13.2 U
120-12-7	ANTHRACENE	ug/kg	12.1 U	133600	130 J	5400
1912-24-9	ATRAZINE	ug/kg	31.2 U	260 U	30.8 U	22.7 U
100-52-7	BENZALDEHYDE	ug/kg	R	R	R	R
56-55-3	BENZO(A)ANTHRACENE	ug/kg	310 J	80200	230 J	8200
50-32-8	BENZO(A)PYRENE	ug/kg	310 J	72200	210 J	7300
205-99-2	BENZO(B)FLUORANTHENE	ug/kg	370 J	59100	240 J	7700
191-24-2	BENZO(G,H,I)PERYLENE	ug/kg	200 J	29200	23.7 U	3100
207-08-9	BENZO(K)FLUORANTHENE	ug/kg	150 J	14100	27.5 U	3800
85-68-7	BENZYL BUTYL PHTHALATE	ug/kg	28.4 U	240 U	28 U	20.7 U
92-52-4	BIPHENYL (DIPHENYL)	ug/kg	22.4 U	74500	22.1 U	16.3 U
111-91-1	BIS(2-CHLOROETHOXY) METHANE	ug/kg	34.1 U	290 U	33.6 U	24.8 U
111-44-4	BIS(2-CHLOROETHYL) ETHER	ug/kg	28.4 U	240 U	28 U	20.7 U
108-60-1	BIS(2-CHLOROISOPROPYL) ETHER	ug/kg	24.5 U	210 U	24.2 U	17.8 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/kg	20.9 UJ	180 U	20.7 UJ	15.2 UJ
105-60-2	CAPROLACTAM	ug/kg	27.5 U	230 U	27.2 U	20 U
86-74-8	CARBAZOLE	ug/kg	13 U	3600 J	12.8 U	1400
218-01-9	CHRYSENE	ug/kg	240 J	73400	180 J	7300
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/kg	17 U	6900	16.8 U	870
132-64-9	DIBENZOFURAN	ug/kg	23.1 U	23200	22.8 U	440
84-66-2	DIETHYL PHTHALATE	ug/kg	9.2 U	77.5 U	9.1 U	6.7 U
131-11-3	DIMETHYL PHTHALATE	ug/kg	1000	130 U	600	660
84-74-2	DI-N-BUTYL PHTHALATE	ug/kg	46.5 U	390 U	45.9 U	33.8 U
117-84-0	DI-N-OCTYLPHTHALATE	ug/kg	6.7 UJ	56.7 U	6.7 UJ	4.9 UJ
206-44-0	FLUORANTHENE	ug/kg	750	145400	410 J	14800
86-73-7	FLUORENE	ug/kg	22.4 U	165000	22.1 U	1100
118-74-1	HEXACHLOROBENZENE	ug/kg	24.1 U	200 U	23.8 U	17.6 U
87-68-3	HEXACHLOROBUTADIENE	ug/kg	21.5 U	180 U	21.2 U	15.6 U
77-47-4	HEXACHLOROCYCLOPENTADIENE	ug/kg	14.4 UJ	120 U	14.2 UJ	10.5 UJ
67-72-1	HEXACHLOROETHANE	ug/kg	26.4 U	220 U	26.1 U	19.2 U
193-39-5	INDENO(1,2,3-C,D)PYRENE	ug/kg	170 J	21600	19.4 U	3200
78-59-1	ISOPHORONE	ug/kg	19.5 U	160 U	19.3 U	14.2 U

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-3 SB-3(12-14)-20131101 E4277-24 12 - 14 ft CTECH E4277 SOIL 11/1/2013 11:25 2/4/2014	SB-4 SB-4(15-17)-20131031 E4277-19 15 - 17 ft CTECH E4277 SOIL 10/31/2013 14:15 2/4/2014	SB-4 SB-4(19-20)-20131031 E4277-20 19 - 20 ft CTECH E4277 SOIL 10/31/2013 14:15 2/4/2014	SB-5 SB-5(5-10)-20131031 E4277-17 5 - 10 ft CTECH E4277 SOIL 10/31/2013 12:40 2/4/2014
CAS NO.	COMPOUND	UNITS:				
	SEMIVOLATILES					
91-20-3	NAPHTHALENE	ug/kg	170 J	1899200	480 J	1200
98-95-3	NITROBENZENE	ug/kg	22.4 U	190 U	22.1 U	16.3 U
621-64-7	N-NITROSODI-N-PROPYLAMINE	ug/kg	29.8 U	250 U	29.4 U	21.7 U
86-30-6	N-NITROSODIPHENYLAMINE	ug/kg	14.2 U	120 U	14 U	10.3 U
87-86-5	PENTACHLOROPHENOL	ug/kg	40.5 U	340 U	40 U	29.5 U
85-01-8	PHENANTHRENE	ug/kg	250 J	547300	380 J	12300
108-95-2	PHENOL	ug/kg	13.7 UJ	110 U	13.5 UJ	9.9 UJ
129-00-0	PYRENE	ug/kg	650	253600	370 J	13000
	INORGANICS					
7429-90-5	ALUMINUM	mg/kg	2300	4000	15500	2700
7440-36-0	ANTIMONY	mg/kg	15 J	1.74 J	0.813 UJ	0.641 UJ
7440-38-2	ARSENIC	mg/kg	11.8	12	10.52	5.42
7440-39-3	BARIUM	mg/kg	369	78.3	54.5	170
7440-41-7	BERYLLIUM	mg/kg	0.091 U	0.291 J	0.643	0.204 J
7440-43-9	CADMIUM	mg/kg	1.31	0.3 J	0.087 U	0.069 U
7440-70-2	CALCIUM	mg/kg	7700	2000	3700	3200
7440-47-3	CHROMIUM, TOTAL	mg/kg	10.95	16.7	35	24.8
7440-48-4	COBALT	mg/kg	16.8	9.46	17.2	5.01
7440-50-8	COPPER	mg/kg	64.9	92.1	28.2	129
7439-89-6	IRON	mg/kg	102200	46400	40100	13000
7439-92-1	LEAD	mg/kg	595 J	292 J	63.9 J	737 J
7439-95-4	MAGNESIUM	mg/kg	721	374	8700	1200
7439-96-5	MANGANESE	mg/kg	1300	113	808	49.9
7439-97-6	MERCURY	mg/kg	0.879	3.11	0.461	0.199
7440-02-0	NICKEL	mg/kg	12.4 J	21.6 J	37.2 J	11.2 J
7440-09-7	POTASSIUM	mg/kg	597 J	597 J	4100 J	584 J
7782-49-2	SELENIUM	mg/kg	1.24 J	3.67	2.49	1.95
7440-22-4	SILVER	mg/kg	5.72	2.97	2.81	0.862
7440-23-5	SODIUM	mg/kg	1500	430	3400	299
7440-28-0	THALLIUM	mg/kg	1.07 J	0.35 U	0.392 U	0.309 U
7440-62-2	VANADIUM	mg/kg	36.3	19.5	58.1	24.5
7440-66-6	ZINC	mg/kg	781	495	219	72
57-12-5	CYANIDE	mg/kg	9.8	34.9	1.84	0.79

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-5 SB-5(10-15)-20131031 E4277-18	SB-6 SB-6(8-10)-20131031 E4277-12	SB-6 SB-6(10-15)-20131031 E4277-15	SB-7 SB-7(8-10)-20131031 E4277-10
CAS NO.	COMPOUND	UNITS:				
	VOLATILES					
71-55-6	1,1,1-TRICHLOROETHANE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/kg	0.61 UJ	0.75 UJ	0.81 U	0.55 UJ
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/kg	1.2 U	1.5 U	1.6 U	1.1 U
75-34-3	1,1-DICHLOROETHANE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
75-35-4	1,1-DICHLOROETHENE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/kg	0.66 UJ	0.81 UJ	0.88 U	0.6 UJ
120-82-1	1,2,4-TRICHLOROBENZENE	ug/kg	0.66 UJ	0.81 UJ	0.88 U	0.6 UJ
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/kg	1.2 UJ	1.4 UJ	1.5 U	1 UJ
106-93-4	1,2-DIBromoETHANE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
95-50-1	1,2-DICHLOROBENZENE	ug/kg	0.66 UJ	0.81 UJ	0.88 U	0.6 UJ
107-06-2	1,2-DICHLOROETHANE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
78-87-5	1,2-DICHLOROPROPANE	ug/kg	0.35 U	0.42 U	0.46 U	0.31 U
541-73-1	1,3-DICHLOROBENZENE	ug/kg	0.49 UJ	0.6 UJ	0.65 U	0.44 UJ
106-46-7	1,4-DICHLOROBENZENE	ug/kg	0.55 UJ	0.66 UJ	0.72 U	0.49 UJ
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/kg	130 U	160 U	180 U	120 U
591-78-6	2-HEXANONE	ug/kg	3.3 U	4.1 U	4.4 U	3 U
67-64-1	ACETONE	ug/kg	3.3 U	4.1 U	63.7	33.7
71-43-2	BENZENE	ug/kg	0.51 U	0.62 U	0.67 U	0.46 U
74-97-5	BROMOCHLOROMETHANE	ug/kg	0.66 U	0.81 U	0.88 UJ	0.6 U
75-27-4	BROMODICHLOROMETHANE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
75-25-2	BROMOFORM	ug/kg	0.98 U	1.2 U	1.3 U	0.89 U
74-83-9	BROMOMETHANE	ug/kg	1.3 U	1.6 U	1.8 U	1.2 U
75-15-0	CARBON DISULFIDE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
56-23-5	CARBON TETRACHLORIDE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
108-90-7	CHLOROBENZENE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
75-00-3	CHLOROETHANE	ug/kg	0.66 UJ	0.81 U	0.88 U	0.6 U
67-66-3	CHLOROFORM	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
74-87-3	CHLOROMETHANE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/kg	0.66 UJ	0.81 U	0.88 U	0.6 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
110-82-7	CYCLOHEXANE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
124-48-1	DIBROMOCHLOROMETHANE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
75-71-8	DICHLORODIFLUOROMETHANE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
100-41-4	ETHYLBENZENE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/kg	0.64 UJ	0.78 UJ	0.84 U	0.58 UJ
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/kg	0.96 U	1.2 U	1.3 U	0.86 U
79-20-9	METHYL ACETATE	ug/kg	1.3 U	1.6 U	1.8 U	1.2 U
78-93-3	METHYL ETHYL KETONE	ug/kg	4.1 U	5 U	5.5 U	3.7 U
108-10-1	METHYL ISOBUTYL KETONE	ug/kg	3.3 U	4.1 U	4.4 U	3 U
108-87-2	METHYLCYCLOHEXANE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
75-09-2	METHYLENE CHLORIDE	ug/kg	4 J	7.8 J	8.8 U	4.4 J
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
100-42-5	STYRENE	ug/kg	0.6 U	0.73 U	0.79 U	0.54 U
1634-04-4	TERT-BUTYL METHYL ETHER	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
108-88-3	TOLUENE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U
75-69-4	TRICHLOROFLUOROMETHANE	ug/kg	0.66 UJ	0.81 U	0.88 U	0.6 U
75-01-4	VINYL CHLORIDE	ug/kg	0.66 U	0.81 U	0.88 U	0.6 U

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-5 SB-5(10-15)-20131031 E4277-18	SB-6 SB-6(8-10)-20131031 E4277-12	SB-6 SB-6(10-15)-20131031 E4277-15	SB-7 SB-7(8-10)-20131031 E4277-10
CAS NO.	COMPOUND	UNITS:				
95-94-3	SEMIVOLATILES	ug/kg	15.7 U	16 U	19.8 U	14.8 U
58-90-2	1,2,4,5-TETRACHLOROBENZENE	ug/kg	15.7 U	16 U	19.8 U	14.8 U
95-95-4	2,3,4,6-TETRACHLOROPHENOL	ug/kg	28 U	28.5 U	35.3 U	26.4 U
88-06-2	2,4,5-TRICHLOROPHENOL	ug/kg	12.2 U	12.4 U	15.4 U	11.5 U
120-83-2	2,4-DICHLOROPHENOL	ug/kg	15.2 U	15.5 U	19.2 U	14.3 U
105-67-9	2,4-DIMETHYLPHENOL	ug/kg	22.6 U	23 U	28.5 U	21.3 U
51-28-5	2,4-DINITROPHENOL	ug/kg	40.5 UJ	41.3 UJ	51.1 UJ	38.2 UJ
121-14-2	2,4-DINITROTOLUENE	ug/kg	11.9 U	12.2 U	15.1 U	11.3 U
606-20-2	2,6-DINITROTOLUENE	ug/kg	16.2 U	16.6 U	20.5 U	15.3 U
91-58-7	2-CHLORONAPHTHALENE	ug/kg	9.1 U	9.3 U	11.5 U	8.6 U
95-57-8	2-CHLOROPHENOL	ug/kg	21 U	21.4 U	26.6 U	19.9 U
91-57-6	2-METHYLNAPHTHALENE	ug/kg	10 U	170 J	130 J	86.9 J
95-48-7	2-METHYLPHENOL (O-CRESOL)	ug/kg	21.6 U	22.1 U	27.3 U	20.4 U
88-74-4	2-NITROANILINE	ug/kg	17.7 U	18 U	22.3 U	16.7 U
88-75-5	2-NITROPHENOL	ug/kg	19.2 U	19.6 U	24.3 U	18.2 U
MEPH3ME	3- AND 4- METHYLPHENOL (TOTAL)	ug/kg	20.7 U	21.1 U	26.1 U	19.5 U
91-94-1	3,3'-DICHLOROBENZIDINE	ug/kg	25.6 U	26.1 U	32.3 U	24.1 U
99-09-2	3-NITROANILINE	ug/kg	25.6 U	26.1 U	32.3 U	24.1 U
534-52-1	4,6-DINITRO-2-METHYLPHENOL	ug/kg	22.8 UJ	23.3 UJ	28.8 UJ	21.5 UJ
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/kg	7.8 U	7.9 U	9.8 U	7.3 U
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/kg	17.7 U	18 U	22.3 U	16.7 U
106-47-8	4-CHLOROANILINE	ug/kg	28.1 U	28.6 U	35.5 U	26.5 U
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/kg	21.6 U	22.1 U	27.3 U	20.4 U
100-01-6	4-NITROANILINE	ug/kg	51.8 U	52.9 U	65.5 U	49 U
100-02-7	4-NITROPHENOL	ug/kg	73.9 U	75.4 U	93.4 U	69.8 U
83-32-9	ACENAPHTHENE	ug/kg	11.2 U	11.5 U	14.2 U	10.6 U
208-96-8	ACENAPHTHYLENE	ug/kg	10 U	340 J	500	230 J
98-86-2	ACETOPHENONE	ug/kg	12.2 U	12.4 U	15.4 U	11.5 U
120-12-7	ANTHRACENE	ug/kg	8.1 U	320 J	670	330 J
1912-24-9	ATRAZINE	ug/kg	21 U	21.4 U	26.6 U	19.9 U
100-52-7	BENZALDEHYDE	ug/kg	R	R	R	R
56-55-3	BENZO(A)ANTHRACENE	ug/kg	130 J	750	1700	1200
50-32-8	BENZO(A)PYRENE	ug/kg	120 J	730	1500	1200
205-99-2	BENZO(B)FLUORANTHENE	ug/kg	150 J	790	1800	1500
191-24-2	BENZO(G,H,I)PERYLENE	ug/kg	16.1 U	550	900	890 J
207-08-9	BENZO(K)FLUORANTHENE	ug/kg	18.8 U	380 J	500	360 J
85-68-7	BENZYL BUTYL PHTHALATE	ug/kg	19.1 U	19.5 U	24.1 U	18.1 U
92-52-4	BIPHENYL (DIPHENYL)	ug/kg	15.1 U	15.4 U	19 U	14.2 U
111-91-1	BIS(2-CHLOROETHOXY) METHANE	ug/kg	22.9 U	23.4 U	29 U	21.7 U
111-44-4	BIS(2-CHLOROETHYL) ETHER	ug/kg	19.1 U	19.5 U	24.1 U	18.1 U
108-60-1	BIS(2-CHLOROISOPROPYL) ETHER	ug/kg	16.5 U	16.8 U	20.8 U	15.6 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/kg	14.1 UJ	14.4 U	17.8 UJ	13.3 U
105-60-2	CAPROLACTAM	ug/kg	18.5 U	18.9 U	23.4 U	17.5 U
86-74-8	CARBAZOLE	ug/kg	8.7 U	8.9 U	120 J	80.1 J
218-01-9	CHRYSENE	ug/kg	110 J	710	1300	1200
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/kg	11.5 U	120 J	240 J	210 J
132-64-9	DIBENZOFURAN	ug/kg	15.5 U	15.8 U	150 J	14.7 U
84-66-2	DIETHYL PHTHALATE	ug/kg	6.2 U	6.3 U	150 J	5.9 U
131-11-3	DIMETHYL PHTHALATE	ug/kg	310 J	450	680	240 J
84-74-2	DI-N-BUTYL PHTHALATE	ug/kg	31.3 U	31.9 U	39.5 U	29.6 U
117-84-0	DI-N-OCTYLPHthalate	ug/kg	4.5 UJ	4.6 U	5.7 UJ	4.3 U
206-44-0	FLUORANTHENE	ug/kg	210 J	1400	2800	1700
86-73-7	FLUORENE	ug/kg	15.1 U	250 J	360 J	150 J
118-74-1	HEXACHLOROBENZENE	ug/kg	16.2 U	16.6 U	20.5 U	15.3 U
87-68-3	HEXACHLOROBUTADIENE	ug/kg	14.5 U	14.7 U	18.3 U	13.7 U
77-47-4	HEXACHLOROCYCLOPENTADIENE	ug/kg	9.7 UJ	9.9 U	12.2 UJ	9.1 U
67-72-1	HEXACHLOROETHANE	ug/kg	17.8 U	18.2 U	22.5 U	16.8 U
193-39-5	INDENO(1,2,3-C,D)PYRENE	ug/kg	13.3 U	450	790	830
78-59-1	ISOPHORONE	ug/kg	13.1 U	13.4 U	16.6 U	12.4 U

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-5 SB-5(10-15)-20131031 E4277-18 10 - 15 ft CTECH E4277 SOIL	SB-6 SB-6(8-10)-20131031 E4277-12 8 - 10 ft CTECH E4277 SOIL	SB-6 SB-6(10-15)-20131031 E4277-15 10 - 15 ft CTECH E4277 SOIL	SB-7 SB-7(8-10)-20131031 E4277-10 8 - 10 ft CTECH E4277 SOIL
CAS NO.	COMPOUND	UNITS:	10/31/2013 12:50 2/4/2014	10/31/2013 10:20 2/4/2014	10/31/2013 10:30 2/4/2014	10/31/2013 8:35 2/4/2014
	SEMIVOLATILES					
91-20-3	NAPHTHALENE	ug/kg	190 J	180 J	300 J	130 J
98-95-3	NITROBENZENE	ug/kg	15.1 U	15.4 U	19 U	14.2 U
621-64-7	N-NITROSODI-N-PROPYLAMINE	ug/kg	20.1 U	20.5 U	25.3 U	19 U
86-30-6	N-NITROSODIPHENYLAMINE	ug/kg	9.6 U	9.7 U	12.1 U	9 U
87-86-5	PENTACHLOROPHENOL	ug/kg	27.2 U	27.8 U	34.4 U	25.7 U
85-01-8	PHENANTHRENE	ug/kg	170 J	1200	2200	1000
108-95-2	PHENOL	ug/kg	9.2 UJ	9.4 U	11.6 UJ	8.7 U
129-00-0	PYRENE	ug/kg	200 J	1500	2600	1800
	INORGANICS					
7429-90-5	ALUMINUM	mg/kg	6300	7300	6000	6800
7440-36-0	ANTIMONY	mg/kg	0.589 UJ	1.36 J	0.971 J	0.76 J
7440-38-2	ARSENIC	mg/kg	7.29	15.9	14.3	4.36
7440-39-3	BARIUM	mg/kg	377	203	420	195
7440-41-7	BERYLLIUM	mg/kg	0.474	0.34	0.29 J	0.317
7440-43-9	CADMIUM	mg/kg	0.619	0.404	4.22	2.36
7440-70-2	CALCIUM	mg/kg	5000	15000	5100	6400
7440-47-3	CHROMIUM, TOTAL	mg/kg	20	62.1	79.2	20.6
7440-48-4	COBALT	mg/kg	7.22	7.31	8.37	9.55
7440-50-8	COPPER	mg/kg	120	79.6	310	120
7439-89-6	IRON	mg/kg	18200	34300	58900	22400
7439-92-1	LEAD	mg/kg	494 J	227 J	870 J	301 J
7439-95-4	MAGNESIUM	mg/kg	2300	3200	1800	2800
7439-96-5	MANGANESE	mg/kg	128	165	134	237
7439-97-6	MERCURY	mg/kg	0.319	0.39	1.46	0.361
7440-02-0	NICKEL	mg/kg	18.5 J	55.1 J	59.4 J	29.4 J
7440-09-7	POTASSIUM	mg/kg	1200 J	1100 J	1600 J	2300 J
7782-49-2	SELENIUM	mg/kg	1.94	2.5	2.25	1.1
7440-22-4	SILVER	mg/kg	1.31	2.41	2.36	0.661
7440-23-5	SODIUM	mg/kg	1600	1300	371	242
7440-28-0	THALLIUM	mg/kg	0.284 U	0.278 U	0.338 U	0.26 U
7440-62-2	VANADIUM	mg/kg	29	41.2	43.4	41.8
7440-66-6	ZINC	mg/kg	259	336	528	315
57-12-5	CYANIDE	mg/kg	4.13	7.32	5.4	0.037 UJ

		Dup of SB-7(8-10)-20131031				
		Location ID:	SB-7	SB-7	SB-8	SB-8
		Sample ID:	SB-17(8-10)-20131031	SB-7(16-18)-20131031	SB-8(15-16)-20131030	SB-8(17-19)-20131030
		Lab Sample Id:	E4277-16	E4277-11	E4277-07	E4277-08
		Depth:	8 - 10 ft	16 - 18 ft	15 - 16 ft	17 - 19 ft
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	E4277	E4277	E4277	E4277
		Matrix:	SOIL	SOIL	SOIL	SOIL
		Sampled:	10/31/2013 8:45	10/31/2013 10:20	10/30/2013 14:55	10/30/2013 14:55
		Validated:	2/4/2014	2/4/2014	2/4/2014	2/4/2014
CAS NO.	COMPOUND	UNITS:				
	VOLATILES					
71-55-6	1,1,1-TRICHLOROETHANE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/kg	0.54 U	0.82 UJ	0.73 UJ	0.75 UJ
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/kg	1 U	1.6 U	1.4 U	1.5 U
75-34-3	1,1-DICHLOROETHANE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
75-35-4	1,1-DICHLOROETHENE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/kg	0.58 U	0.9 UJ	0.79 UJ	0.82 UJ
120-82-1	1,2,4-TRICHLOROBENZENE	ug/kg	0.58 U	0.9 UJ	0.79 UJ	0.82 UJ
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/kg	1 U	1.6 UJ	1.4 U	1.4 U
106-93-4	1,2-DIBromoETHANE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
95-50-1	1,2-DICHLOROBENZENE	ug/kg	0.58 U	0.9 UJ	0.79 U	0.82 U
107-06-2	1,2-DICHLOROETHANE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
78-87-5	1,2-DICHLOROPROPANE	ug/kg	0.3 U	0.47 U	0.41 UJ	0.42 UJ
541-73-1	1,3-DICHLOROBENZENE	ug/kg	0.43 U	0.66 UJ	0.58 U	0.6 U
106-46-7	1,4-DICHLOROBENZENE	ug/kg	0.48 U	0.74 UJ	0.65 U	0.67 U
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/kg	120 U	180 U	R	R
591-78-6	2-HEXANONE	ug/kg	2.9 U	4.5 U	3.9 U	4.1 U
67-64-1	ACETONE	ug/kg	61.6	39.1 J	52.5	4.1 U
71-43-2	BENZENE	ug/kg	0.44 U	0.68 U	13.7	0.62 U
74-97-5	BROMOCHLOROMETHANE	ug/kg	0.58 UJ	0.9 U	0.79 U	0.82 U
75-27-4	BROMODICHLOROMETHANE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
75-25-2	BROMOFORM	ug/kg	0.86 U	1.3 U	1.2 UJ	1.2 UJ
74-83-9	BROMOMETHANE	ug/kg	1.2 U	1.8 U	1.6 U	1.6 U
75-15-0	CARBON DISULFIDE	ug/kg	0.58 U	1.9 J	2.3 J	3 J
56-23-5	CARBON TETRACHLORIDE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
108-90-7	CHLOROBENZENE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
75-00-3	CHLOROETHANE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
67-66-3	CHLOROFORM	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
74-87-3	CHLOROMETHANE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
110-82-7	CYCLOHEXANE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
124-48-1	DIBROMOCHLOROMETHANE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
75-71-8	DICHLORODIFLUOROMETHANE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
100-41-4	ETHYLBENZENE	ug/kg	0.58 U	0.9 U	4.5 J	0.82 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/kg	0.56 U	0.86 UJ	2.5 J	0.78 U
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/kg	0.84 U	1.3 U	26.6	1.2 U
79-20-9	METHYL ACETATE	ug/kg	1.2 U	1.8 U	1.6 UJ	1.6 UJ
78-93-3	METHYL ETHYL KETONE	ug/kg	3.6 U	5.6 U	4.9 U	5.1 U
108-10-1	METHYL ISOBUTYL KETONE	ug/kg	2.9 U	4.5 U	6.1 J	4.1 U
108-87-2	METHYLCYCLOHEXANE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
75-09-2	METHYLENE CHLORIDE	ug/kg	5.8 U	6 J	2.5 J	2.2 J
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/kg	0.58 U	0.9 U	13	0.82 U
100-42-5	STYRENE	ug/kg	0.52 U	0.81 U	3.7 J	0.73 U
1634-04-4	TERT-BUTYL METHYL ETHER	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
108-88-3	TOLUENE	ug/kg	0.58 U	0.9 U	9.3	0.82 U
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
75-69-4	TRICHLOROFUOROMETHANE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U
75-01-4	VINYL CHLORIDE	ug/kg	0.58 U	0.9 U	0.79 U	0.82 U

		Dup of SB-7(8-10)-20131031				
Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-7 SB-17(8-10)-20131031 E4277-16	SB-7 SB-7(16-18)-20131031 E4277-11	SB-8 SB-8(15-16)-20131030 E4277-07	SB-8 SB-8(17-19)-20131030 E4277-08
CAS NO.	COMPOUND	UNITS:				
	SEMIVOLATILES					
95-94-3	1,2,4,5-TETRACHLOROBENZENE	ug/kg	15.3 U	18.3 U	19.7 U	20.1 U
58-90-2	2,3,4,6-TETRACHLOROPHENOL	ug/kg	15.3 U	18.3 U	19.7 U	20.1 U
95-95-4	2,4,5-TRICHLOROPHENOL	ug/kg	27.3 U	32.7 U	35.1 U	36 U
88-06-2	2,4,6-TRICHLOROPHENOL	ug/kg	11.9 U	14.2 U	15.3 U	15.7 U
120-83-2	2,4-DICHLOROPHENOL	ug/kg	14.8 U	17.7 U	19.1 U	19.5 U
105-67-9	2,4-DIMETHYLPHENOL	ug/kg	22 U	26.4 U	28.4 U	29 U
51-28-5	2,4-DINITROPHENOL	ug/kg	39.5 UJ	47.3 UJ	50.9 UJ	52.1 UJ
121-14-2	2,4-DINITROTOLUENE	ug/kg	11.7 U	14 U	15 U	15.4 U
606-20-2	2,6-DINITROTOLUENE	ug/kg	15.9 U	19 U	20.4 U	20.9 U
91-58-7	2-CHLORONAPHTHALENE	ug/kg	8.9 U	10.6 U	11.4 U	11.7 U
95-57-8	2-CHLOROPHENOL	ug/kg	20.5 U	24.6 U	26.4 U	27 U
91-57-6	2-METHYLNAPHTHALENE	ug/kg	150 J	11.7 U	1900	12.9 U
95-48-7	2-METHYLPHENOL (O-CRESOL)	ug/kg	21.1 U	25.3 U	27.2 U	27.8 U
88-74-4	2-NITROANILINE	ug/kg	17.3 U	20.7 U	22.2 U	22.7 U
88-75-5	2-NITROPHENOL	ug/kg	18.8 U	22.5 U	24.2 U	24.7 U
MEPH3ME	3- AND 4- METHYLPHENOL (TOTAL)	ug/kg	20.2 U	220 J	140 J	26.6 U
91-94-1	3,3'-DICHLOROBENZIDINE	ug/kg	25 U	29.9 U	32.1 U	32.9 U
99-09-2	3-NITROANILINE	ug/kg	25 U	29.9 U	32.1 U	32.9 U
534-52-1	4,6-DINITRO-2-METHYLPHENOL	ug/kg	22.3 UJ	26.7 UJ	28.7 U	29.4 UJ
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/kg	7.6 U	9.1 U	9.8 U	10 U
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/kg	17.3 U	20.7 U	22.2 U	22.7 U
106-47-8	4-CHLOROANILINE	ug/kg	27.4 U	32.8 U	35.3 U	36.1 U
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/kg	21.1 U	25.3 U	27.2 U	27.8 U
100-01-6	4-NITROANILINE	ug/kg	50.6 U	60.6 U	65.2 U	66.7 U
100-02-7	4-NITROPHENOL	ug/kg	72.2 U	86.5 U	92.9 U	95.1 U
83-32-9	ACENAPHTHENE	ug/kg	11 U	13.1 U	700	14.4 U
208-96-8	ACENAPHTHYLENE	ug/kg	320 J	11.7 U	2500	12.9 U
98-86-2	ACETOPHENONE	ug/kg	11.9 U	14.2 U	15.3 U	15.7 U
120-12-7	ANTHRACENE	ug/kg	580 J	200 J	3900	10.5 U
1912-24-9	ATRAZINE	ug/kg	20.5 U	24.6 U	26.4 U	27 UJ
100-52-7	BENZALDEHYDE	ug/kg	R	R	R	R
56-55-3	BENZO(A)ANTHRACENE	ug/kg	1700	990	4900	24.4 U
50-32-8	BENZO(A)PYRENE	ug/kg	1900	890	3100	11.1 U
205-99-2	BENZO(B)FLUORANTHENE	ug/kg	2000	1100	2700 J	16.8 UJ
191-24-2	BENZO(G,H,I)PERYLENE	ug/kg	1600 J	530	1900 J	20.7 U
207-08-9	BENZO(K)FLUORANTHENE	ug/kg	900 J	390 J	1900	24.1 UJ
85-68-7	BENZYL BUTYL PHTHALATE	ug/kg	18.7 U	22.3 U	24 U	24.6 U
92-52-4	BIPHENYL (DIPHENYL)	ug/kg	14.7 U	17.6 U	630	19.4 U
111-91-1	BIS(2-CHLOROETHOXY) METHANE	ug/kg	22.4 U	26.8 U	28.8 U	29.5 U
111-44-4	BIS(2-CHLOROETHYL) ETHER	ug/kg	18.7 U	22.3 U	24 U	24.6 U
108-60-1	BIS(2-CHLOROISOPROPYL) ETHER	ug/kg	16.1 U	19.3 U	20.7 U	21.2 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/kg	110 J	16.5 U	17.7 U	18.1 U
105-60-2	CAPROLACTAM	ug/kg	18.1 U	21.6 U	23.3 U	23.8 U
86-74-8	CARBAZOLE	ug/kg	130 J	10.2 U	970 J	11.2 U
218-01-9	CHRYSENE	ug/kg	1500	890	2700	23.2 U
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/kg	250 J	140 J	760	14.8 U
132-64-9	DIBENZOFURAN	ug/kg	84.7 J	18.2 U	1400	20 U
84-66-2	DIETHYL PHTHALATE	ug/kg	6.1 U	7.3 U	7.8 U	8 U
131-11-3	DIMETHYL PHTHALATE	ug/kg	610 J	430 J	720	610
84-74-2	DI-N-BUTYL PHTHALATE	ug/kg	30.6 U	36.6 U	39.3 U	40.3 U
117-84-0	DI-N-OCTYLPHTHALATE	ug/kg	4.4 UJ	5.3 U	5.7 U	5.8 U
206-44-0	FLUORANTHENE	ug/kg	2700	1300	9200	10.3 U
86-73-7	FLUORENE	ug/kg	250 J	17.6 U	3200	19.4 U
118-74-1	HEXACHLOROBENZENE	ug/kg	15.9 U	19 U	20.4 U	20.9 U
87-68-3	HEXACHLOROBUTADIENE	ug/kg	14.1 U	16.9 U	18.2 U	18.6 U
77-47-4	HEXACHLOROCYCLOPENTADIENE	ug/kg	9.4 UJ	11.3 U	12.2 U	12.4 U
67-72-1	HEXACHLOROETHANE	ug/kg	17.4 U	20.8 U	22.4 U	22.9 U
193-39-5	INDENO(1,2,3-C,D)PYRENE	ug/kg	1300	490	1600	17.1 U
78-59-1	ISOPHORONE	ug/kg	12.8 U	15.4 U	16.5 U	16.9 U

		Dup of SB-7(8-10)-20131031				
Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-7 SB-17(8-10)-20131031 E4277-16 8 - 10 ft CTECH E4277 SOIL 10/31/2013 8:45 2/4/2014	SB-7 SB-7(16-18)-20131031 E4277-11 16 - 18 ft CTECH E4277 SOIL 10/31/2013 10:20 2/4/2014	SB-8 SB-8(15-16)-20131030 E4277-07 15 - 16 ft CTECH E4277 SOIL 10/30/2013 14:55 2/4/2014	SB-8 SB-8(17-19)-20131030 E4277-08 17 - 19 ft CTECH E4277 SOIL 10/30/2013 14:55 2/4/2014
CAS NO.	COMPOUND	UNITS:				
	SEMIVOLATILES					
91-20-3	NAPHTHALENE	ug/kg	270 J	16.1 U	6500	17.7 U
98-95-3	NITROBENZENE	ug/kg	14.7 U	17.6 U	18.9 U	19.4 U
621-64-7	N-NITROSODI-N-PROPYLAMINE	ug/kg	19.6 U	23.5 U	25.2 U	25.8 U
86-30-6	N-NITROSODIPHENYLAMINE	ug/kg	9.3 U	11.2 U	12 U	12.3 U
87-86-5	PENTACHLOROPHENOL	ug/kg	26.6 U	31.8 U	34.2 U	35 U
85-01-8	PHENANTHRENE	ug/kg	1700	510	15000	13.8 U
108-95-2	PHENOL	ug/kg	9 UJ	10.8 U	140 J	11.8 U
129-00-0	PYRENE	ug/kg	2800	1400	9700	12.3 U
	INORGANICS					
7429-90-5	ALUMINUM	mg/kg	7500	2400	4100	2000
7440-36-0	ANTIMONY	mg/kg	1.11 J	3.91 J	4.33 J	3.23 J
7440-38-2	ARSENIC	mg/kg	5.98	12.8	13.4	5.75
7440-39-3	BARIUM	mg/kg	201	240	275	157
7440-41-7	BERYLLIUM	mg/kg	0.36	0.097 J	0.226 J	0.183 J
7440-43-9	CADMIUM	mg/kg	2.58	5.08	1.42	0.081 U
7440-70-2	CALCIUM	mg/kg	9800	14200	6500	18900
7440-47-3	CHROMIUM, TOTAL	mg/kg	21.9	17	20.5	19.1
7440-48-4	COBALT	mg/kg	10.81	6.35	7.93	7.17
7440-50-8	COPPER	mg/kg	126	414	390	63.6
7439-89-6	IRON	mg/kg	28900	52900	24000	21000
7439-92-1	LEAD	mg/kg	283 J	719 J	884 J	261 J
7439-95-4	MAGNESIUM	mg/kg	5100	1400	2500	3900
7439-96-5	MANGANESE	mg/kg	342	336	200	185
7439-97-6	MERCURY	mg/kg	0.454	1.12	2.18	0.131
7440-02-0	NICKEL	mg/kg	37.5 J	212 J	96.2 J	20.1 J
7440-09-7	POTASSIUM	mg/kg	2400 J	419 J	799 J	410 J
7782-49-2	SELENIUM	mg/kg	0.856 J	2.62	2.35	1.66
7440-22-4	SILVER	mg/kg	0.99	1.86	2.08	1.58
7440-23-5	SODIUM	mg/kg	245	684	387	460
7440-28-0	THALLIUM	mg/kg	0.259 U	0.32 U	0.35 U	0.363 U
7440-62-2	VANADIUM	mg/kg	50.9	20.8	24.1	23.1
7440-66-6	ZINC	mg/kg	372	719	3200	197
57-12-5	CYANIDE	mg/kg	1.3 J	1.78	2	0.109 J

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-9 SB-9(5-10)-20131030 E4277-05 5 - 10 ft CTECH E4277 SOIL 10/30/2013 13:30 2/4/2014	SB-9 SB-9(10-15)-20131030 E4277-06 10 - 15 ft CTECH E4277 SOIL 10/30/2013 13:40 2/4/2014	SB-10 SB-10(5-10)-20131030 E4277-03 5 - 10 ft CTECH E4277 SOIL 10/30/2013 11:20 2/4/2014	SB-10 SB-10(10-12)-20131030 E4277-04 10 - 12 ft CTECH E4277 SOIL 10/30/2013 11:30 2/4/2014
CAS NO.	COMPOUND	UNITS:				
	VOLATILES					
71-55-6	1,1,1-TRICHLOROETHANE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/kg	0.47 UJ	0.77 UJ	0.58 UJ	0.61 UJ
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/kg	0.92 U	1.5 U	1.1 U	1.2 U
75-34-3	1,1-DICHLOROETHANE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
75-35-4	1,1-DICHLOROETHENE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/kg	0.51 UJ	0.83 UJ	0.63 UJ	0.67 UJ
120-82-1	1,2,4-TRICHLOROBENZENE	ug/kg	0.51 UJ	0.83 UJ	0.63 UJ	0.67 UJ
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/kg	0.89 U	1.5 U	1.1 UJ	1.2 U
106-93-4	1,2-DIBromoETHANE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
95-50-1	1,2-DICHLOROBENZENE	ug/kg	0.51 U	0.83 U	0.63 UJ	0.67 U
107-06-2	1,2-DICHLOROETHANE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
78-87-5	1,2-DICHLOROPROPANE	ug/kg	0.27 UJ	0.43 UJ	0.33 UJ	0.35 UJ
541-73-1	1,3-DICHLOROBENZENE	ug/kg	0.38 U	0.62 U	0.47 UJ	0.49 U
106-46-7	1,4-DICHLOROBENZENE	ug/kg	0.42 U	0.68 U	0.52 UJ	0.55 U
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/kg	R	R	R	R
591-78-6	2-HEXANONE	ug/kg	2.6 U	4.2 U	3.2 U	3.3 U
67-64-1	ACETONE	ug/kg	23.6 J	81.4	47.6	18 J
71-43-2	BENZENE	ug/kg	9.7	0.63 U	0.48 U	0.51 U
74-97-5	BROMOCHLOROMETHANE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
75-27-4	BROMODICHLOROMETHANE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
75-25-2	BROMOFORM	ug/kg	0.76 UJ	1.2 UJ	0.94 UJ	0.99 UJ
74-83-9	BROMOMETHANE	ug/kg	1 U	1.7 U	1.3 U	1.3 U
75-15-0	CARBON DISULFIDE	ug/kg	3.5 J	1.7 J	2 J	2.6 J
56-23-5	CARBON TETRACHLORIDE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
108-90-7	CHLOROBENZENE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
75-00-3	CHLOROETHANE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
67-66-3	CHLOROFORM	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
74-87-3	CHLOROMETHANE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
110-82-7	CYCLOHEXANE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
124-48-1	DIBROMOCHLOROMETHANE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
75-71-8	DICHLORODIFLUOROMETHANE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
100-41-4	ETHYLBENZENE	ug/kg	4 J	0.83 U	0.63 U	0.67 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/kg	0.49 U	0.8 U	0.61 UJ	0.64 U
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/kg	76.8	1.2 U	0.91 U	0.96 U
79-20-9	METHYL ACETATE	ug/kg	1 UJ	1.7 UJ	1.3 UJ	1.3 UJ
78-93-3	METHYL ETHYL KETONE	ug/kg	3.2 U	5.2 U	3.9 U	4.2 U
108-10-1	METHYL ISOBUTYL KETONE	ug/kg	5.4 J	4.2 U	3.2 U	3.3 U
108-87-2	METHYLCYCLOHEXANE	ug/kg	1.4 J	0.83 U	0.63 U	0.67 U
75-09-2	METHYLENE CHLORIDE	ug/kg	3.5 J	2.8 J	4.8 J	5 J
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/kg	36.7	0.83 U	0.63 U	0.67 U
100-42-5	STYRENE	ug/kg	9.4	0.75 U	0.57 U	0.6 U
1634-04-4	TERT-BUTYL METHYL ETHER	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
108-88-3	TOLUENE	ug/kg	21.3	0.83 U	0.63 U	0.67 U
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
75-69-4	TRICHLOROFUOROMETHANE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U
75-01-4	VINYL CHLORIDE	ug/kg	0.51 U	0.83 U	0.63 U	0.67 U

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-9 SB-9(5-10)-20131030 E4277-05 5 - 10 ft CTECH E4277 SOIL 10/30/2013 13:30 2/4/2014	SB-9 SB-9(10-15)-20131030 E4277-06 10 - 15 ft CTECH E4277 SOIL 10/30/2013 13:40 2/4/2014	SB-10 SB-10(5-10)-20131030 E4277-03 5 - 10 ft CTECH E4277 SOIL 10/30/2013 11:20 2/4/2014	SB-10 SB-10(10-12)-20131030 E4277-04 10 - 12 ft CTECH E4277 SOIL 10/30/2013 11:30 2/4/2014
CAS NO.	COMPOUND	UNITS:				
	SIMEVOLATILES					
95-94-3	1,2,4,5-TETRACHLOROBENZENE	ug/kg	15.7 U	16.6 U	16.8 U	17.1 U
58-90-2	2,3,4,6-TETRACHLOROPHENOL	ug/kg	15.7 U	16.6 U	16.8 U	17.1 U
95-95-4	2,4,5-TRICHLOROPHENOL	ug/kg	28.1 U	29.7 U	30 U	30.5 U
88-06-2	2,4,6-TRICHLOROPHENOL	ug/kg	12.2 U	12.9 U	13.1 U	13.3 U
120-83-2	2,4-DICHLOROPHENOL	ug/kg	15.2 U	16.1 U	16.3 U	16.5 U
105-67-9	2,4-DIMETHYLPHENOL	ug/kg	110 J	24 U	24.3 U	24.6 U
51-28-5	2,4-DINITROPHENOL	ug/kg	40.6 UJ	43 UJ	43.5 UJ	44.2 UJ
121-14-2	2,4-DINITROTOLUENE	ug/kg	12 U	12.7 U	12.8 U	13 U
606-20-2	2,6-DINITROTOLUENE	ug/kg	16.3 U	17.2 U	17.5 U	17.7 U
91-58-7	2-CHLORONAPHTHALENE	ug/kg	9.1 U	9.6 U	9.8 U	9.9 U
95-57-8	2-CHLOROPHENOL	ug/kg	21.1 U	22.3 U	22.6 U	22.9 U
91-57-6	2-METHYLNAPHTHALENE	ug/kg	1300	10.6 U	120 J	10.9 U
95-48-7	2-METHYLPHENOL (O-CRESOL)	ug/kg	120 J	22.9 U	23.2 U	23.6 U
88-74-4	2-NITROANILINE	ug/kg	17.7 U	18.8 U	19 U	19.3 U
88-75-5	2-NITROPHENOL	ug/kg	19.3 U	20.4 U	20.7 U	21 U
MEPH3ME	3- AND 4- METHYLPHENOL (TOTAL)	ug/kg	330 J	21.9 U	22.2 U	22.5 U
91-94-1	3,3'-DICHLOROBENZIDINE	ug/kg	25.7 U	27.1 U	27.5 U	27.9 U
99-09-2	3-NITROANILINE	ug/kg	25.7 U	27.1 U	27.5 U	27.9 U
534-52-1	4,6-DINITRO-2-METHYLPHENOL	ug/kg	22.9 U	24.2 U	24.5 U	24.9 U
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/kg	7.8 U	8.2 U	8.3 U	8.5 U
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/kg	17.7 U	18.8 U	19 U	19.3 U
106-47-8	4-CHLOROANILINE	ug/kg	28.2 U	29.8 U	30.2 U	30.6 U
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/kg	21.7 U	22.9 U	23.2 U	23.6 U
100-01-6	4-NITROANILINE	ug/kg	52 U	55 U	55.7 U	56.5 U
100-02-7	4-NITROPHENOL	ug/kg	74.2 U	78.4 U	79.5 U	80.6 U
83-32-9	ACENAPHTHENE	ug/kg	470	11.9 U	12.1 U	210 J
208-96-8	ACENAPHTHYLENE	ug/kg	3600	240 J	170 J	94.2 J
98-86-2	ACETOPHENONE	ug/kg	12.2 U	12.9 U	13.1 U	13.3 U
120-12-7	ANTHRACENE	ug/kg	13500	480	160 J	920
1912-24-9	ATRAZINE	ug/kg	21.1 U	22.3 U	22.6 U	22.9 U
100-52-7	BENZALDEHYDE	ug/kg	R	R	99.3 J	R
56-55-3	BENZO(A)ANTHRACENE	ug/kg	22000	950	380 J	1400
50-32-8	BENZO(A)PYRENE	ug/kg	12300	770	310 J	920
205-99-2	BENZO(B)FLUORANTHENE	ug/kg	15300 J	860 J	310 J	1100 J
191-24-2	BENZO(G,H,I)PERYLENE	ug/kg	6800 J	760 J	210 J	500 J
207-08-9	BENZO(K)FLUORANTHENE	ug/kg	4900	290 J	180 J	500
85-68-7	BENZYL BUTYL PHTHALATE	ug/kg	19.2 U	20.3 U	20.5 U	20.8 U
92-52-4	BIPHENYL (DIPHENYL)	ug/kg	400	16 U	16.2 U	16.4 U
111-91-1	BIS(2-CHLOROETHoxy) METHANE	ug/kg	23 U	24.3 U	24.7 U	25 U
111-44-4	BIS(2-CHLOROETHYL) ETHER	ug/kg	19.2 U	20.3 U	20.5 U	20.8 U
108-60-1	BIS(2-CHLOROISOPROPYL) ETHER	ug/kg	16.5 U	17.5 U	17.7 U	18 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/kg	14.1 U	15 U	15.2 U	15.4 U
105-60-2	CAPROLACTAM	ug/kg	18.6 U	19.6 U	19.9 U	20.2 U
86-74-8	CARBAZOLE	ug/kg	4100 J	110 J	9.4 U	130 J
218-01-9	CHRYSENE	ug/kg	12900	750	360 J	1100
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/kg	3000	190 J	12.3 U	240 J
132-64-9	DIBENZOFURAN	ug/kg	2200	16.5 U	16.7 U	130 J
84-66-2	DIETHYL PHTHALATE	ug/kg	6.2 U	6.6 U	6.7 U	6.8 U
131-11-3	DIMETHYL PHTHALATE	ug/kg	410	590	590	540
84-74-2	DI-N-BUTYL PHTHALATE	ug/kg	31.4 U	33.2 U	33.6 U	34.1 U
117-84-0	DI-N-OCTYLPHthalate	ug/kg	4.6 U	4.8 U	4.9 U	5 U
206-44-0	FLUORANTHENE	ug/kg	46700	1700	570	2600
86-73-7	FLUORENE	ug/kg	5200	230 J	16.2 U	310 J
118-74-1	HEXACHLOROBENZENE	ug/kg	16.3 U	17.2 U	17.5 U	17.7 U
87-68-3	HEXACHLOROBUTADIENE	ug/kg	14.5 U	15.3 U	15.5 U	15.8 U
77-47-4	HEXACHLOROCYCLOPENTADIENE	ug/kg	9.7 U	10.3 U	10.4 U	10.6 U
67-72-1	HEXACHLOROETHANE	ug/kg	17.9 U	18.9 U	19.1 U	19.4 U
193-39-5	INDENO(1,2,3-C,D)PYRENE	ug/kg	7300	550	180 J	490
78-59-1	ISOPHORONE	ug/kg	13.2 U	13.9 U	14.1 U	14.3 U

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-9 SB-9(5-10)-20131030 E4277-05 5 - 10 ft CTECH E4277 SOIL 10/30/2013 13:30 2/4/2014	SB-9 SB-9(10-15)-20131030 E4277-06 10 - 15 ft CTECH E4277 SOIL 10/30/2013 13:40 2/4/2014	SB-10 SB-10(5-10)-20131030 E4277-03 5 - 10 ft CTECH E4277 SOIL 10/30/2013 11:20 2/4/2014	SB-10 SB-10(10-12)-20131030 E4277-04 10 - 12 ft CTECH E4277 SOIL 10/30/2013 11:30 2/4/2014
CAS NO.	COMPOUND	UNITS:				
	SEMIVOLATILES					
91-20-3	NAPHTHALENE	ug/kg	2300	200 J	150 J	15 U
98-95-3	NITROBENZENE	ug/kg	15.1 U	16 U	16.2 U	16.4 U
621-64-7	N-NITROSODI-N-PROPYLAMINE	ug/kg	20.1 U	21.3 U	21.6 U	21.9 U
86-30-6	N-NITROSODIPHENYLAMINE	ug/kg	9.6 U	10.1 U	10.3 U	10.4 U
87-86-5	PENTACHLOROPHENOL	ug/kg	27.3 U	28.9 U	29.3 U	29.7 U
85-01-8	PHENANTHRENE	ug/kg	39000	1200	470	2100
108-95-2	PHENOL	ug/kg	240 J	9.8 U	9.9 U	10 U
129-00-0	PYRENE	ug/kg	33200	1700	650	2200
	INORGANICS					
7429-90-5	ALUMINUM	mg/kg	7600	5200	5400	7800
7440-36-0	ANTIMONY	mg/kg	1.19 J	1.15 J	0.996 J	0.732 J
7440-38-2	ARSENIC	mg/kg	7.18	6.71	7.74	4.85
7440-39-3	BARIUM	mg/kg	136	99.1	137	135
7440-41-7	BERYLLIUM	mg/kg	0.296 J	0.323	0.308 J	0.25 J
7440-43-9	CADMIUM	mg/kg	0.405	0.862	0.064 U	0.14 J
7440-70-2	CALCIUM	mg/kg	21500	49900	5000	22900
7440-47-3	CHROMIUM, TOTAL	mg/kg	21.9	17.2	13.1	19.3
7440-48-4	COBALT	mg/kg	20.9	9.24	5.8	9.87
7440-50-8	COPPER	mg/kg	72	70.3	84.1	56.3
7439-89-6	IRON	mg/kg	32600	17000	17700	22200
7439-92-1	LEAD	mg/kg	232 J	161 J	231 J	173 J
7439-95-4	MAGNESIUM	mg/kg	7600	18200	2500	7600
7439-96-5	MANGANESE	mg/kg	237	176	483	229
7439-97-6	MERCURY	mg/kg	0.826	0.573	0.373	0.303
7440-02-0	NICKEL	mg/kg	43.2 J	32.3 J	18.6 J	26.5 J
7440-09-7	POTASSIUM	mg/kg	2200 J	912 J	801 J	1800 J
7782-49-2	SELENIUM	mg/kg	2.53	1.27	2.09	1.2
7440-22-4	SILVER	mg/kg	1.87	1.41	1.3	1.34
7440-23-5	SODIUM	mg/kg	279	258	481	623
7440-28-0	THALLIUM	mg/kg	0.275 U	0.29 U	0.29 U	0.288 U
7440-62-2	VANADIUM	mg/kg	88	25.6	32.5	34.8
7440-66-6	ZINC	mg/kg	210	246	233	151
57-12-5	CYANIDE	mg/kg	5.36	0.104 J	0.609	0.522

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-11 SB-11(5-10)-20131030 E4277-01 5 - 10 ft CTECH E4277 SOIL 10/30/2013 9:45 2/4/2014	SB-11 SB-11(15-20)-20131030 E4277-02 15 - 20 ft CTECH E4277 SOIL 10/30/2013 10:05 2/4/2014	FIELDQC TB110113-20131101 E4277-28 -	FIELDQC TB110413-20131104 E4340-07 -
CAS NO.	COMPOUND	UNITS:			11/1/2013 10:00 2/4/2014 ug/l	11/4/2013 10:00 2/4/2014 ug/l
	VOLATILES					
71-55-6	1,1,1-TRICHLOROETHANE	ug/kg	0.61 U	0.79 U	0.4 U	0.4 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/kg	0.56 UJ	0.73 UJ	0.31 U	0.31 U
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/kg	0.61 U	0.79 U	0.45 U	0.45 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/kg	1.1 U	1.4 U	0.38 U	0.38 U
75-34-3	1,1-DICHLOROETHANE	ug/kg	0.61 U	0.79 U	0.36 UJ	0.36 UJ
75-35-4	1,1-DICHLOROETHENE	ug/kg	0.61 U	0.79 U	0.47 U	0.47 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/kg	0.61 UJ	0.79 UJ	0.2 U	0.2 U
120-82-1	1,2,4-TRICHLOROBENZENE	ug/kg	0.61 UJ	0.79 UJ	0.2 U	0.2 U
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/kg	1.1 UJ	1.4 U	0.46 UJ	0.46 UJ
106-93-4	1,2-DIBROMOETHANE	ug/kg	0.61 U	0.79 U	0.41 UJ	0.41 UJ
95-50-1	1,2-DICHLOROBENZENE	ug/kg	0.61 UJ	0.79 U	0.45 U	0.45 U
107-06-2	1,2-DICHLOROETHANE	ug/kg	0.61 U	0.79 U	0.48 U	0.48 U
78-87-5	1,2-DICHLOROPROPANE	ug/kg	0.32 UJ	0.41 UJ	0.46 U	0.46 U
541-73-1	1,3-DICHLOROBENZENE	ug/kg	0.45 UJ	0.58 U	0.43 U	0.43 U
106-46-7	1,4-DICHLOROBENZENE	ug/kg	0.5 UJ	0.65 U	0.32 U	0.32 U
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/kg	R	R	50 U	50 U
591-78-6	2-HEXANONE	ug/kg	3 U	4 U	1.9 UJ	1.9 UJ
67-64-1	ACETONE	ug/kg	29.5 J	36.2 J	0.5 U	0.5 U
71-43-2	BENZENE	ug/kg	0.46 U	0.6 U	0.32 U	0.32 U
74-97-5	BROMOCHLOROMETHANE	ug/kg	0.61 U	0.79 U	0.2 U	0.2 U
75-27-4	BROMODICHLOROMETHANE	ug/kg	0.61 U	0.79 U	0.36 U	0.36 U
75-25-2	BROMOFORM	ug/kg	0.9 UJ	1.2 UJ	0.47 UJ	0.47 UJ
74-83-9	BROMOMETHANE	ug/kg	1.2 U	1.6 U	0.2 U	0.2 U
75-15-0	CARBON DISULFIDE	ug/kg	0.61 U	3.6 J	0.2 U	0.2 U
56-23-5	CARBON TETRACHLORIDE	ug/kg	0.61 U	0.79 U	0.2 U	0.2 U
108-90-7	CHLOROBENZENE	ug/kg	0.61 U	0.79 U	0.49 U	0.49 U
75-00-3	CHLOROETHANE	ug/kg	0.61 U	0.79 U	0.2 UJ	0.2 UJ
67-66-3	CHLOROFORM	ug/kg	0.61 U	0.79 U	0.34 U	0.34 U
74-87-3	CHLOROMETHANE	ug/kg	0.61 U	0.79 U	0.2 U	0.2 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/kg	0.61 U	0.79 U	0.35 U	0.35 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/kg	0.61 U	0.79 U	0.31 U	0.31 U
110-82-7	CYCLOHEXANE	ug/kg	0.61 U	0.79 U	0.2 U	0.2 U
124-48-1	DIBROMOCHLOROMETHANE	ug/kg	0.61 U	0.79 U	0.2 UJ	0.2 UJ
75-71-8	DICHLORODIFLUOROMETHANE	ug/kg	0.61 U	0.79 U	0.2 U	0.2 U
100-41-4	ETHYLBENZENE	ug/kg	0.61 U	0.79 U	0.2 U	0.2 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/kg	0.59 UJ	0.76 U	0.45 U	0.45 U
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/kg	0.88 U	1.1 U	0.95 U	0.95 U
79-20-9	METHYL ACETATE	ug/kg	1.2 UJ	1.6 UJ	0.2 U	0.2 U
78-93-3	METHYL ETHYL KETONE	ug/kg	3.8 U	4.9 U	1.3 U	1.3 U
108-10-1	METHYL ISOBUTYL KETONE	ug/kg	3 U	4 U	2.1 U	2.1 U
108-87-2	METHYLCYCLOHEXANE	ug/kg	0.61 U	0.79 U	0.2 U	0.2 U
75-09-2	METHYLENE CHLORIDE	ug/kg	4.2 J	5.6 J	0.41 U	0.41 U
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/kg	0.61 U	0.79 U	0.43 U	0.43 U
100-42-5	STYRENE	ug/kg	0.55 U	0.71 U	0.36 U	0.36 U
1634-04-4	TERT-BUTYL METHYL ETHER	ug/kg	0.61 U	0.79 U	0.35 U	0.35 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/kg	0.61 U	0.79 U	0.27 U	0.27 U
108-88-3	TOLUENE	ug/kg	0.61 U	0.79 U	0.37 U	0.37 U
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/kg	0.61 U	0.79 U	0.41 U	0.41 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/kg	0.61 U	0.79 U	0.29 U	0.29 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/kg	0.61 U	0.79 U	0.28 U	0.28 U
75-69-4	TRICHLOROFUOROMETHANE	ug/kg	0.61 U	0.79 U	0.35 U	0.35 U
75-01-4	VINYL CHLORIDE	ug/kg	0.61 U	0.79 U	0.34 UJ	0.34 UJ

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-11 SB-11(5-10)-20131030 E4277-01 5 - 10 ft CTECH E4277 SOIL 10/30/2013 9:45 2/4/2014	SB-11 SB-11(15-20)-20131030 E4277-02 15 - 20 ft CTECH E4277 SOIL 10/30/2013 10:05 2/4/2014	FIELDQC TB110113-20131101 E4277-28 -	FIELDQC TB110413-20131104 E4340-07 -
CAS NO.	COMPOUND	UNITS:			11/1/2013 10:00 2/4/2014 ug/l	11/4/2013 10:00 2/4/2014 ug/l
	SEMIVOLATILES					
95-94-3	1,2,4,5-TETRACHLOROBENZENE	ug/kg	16.3 U	19.5 U		
58-90-2	2,3,4,6-TETRACHLOROPHENOL	ug/kg	16.3 U	19.5 U		
95-95-4	2,4,5-TRICHLOROPHENOL	ug/kg	29.2 U	34.9 U		
88-06-2	2,4,6-TRICHLOROPHENOL	ug/kg	12.7 U	15.2 U		
120-83-2	2,4-DICHLOROPHENOL	ug/kg	15.8 U	18.9 U		
105-67-9	2,4-DIMETHYLPHENOL	ug/kg	23.6 U	28.2 U		
51-28-5	2,4-DINITROPHENOL	ug/kg	42.3 UJ	50.6 UJ		
121-14-2	2,4-DINITROTOLUENE	ug/kg	12.5 U	14.9 U		
606-20-2	2,6-DINITROTOLUENE	ug/kg	17 U	20.3 U		
91-58-7	2-CHLORONAPHTHALENE	ug/kg	9.5 U	11.3 U		
95-57-8	2-CHLOROPHENOL	ug/kg	22 U	26.3 U		
91-57-6	2-METHYLNAPHTHALENE	ug/kg	10.5 U	12.5 U		
95-48-7	2-METHYLPHENOL (O-CRESOL)	ug/kg	22.6 U	27 U		
88-74-4	2-NITROANILINE	ug/kg	18.5 U	22.1 U		
88-75-5	2-NITROPHENOL	ug/kg	20.1 U	24 U		
MEPH3ME	3- AND 4- METHYLPHENOL (TOTAL)	ug/kg	21.6 U	25.8 U		
91-94-1	3,3'-DICHLOROBENZIDINE	ug/kg	26.7 U	31.9 U		
99-09-2	3-NITROANILINE	ug/kg	26.7 U	31.9 U		
534-52-1	4,6-DINITRO-2-METHYLPHENOL	ug/kg	23.8 U	28.5 U		
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/kg	8.1 U	9.7 U		
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/kg	18.5 U	22.1 U		
106-47-8	4-CHLOROANILINE	ug/kg	29.3 U	35.1 U		
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/kg	22.6 U	27 U		
100-01-6	4-NITROANILINE	ug/kg	54.2 U	64.7 U		
100-02-7	4-NITROPHENOL	ug/kg	77.2 U	92.3 U		
83-32-9	ACENAPHTHENE	ug/kg	11.7 U	14 U		
208-96-8	ACENAPHTHYLENE	ug/kg	10.5 U	12.5 U		
98-86-2	ACETOPHENONE	ug/kg	12.7 U	15.2 U		
120-12-7	ANTHRACENE	ug/kg	140 J	120 J		
1912-24-9	ATRAZINE	ug/kg	22 U	26.3 U		
100-52-7	BENZALDEHYDE	ug/kg	R	R		
56-55-3	BENZO(A)ANTHRACENE	ug/kg	720	240 J		
50-32-8	BENZO(A)PYRENE	ug/kg	660	190 J		
205-99-2	BENZO(B)FLUORANTHENE	ug/kg	760 J	180 J		
191-24-2	BENZO(G,H,I)PERYLENE	ug/kg	480 J	120 J		
207-08-9	BENZO(K)FLUORANTHENE	ug/kg	300 J	23.4 U		
85-68-7	BENZYL BUTYL PHTHALATE	ug/kg	20 U	23.9 U		
92-52-4	BIPHENYL (DIPHENYL)	ug/kg	15.7 U	18.8 U		
111-91-1	BIS(2-CHLOROETHOXY) METHANE	ug/kg	24 U	28.6 U		
111-44-4	BIS(2-CHLOROETHYL) ETHER	ug/kg	20 U	23.9 U		
108-60-1	BIS(2-CHLOROISOPROPYL) ETHER	ug/kg	17.2 U	20.6 U		
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/kg	88.6 J	17.6 U		
105-60-2	CAPROLACTAM	ug/kg	19.3 U	23.1 U		
86-74-8	CARBAZOLE	ug/kg	9.1 U	10.9 U		
218-01-9	CHRYSENE	ug/kg	500	170 J		
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/kg	190 J	14.3 U		
132-64-9	DIBENZOFURAN	ug/kg	16.2 U	19.4 U		
84-66-2	DIETHYL PHTHALATE	ug/kg	6.5 U	7.8 U		
131-11-3	DIMETHYL PHTHALATE	ug/kg	650	720		
84-74-2	DI-N-BUTYL PHTHALATE	ug/kg	32.7 U	39.1 U		
117-84-0	DI-N-OCTYLPHthalate	ug/kg	4.7 U	5.7 U		
206-44-0	FLUORANTHENE	ug/kg	820	370 J		
86-73-7	FLUORENE	ug/kg	15.7 U	18.8 U		
118-74-1	HEXACHLOROBENZENE	ug/kg	17 U	20.3 U		
87-68-3	HEXACHLOROBUTADIENE	ug/kg	15.1 U	18 U		
77-47-4	HEXACHLOROCYCLOPENTADIENE	ug/kg	10.1 U	12.1 U		
67-72-1	HEXACHLOROETHANE	ug/kg	18.6 U	22.2 U		
193-39-5	INDENO(1,2,3-C,D)PYRENE	ug/kg	420	99.9 J		
78-59-1	ISOPHORONE	ug/kg	13.7 U	16.4 U		

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	SB-11 SB-11(5-10)-20131030 E4277-01 5 - 10 ft CTECH E4277 SOIL 10/30/2013 9:45 2/4/2014	SB-11 SB-11(15-20)-20131030 E4277-02 15 - 20 ft CTECH E4277 SOIL 10/30/2013 10:05 2/4/2014	FIELDQC TB110113-20131101 E4277-28 - CTECH E4277 SOIL 11/1/2013 10:00 2/4/2014 ug/l	FIELDQC TB110413-20131104 E4340-07 - CTECH E4340 WATER 11/4/2013 10:00 2/4/2014 ug/l
CAS NO.	COMPOUND	UNITS:				
	SEMIVOLATILES					
91-20-3	NAPHTHALENE	ug/kg	14.3 U	17.2 U		
98-95-3	NITROBENZENE	ug/kg	15.7 U	18.8 U		
621-64-7	N-NITROSODI-N-PROPYLAMINE	ug/kg	21 U	25.1 U		
86-30-6	N-NITROSODIPHENYLAMINE	ug/kg	10 U	11.9 U		
87-86-5	PENTACHLOROPHENOL	ug/kg	28.5 U	34 U		
85-01-8	PHENANTHRENE	ug/kg	410	380 J		
108-95-2	PHENOL	ug/kg	9.6 U	11.5 U		
129-00-0	PYRENE	ug/kg	830	420 J		
	INORGANICS					
7429-90-5	ALUMINUM	mg/kg	4300	9600		
7440-36-0	ANTIMONY	mg/kg	1.5 J	2.26 J		
7440-38-2	ARSENIC	mg/kg	5.77	20.8		
7440-39-3	BARIUM	mg/kg	237	207		
7440-41-7	BERYLLIUM	mg/kg	0.348	0.271 J		
7440-43-9	CADMIUM	mg/kg	0.063 U	0.839		
7440-70-2	CALCIUM	mg/kg	4300	7300		
7440-47-3	CHROMIUM, TOTAL	mg/kg	12.6	34		
7440-48-4	COBALT	mg/kg	11.7	13.9		
7440-50-8	COPPER	mg/kg	346	130		
7439-89-6	IRON	mg/kg	22300	77700		
7439-92-1	LEAD	mg/kg	884 J	338 J		
7439-95-4	MAGNESIUM	mg/kg	1300	4700		
7439-96-5	MANGANESE	mg/kg	128	564		
7439-97-6	MERCURY	mg/kg	0.709	2.67		
7440-02-0	NICKEL	mg/kg	14.5 J	33.2 J		
7440-09-7	POTASSIUM	mg/kg	871 J	2300 J		
7782-49-2	SELENIUM	mg/kg	1.97	3.85		
7440-22-4	SILVER	mg/kg	1.64	5.38		
7440-23-5	SODIUM	mg/kg	508	2300		
7440-28-0	THALLIUM	mg/kg	0.283 U	0.353 U		
7440-62-2	VANADIUM	mg/kg	19.4	41		
7440-66-6	ZINC	mg/kg	112	288		
57-12-5	CYANIDE	mg/kg	8.79	9.13		

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	FIELDQC FB110113-20131101 E4277-27 - CTECH E4277 SOIL 11/1/2013 14:20 2/4/2014	FIELDQC FB110413-20131104 E4340-06 - CTECH E4340 SOIL 11/4/2013 11:45 2/4/2014
CAS NO.	COMPOUND	UNITS:	ug/l	ug/l
	VOLATILES			
71-55-6	1,1,1-TRICHLOROETHANE	ug/kg	0.4 U	0.4 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/kg	0.31 U	0.31 U
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/kg	0.45 U	0.45 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/kg	0.38 U	0.38 U
75-34-3	1,1-DICHLOROETHANE	ug/kg	0.36 UJ	0.36 UJ
75-35-4	1,1-DICHLOROETHENE	ug/kg	0.47 U	0.47 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/kg	0.2 U	0.2 U
120-82-1	1,2,4-TRICHLOROBENZENE	ug/kg	0.2 U	0.2 U
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/kg	0.46 UJ	0.46 UJ
106-93-4	1,2-DIBROMOETHANE	ug/kg	0.41 UJ	0.41 UJ
95-50-1	1,2-DICHLOROBENZENE	ug/kg	0.45 U	0.45 U
107-06-2	1,2-DICHLOROETHANE	ug/kg	0.48 U	0.48 U
78-87-5	1,2-DICHLOROPROPANE	ug/kg	0.46 U	0.46 U
541-73-1	1,3-DICHLOROBENZENE	ug/kg	0.43 U	0.43 U
106-46-7	1,4-DICHLOROBENZENE	ug/kg	0.32 U	0.32 U
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/kg	50 U	50 U
591-78-6	2-HEXANONE	ug/kg	1.9 UJ	1.9 UJ
67-64-1	ACETONE	ug/kg	0.5 U	0.5 U
71-43-2	BENZENE	ug/kg	0.32 U	0.32 U
74-97-5	BROMOCHLOROMETHANE	ug/kg	0.2 U	0.2 U
75-27-4	BROMODICHLOROMETHANE	ug/kg	0.36 U	0.36 U
75-25-2	BROMOFORM	ug/kg	0.47 UJ	0.47 UJ
74-83-9	BROMOMETHANE	ug/kg	0.2 U	0.2 U
75-15-0	CARBON DISULFIDE	ug/kg	0.2 U	0.2 U
56-23-5	CARBON TETRACHLORIDE	ug/kg	0.2 U	0.2 U
108-90-7	CHLOROBENZENE	ug/kg	0.49 U	0.49 U
75-00-3	CHLOROETHANE	ug/kg	0.2 UJ	0.2 UJ
67-66-3	CHLOROFORM	ug/kg	0.34 U	0.34 U
74-87-3	CHLOROMETHANE	ug/kg	0.2 U	0.2 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/kg	0.35 U	0.35 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/kg	0.31 U	0.31 U
110-82-7	CYCLOHEXANE	ug/kg	0.2 U	0.2 U
124-48-1	DIBROMOCHLOROMETHANE	ug/kg	0.2 UJ	0.2 UJ
75-71-8	DICHLORODIFLUOROMETHANE	ug/kg	0.2 U	0.2 U
100-41-4	ETHYLBENZENE	ug/kg	0.2 U	0.2 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/kg	0.45 U	0.45 U
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/kg	0.95 U	0.95 U
79-20-9	METHYL ACETATE	ug/kg	0.2 U	0.2 U
78-93-3	METHYL ETHYL KETONE	ug/kg	1.3 U	1.3 U
108-10-1	METHYL ISOBUTYL KETONE	ug/kg	2.1 U	2.1 U
108-87-2	METHYLCYCLOHEXANE	ug/kg	0.2 U	0.2 U
75-09-2	METHYLENE CHLORIDE	ug/kg	0.41 U	0.41 U
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/kg	0.43 U	0.43 U
100-42-5	STYRENE	ug/kg	0.36 U	0.36 U
1634-04-4	TERT-BUTYL METHYL ETHER	ug/kg	0.35 U	0.35 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/kg	0.27 U	0.27 U
108-88-3	TOLUENE	ug/kg	0.37 U	0.37 U
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/kg	0.41 U	0.41 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/kg	0.29 U	0.29 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/kg	0.28 U	0.28 U
75-69-4	TRICHLOROFUOROMETHANE	ug/kg	0.35 U	0.35 U
75-01-4	VINYL CHLORIDE	ug/kg	0.34 UJ	0.34 UJ

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	FIELDQC FB110113-20131101 E4277-27 - CTECH E4277 SOIL 11/1/2013 14:20 2/4/2014	FIELDQC FB110413-20131104 E4340-06 - CTECH E4340 SOIL 11/4/2013 11:45 2/4/2014
CAS NO.	COMPOUND	UNITS:	ug/l	ug/l
	SEMICVOLATILES			
95-94-3	1,2,4,5-TETRACHLOROBENZENE	ug/kg	0.2 U	0.2 U
58-90-2	2,3,4,6-TETRACHLOROPHENOL	ug/kg	0.2 U	0.2 U
95-95-4	2,4,5-TRICHLOROPHENOL	ug/kg	0.4 U	0.4 U
88-06-2	2,4,6-TRICHLOROPHENOL	ug/kg	0.57 U	0.56 U
120-83-2	2,4-DICHLOROPHENOL	ug/kg	0.67 U	0.66 U
105-67-9	2,4-DIMETHYLPHENOL	ug/kg	0.72 U	0.71 U
51-28-5	2,4-DINITROPHENOL	ug/kg	2.1 UJ	2.1 UJ
121-14-2	2,4-DINITROTOLUENE	ug/kg	1 U	1 U
606-20-2	2,6-DINITROTOLUENE	ug/kg	0.32 U	0.32 U
91-58-7	2-CHLORONAPHTHALENE	ug/kg	0.16 U	0.16 U
95-57-8	2-CHLOROPHENOL	ug/kg	0.55 U	0.54 U
91-57-6	2-METHYLNAPHTHALENE	ug/kg	0.32 U	0.32 U
95-48-7	2-METHYLPHENOL (O-CRESOL)	ug/kg	0.24 U	0.24 U
88-74-4	2-NITROANILINE	ug/kg	0.49 U	0.49 U
88-75-5	2-NITROPHENOL	ug/kg	0.53 U	0.52 U
MEPH3ME	3- AND 4- METHYLPHENOL (TOTAL)	ug/kg	0.38 U	0.38 U
91-94-1	3,3'-DICHLOROBENZIDINE	ug/kg	1 U	1 U
99-09-2	3-NITROANILINE	ug/kg	1 U	1 U
534-52-1	4,6-DINITRO-2-METHYLPHENOL	ug/kg	0.75 U	0.74 U
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/kg	0.23 U	0.23 U
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/kg	0.4 U	0.4 U
106-47-8	4-CHLOROANILINE	ug/kg	1 U	1 U
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/kg	0.21 U	0.21 U
100-01-6	4-NITROANILINE	ug/kg	1.4 U	1.4 U
100-02-7	4-NITROPHENOL	ug/kg	2 U	2 U
83-32-9	ACENAPHTHENE	ug/kg	0.21 U	0.21 U
208-96-8	ACENAPHTHYLENE	ug/kg	0.71 U	0.7 U
98-86-2	ACETOPHENONE	ug/kg	0.14 U	0.14 U
120-12-7	ANTHRACENE	ug/kg	0.16 U	0.16 U
1912-24-9	ATRAZINE	ug/kg	0.4 U	0.4 U
100-52-7	BENZALDEHYDE	ug/kg	R	R
56-55-3	BENZO(A)ANTHRACENE	ug/kg	0.16 U	0.16 U
50-32-8	BENZO(A)PYRENE	ug/kg	0.14 U	0.14 U
205-99-2	BENZO(B)FLUORANTHENE	ug/kg	0.29 U	0.29 U
191-24-2	BENZO(G,H,I)PERYLENE	ug/kg	0.29 U	0.29 U
207-08-9	BENZO(K)FLUORANTHENE	ug/kg	0.18 U	0.18 UJ
85-68-7	BENZYL BUTYL PHTHALATE	ug/kg	0.19 U	0.19 U
92-52-4	BIPHENYL (DIPHENYL)	ug/kg	0.15 U	0.15 U
111-91-1	BIS(2-CHLOROETHOXY) METHANE	ug/kg	0.56 U	0.55 U
111-44-4	BIS(2-CHLOROETHYL) ETHER	ug/kg	0.56 U	0.55 U
108-60-1	BIS(2-CHLOROISOPROPYL) ETHER	ug/kg	0.17 U	0.17 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/kg	0.16 U	0.16 U
105-60-2	CAPROLACTAM	ug/kg	1 U	1 U
86-74-8	CARBAZOLE	ug/kg	0.22 U	0.22 U
218-01-9	CHRYSENE	ug/kg	0.18 U	0.18 U
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/kg	0.42 U	0.42 U
132-64-9	DIBENZOFURAN	ug/kg	0.24 U	0.24 U
84-66-2	DIETHYL PHTHALATE	ug/kg	0.38 U	0.38 U
131-11-3	DIMETHYL PHTHALATE	ug/kg	0.22 U	0.22 U
84-74-2	DI-N-BUTYL PHTHALATE	ug/kg	1 U	1 U
117-84-0	DI-N-OCTYLPHthalate	ug/kg	0.52 U	0.51 U
206-44-0	FLUORANTHENE	ug/kg	0.4 U	0.4 U
86-73-7	FLUORENE	ug/kg	0.31 U	0.31 U
118-74-1	HEXACHLOROBENZENE	ug/kg	0.18 U	0.18 U
87-68-3	HEXACHLOROBUTADIENE	ug/kg	0.25 U	0.25 U
77-47-4	HEXACHLOROCYCLOPENTADIENE	ug/kg	0.24 U	0.24 UJ
67-72-1	HEXACHLOROETHANE	ug/kg	0.25 U	0.25 U
193-39-5	INDENO(1,2,3-C,D)PYRENE	ug/kg	0.15 U	0.15 U
78-59-1	ISOPHORONE	ug/kg	0.3 U	0.3 U

Con Ed - Halleck Street Validated Soil Analytical Data SDG: E4277 & E4340		Location ID: Sample ID: Lab Sample Id: Depth: Source: SDG: Matrix: Sampled: Validated:	FIELDQC FB110113-20131101 E4277-27 - CTECH E4277 SOIL 11/1/2013 14:20 2/4/2014	FIELDQC FB110413-20131104 E4340-06 - CTECH E4340 SOIL 11/4/2013 11:45 2/4/2014
CAS NO.	COMPOUND	UNITS:	ug/l	ug/l
	SEMIVOLATILES			
91-20-3	NAPHTHALENE	ug/kg	0.12 U	0.12 U
98-95-3	NITROBENZENE	ug/kg	0.69 UJ	0.68 U
621-64-7	N-NITROSODI-N-PROPYLAMINE	ug/kg	0.2 U	0.2 U
86-30-6	N-NITROSODIPHENYLAMINE	ug/kg	0.61 U	0.6 U
87-86-5	PENTACHLOROPHENOL	ug/kg	1 UJ	1 U
85-01-8	PHENANTHRENE	ug/kg	0.26 U	0.26 U
108-95-2	PHENOL	ug/kg	0.21 U	0.21 U
129-00-0	PYRENE	ug/kg	0.2 U	0.2 U
	INORGANICS			
7429-90-5	ALUMINUM	mg/kg	6.5 U	6.5 U
7440-36-0	ANTIMONY	mg/kg	8 U	8 U
7440-38-2	ARSENIC	mg/kg	4.2 U	4.2 U
7440-39-3	BARIUM	mg/kg	4 U	4 U
7440-41-7	BERYLLIUM	mg/kg	0.7 U	0.7 U
7440-43-9	CADMIUM	mg/kg	0.5 U	0.5 U
7440-70-2	CALCIUM	mg/kg	86.5 J	31.8 U
7440-47-3	CHROMIUM, TOTAL	mg/kg	1.1 U	1.1 U
7440-48-4	COBALT	mg/kg	5.8 U	5.8 U
7440-50-8	COPPER	mg/kg	2 U	2 U
7439-89-6	IRON	mg/kg	20.4 U	20.4 U
7439-92-1	LEAD	mg/kg	2.6 U	2.6 U
7439-95-4	MAGNESIUM	mg/kg	32.5 U	32.5 U
7439-96-5	MANGANESE	mg/kg	1.7 U	1.7 U
7439-97-6	MERCURY	mg/kg	0.1 U	0.109 J
7440-02-0	NICKEL	mg/kg	4.2 U	4.2 U
7440-09-7	POTASSIUM	mg/kg	38.8 U	38.8 U
7782-49-2	SELENIUM	mg/kg	4.8 U	4.8 U
7440-22-4	SILVER	mg/kg	1.5 U	1.5 U
7440-23-5	SODIUM	mg/kg	41.3 J	13.9 U
7440-28-0	THALLIUM	mg/kg	2.4 U	2.4 U
7440-62-2	VANADIUM	mg/kg	6.1 U	6.1 U
7440-66-6	ZINC	mg/kg	6.5 U	6.5 U
57-12-5	CYANIDE	mg/kg	3 U	3 U

**ATTACHMENT A-2**

**VALIDATED LABORATORY DATA FOR GROUNDWATER SAMPLES**

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**PARSONS**

Con Ed - Halleck Street Validated Groundwater Analytical Data SDG: E4638		Location ID: Sample ID: Lab Sample Id: Source: SDG: Matrix: Sampled: Validated:	MW-1 MW-1-20131127 E4638-06 CTECH E4638 WATER 11/27/2013 12:20 2/4/2014	MW-2 MW-2-20131127 E4638-05 CTECH E4638 WATER 11/27/2013 10:50 2/4/2014	MW-3 MW-3-20131127 E4638-01 CTECH E4638 WATER 11/27/2013 8:45 2/4/2014	MW-3 MW-3A-20131127 E4638-04 CTECH E4638 WATER 11/27/2013 9:10 2/4/2014
CAS NO.	COMPOUND	UNITS:				
	VOLATILES					
71-55-6	1,1,1-TRICHLOROETHANE	ug/l	0.4 U	0.4 U	0.4 U	0.4 U
79-34-5	1,1,2,2-TETRACHLOROETHANE	ug/l	0.31 U	0.31 U	0.31 U	0.31 U
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/l	0.45 U	0.45 U	0.45 U	0.45 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/l	0.38 U	0.38 U	0.38 U	0.38 U
75-34-3	1,1-DICHLOROETHANE	ug/l	0.36 U	0.36 U	0.36 U	0.36 U
75-35-4	1,1-DICHLOROETHENE	ug/l	0.47 U	0.47 U	0.47 U	0.47 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
120-82-1	1,2,4-TRICHLOROBENZENE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/l	0.46 U	0.46 U	0.46 U	0.46 U
106-93-4	1,2-DIBROMOETHANE	ug/l	0.41 U	0.41 U	0.41 U	0.41 U
95-50-1	1,2-DICHLOROBENZENE	ug/l	0.45 U	0.45 U	0.45 U	0.45 U
107-06-2	1,2-DICHLOROETHANE	ug/l	0.48 U	0.48 U	0.48 U	0.48 U
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/l	0.95 U	0.95 U	0.95 U	0.95 U
78-87-5	1,2-DICHLOROPROPANE	ug/l	0.46 U	0.46 U	0.46 U	0.46 U
541-73-1	1,3-DICHLOROBENZENE	ug/l	0.43 U	0.43 U	0.43 U	0.43 U
106-46-7	1,4-DICHLOROBENZENE	ug/l	0.32 U	0.32 U	0.32 U	0.32 U
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/l	50 U	50 U	50 U	50 U
591-78-6	2-HEXANONE	ug/l	1.9 U	1.9 U	1.9 U	1.9 U
67-64-1	ACETONE	ug/l	0.5 U	0.5 U	0.5 U	0.5 U
71-43-2	BENZENE	ug/l	0.32 U	0.32 U	0.32 U	0.32 U
74-97-5	BROMOCHLOROMETHANE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
75-27-4	BROMODICHLOROMETHANE	ug/l	0.36 U	0.36 U	0.36 U	0.36 U
75-25-2	BROMOFORM	ug/l	0.47 U	0.47 U	0.47 U	0.47 U
74-83-9	BROMOMETHANE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
75-15-0	CARBON DISULFIDE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
56-23-5	CARBON TETRACHLORIDE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
108-90-7	CHLOROBENZENE	ug/l	0.49 U	0.49 U	0.49 U	0.49 U
75-00-3	CHLOROETHANE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
67-66-3	CHLOROFORM	ug/l	0.34 U	0.34 U	0.34 U	0.34 U
74-87-3	CHLOROMETHANE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/l	0.35 U	0.35 U	0.35 U	0.35 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/l	0.31 U	0.31 U	0.31 U	0.31 U
110-82-7	CYCLOHEXANE	ug/l	0.2 UJ	0.2 UJ	0.2 UJ	0.2 UJ
124-48-1	DIBROMOCHLOROMETHANE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
75-71-8	DICHLORODIFLUOROMETHANE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
100-41-4	ETHYLBENZENE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/l	0.45 U	0.45 U	0.45 U	0.45 U
79-20-9	METHYL ACETATE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
78-93-3	METHYL ETHYL KETONE	ug/l	1.3 U	1.3 U	1.3 U	1.3 U
108-10-1	METHYL ISOBUTYL KETONE	ug/l	2.1 U	2.1 U	2.1 U	2.1 U
108-87-2	METHYLCYCLOHEXANE	ug/l	0.2 U	0.2 U	0.2 U	0.2 U
75-09-2	METHYLENE CHLORIDE	ug/l	0.41 U	0.41 U	0.41 U	0.41 U
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/l	0.43 U	0.43 U	0.43 U	0.43 U
100-42-5	STYRENE	ug/l	0.36 U	0.36 U	0.36 U	0.36 U
1634-04-4	TERT-BUTYL METHYL ETHER	ug/l	0.35 U	6.8	0.35 U	0.56 J
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/l	0.27 U	0.27 U	0.27 U	0.27 U
108-88-3	TOLUENE	ug/l	0.37 U	0.37 U	0.37 U	0.37 U
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/l	0.41 U	0.41 U	0.41 U	0.41 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/l	0.29 U	0.29 U	0.29 U	0.29 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/l	0.28 U	0.28 U	0.28 U	0.28 U
75-69-4	TRICHLOROFLUOROMETHANE	ug/l	0.35 U	0.35 U	0.35 U	0.35 U
75-01-4	VINYL CHLORIDE	ug/l	0.34 U	0.34 U	0.34 U	0.34 U

Con Ed - Halleck Street		Location ID:	MW-1	MW-2	MW-3	MW-3
Validated Groundwater Analytical Data		Sample ID:	MW-1-20131127	MW-2-20131127	MW-3-20131127	MW-3A-20131127
SDG: E4638		Lab Sample Id:	E4638-06	E4638-05	E4638-01	E4638-04
		Source:	CTECH	CTECH	CTECH	CTECH
		SDG:	E4638	E4638	E4638	E4638
		Matrix:	WATER	WATER	WATER	WATER
		Sampled:	11/27/2013 12:20	11/27/2013 10:50	11/27/2013 8:45	11/27/2013 9:10
		Validated:	2/4/2014	2/4/2014	2/4/2014	2/4/2014
CAS NO.	COMPOUND	UNITS:				
		SEMIVOLATILES				
95-94-3	1,2,4,5-TETRACHLOROBENZENE	ug/l	0.21 U	0.22 U	0.24 U	0.2 U
58-90-2	2,3,4,6-TETRACHLOROPHENOL	ug/l	0.21 U	0.22 U	0.24 U	0.2 U
95-95-4	2,4,5-TRICHLOROPHENOL	ug/l	0.42 U	0.43 U	0.47 U	0.41 U
88-06-2	2,4,6-TRICHLOROPHENOL	ug/l	0.58 U	0.6 U	0.66 U	0.57 U
120-83-2	2,4-DICHLOROPHENOL	ug/l	0.69 U	0.71 U	0.78 U	0.67 U
105-67-9	2,4-DIMETHYLPHENOL	ug/l	0.74 U	0.76 U	0.84 U	0.72 U
51-28-5	2,4-DINITROPHENOL	ug/l	2.2 UJ	2.3 UJ	2.5 UJ	2.1 UJ
121-14-2	2,4-DINITROTOLUENE	ug/l	1 U	1.1 U	1.2 U	1 U
606-20-2	2,6-DINITROTOLUENE	ug/l	0.33 UJ	0.34 UJ	0.38 UJ	0.33 UJ
91-58-7	2-CHLORONAPHTHALENE	ug/l	0.17 U	0.17 U	0.19 U	0.16 U
95-57-8	2-CHLOROPHENOL	ug/l	0.56 U	0.58 U	0.64 U	0.55 U
91-57-6	2-METHYLNAPHTHALENE	ug/l	0.33 U	0.34 U	0.38 U	0.33 U
95-48-7	2-METHYLPHENOL (O-CRESOL)	ug/l	0.25 U	0.26 U	0.28 U	0.24 U
88-74-4	2-NITROANILINE	ug/l	0.51 UJ	0.53 UJ	0.58 UJ	0.5 UJ
88-75-5	2-NITROPHENOL	ug/l	0.54 U	0.56 U	0.61 U	0.53 U
MEPH3MEPH	3- AND 4- METHYLPHENOL (TOTAL)	ug/l	0.4 U	0.41 U	0.45 U	0.39 U
91-94-1	3,3'-DICHLOROBENZIDINE	ug/l	1 U	1.1 U	1.2 U	1 U
99-09-2	3-NITROANILINE	ug/l	1 U	1.1 U	1.2 U	1 U
534-52-1	4,6-DINITRO-2-METHYLPHENOL	ug/l	0.77 U	0.8 U	0.87 U	0.76 U
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/l	0.24 U	0.25 U	0.27 U	0.23 U
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/l	0.42 U	0.43 U	0.47 U	0.41 U
106-47-8	4-CHLOROANILINE	ug/l	1 U	1.1 U	1.2 U	1 U
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/l	0.22 U	0.23 U	0.25 U	0.21 U
100-01-6	4-NITROANILINE	ug/l	1.4 UJ	1.5 UJ	1.6 UJ	1.4 UJ
100-02-7	4-NITROPHENOL	ug/l	2.1 U	2.2 U	2.4 U	2 U
83-32-9	ACENAPHTHENE	ug/l	0.22 U	0.23 U	0.25 U	0.21 U
208-96-8	ACENAPHTHYLENE	ug/l	0.73 U	0.75 U	0.82 U	0.71 U
98-86-2	ACETOPHENONE	ug/l	0.15 U	0.15 U	0.16 U	0.14 U
120-12-7	ANTHRACENE	ug/l	0.17 U	0.17 U	0.19 U	0.16 U
1912-24-9	ATRAZINE	ug/l	0.42 U	0.43 U	0.47 U	0.41 U
100-52-7	BENZALDEHYDE	ug/l	0.8 U	0.83 U	0.91 U	0.79 U
56-55-3	BENZO(A)ANTHRACENE	ug/l	0.17 U	0.17 U	0.19 U	0.16 U
50-32-8	BENZO(A)PYRENE	ug/l	0.15 U	0.15 U	0.16 U	0.14 U
205-99-2	BENZO(B)FLUORANTHENE	ug/l	0.3 U	0.31 U	0.34 U	0.3 U
191-24-2	BENZO(G,H,I)PERYLENE	ug/l	0.3 U	0.31 U	0.34 U	0.3 U
207-08-9	BENZO(K)FLUORANTHENE	ug/l	0.19 U	0.19 U	0.21 U	0.18 U
85-68-7	BENZYL BUTYL PHTHALATE	ug/l	0.2 U	0.2 U	0.22 U	0.19 U
92-52-4	BIPHENYL (DIPHENYL)	ug/l	0.16 U	0.16 U	0.18 U	0.15 U
111-91-1	BIS(2-CHLOROETHoxy) METHANE	ug/l	0.57 U	0.59 U	0.65 U	0.56 U
111-44-4	BIS(2-CHLOROETHYL) ETHER	ug/l	0.57 U	0.59 U	0.65 U	0.56 U
108-60-1	BIS(2-CHLOROISOPROPYL) ETHER	ug/l	0.18 U	0.18 U	0.2 U	0.17 U
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/l	0.17 UJ	0.17 UJ	0.19 UJ	0.16 UJ
105-60-2	CAPROLACTAM	ug/l	1 U	1.1 U	1.2 U	1 U
86-74-8	CARBAZOLE	ug/l	0.23 U	0.24 U	0.26 U	0.22 U
218-01-9	CHRYSENE	ug/l	0.19 U	0.19 U	0.21 U	0.18 U
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	0.44 U	0.45 U	0.49 U	0.43 U
132-64-9	DIBENZOFURAN	ug/l	0.25 U	0.26 U	0.28 U	0.24 U
84-66-2	DIETHYL PHTHALATE	ug/l	0.4 U	0.41 U	0.45 U	0.39 U
131-11-3	DIMETHYL PHTHALATE	ug/l	0.23 U	0.24 U	0.26 U	8.6 J
84-74-2	DI-N-BUTYL PHTHALATE	ug/l	1 U	1.1 U	1.2 U	1 U
117-84-0	DI-N-OCTYLPHthalate	ug/l	0.53 UJ	0.55 UJ	0.6 UJ	0.52 UJ

Con Ed - Halleck Street Validated Groundwater Analytical Data SDG: E4638		Location ID: Sample ID: Lab Sample Id: Source: SDG: Matrix: Sampled: Validated:	MW-1 MW-1-20131127 E4638-06 CTECH E4638 WATER 11/27/2013 12:20 2/4/2014	MW-2 MW-2-20131127 E4638-05 CTECH E4638 WATER 11/27/2013 10:50 2/4/2014	MW-3 MW-3-20131127 E4638-01 CTECH E4638 WATER 11/27/2013 8:45 2/4/2014	MW-3 MW-3A-20131127 E4638-04 CTECH E4638 WATER 11/27/2013 9:10 2/4/2014
CAS NO.	COMPOUND	UNITS:				
	<b>SEMOVOLATILES</b>					
206-44-0	FLUORANTHENE	ug/l	0.42 U	0.43 U	0.47 U	0.41 U
86-73-7	FLUORENE	ug/l	0.32 U	0.33 U	0.36 U	0.32 U
118-74-1	HEXACHLOROBENZENE	ug/l	0.19 U	0.19 U	0.21 U	0.18 U
87-68-3	HEXACHLOROBUTADIENE	ug/l	0.26 U	0.27 U	0.29 U	0.26 U
77-47-4	HEXACHLOROCYCLOPENTADIENE	ug/l	0.25 UJ	0.26 UJ	0.28 UJ	0.24 UJ
67-72-1	HEXACHLOROETHANE	ug/l	0.26 U	0.27 U	0.29 U	0.26 U
193-39-5	INDENO(1,2,3-C,D)PYRENE	ug/l	0.16 U	0.16 U	0.18 U	0.15 U
78-59-1	ISOPHORONE	ug/l	0.31 U	0.32 U	0.35 U	0.31 U
91-20-3	NAPHTHALENE	ug/l	0.13 U	0.13 U	0.14 U	0.12 U
98-95-3	NITROBENZENE	ug/l	0.71 U	0.73 U	0.8 U	0.69 U
621-64-7	N-NITROSODI-N-PROPYLAMINE	ug/l	0.21 U	0.22 U	0.24 U	0.2 U
86-30-6	N-NITROSODIPHENYLAMINE	ug/l	0.63 U	0.65 U	0.71 U	0.61 U
87-86-5	PENTACHLOROPHENOL	ug/l	1 U	1.1 U	1.2 U	1 U
85-01-8	PHENANTHRENE	ug/l	0.27 U	0.28 U	0.31 U	0.27 U
108-95-2	PHENOL	ug/l	0.22 U	0.23 U	0.25 U	0.21 U
129-00-0	PYRENE	ug/l	0.21 U	0.22 U	0.24 U	0.2 U
	<b>INORGANICS</b>					
7429-90-5	ALUMINUM	ug/l	920	252	48.2 J	53.6
7440-36-0	ANTIMONY	ug/l	8 U	8 U	8 U	8 U
7440-38-2	ARSENIC	ug/l	4.2 U	8.81 J	4.2 U	4.2 U
7440-39-3	BARIUM	ug/l	319	84.3	1400	1400
7440-41-7	BERYLLIUM	ug/l	0.7 U	0.7 U	0.7 U	0.7 U
7440-43-9	CADMIUM	ug/l	0.833 J	0.5 U	0.5 U	0.5 U
7440-70-2	CALCIUM	ug/l	148200	142000	175900	170000
7440-47-3	CHROMIUM, TOTAL	ug/l	34.8	1.97 J	1.1 UJ	5.67 J
7440-48-4	COBALT	ug/l	5.8 U	5.8 U	5.8 U	5.8 U
7440-50-8	COPPER	ug/l	2 U	2 U	2 U	2 U
7439-89-6	IRON	ug/l	50500	28100	33500	31300
7439-92-1	LEAD	ug/l	3.31 J	25.3	6.64	6.74
7439-95-4	MAGNESIUM	ug/l	34600	18400	58900	59400
7439-96-5	MANGANESE	ug/l	2900	472	719	698
7439-97-6	MERCURY	ug/l	0.1 U	0.1 U	0.1 U	0.1 U
7440-02-0	NICKEL	ug/l	13.4 J	4.2 U	4.2 U	4.2 U
7440-09-7	POTASSIUM	ug/l	24100	18500	41000	40000
7782-49-2	SELENIUM	ug/l	5.78 J	4.8 U	4.8 U	4.8 U
7440-22-4	SILVER	ug/l	1.5 U	1.5 U	1.5 U	1.5 U
7440-23-5	SODIUM	ug/l	610600	186000	635100	628900
7440-28-0	THALLIUM	ug/l	3.56 J	2.4 U	2.4 U	2.4 U
7440-62-2	VANADIUM	ug/l	6.1 U	6.1 U	6.1 U	6.1 U
7440-66-6	ZINC	ug/l	11.2 J	12.3 J	6.5 U	6.5 U
57-12-5	CYANIDE	ug/l	12	28	687	681

Con Ed - Halleck Street		Location ID:	FIELDQC
Validated Groundwater Analytical Data		Sample ID:	TB112713-20131127
SDG: E4638		Lab Sample Id:	E4638-08
		Source:	CTECH
		SDG:	E4638
		Matrix:	WATER
		Sampled:	11/27/2013 8:00
		Validated:	2/4/2014
CAS NO.	COMPOUND	UNITS:	
	VOLATILES		
71-55-6	1,1,1-TRICHLOROETHANE	ug/l	0.4 U
79-34-5	1,1,2-TETRACHLOROETHANE	ug/l	0.31 U
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	ug/l	0.45 U
79-00-5	1,1,2-TRICHLOROETHANE	ug/l	0.38 U
75-34-3	1,1-DICHLOROETHANE	ug/l	0.36 U
75-35-4	1,1-DICHLOROETHENE	ug/l	0.47 U
87-61-6	1,2,3-TRICHLOROBENZENE	ug/l	0.2 U
120-82-1	1,2,4-TRICHLOROBENZENE	ug/l	0.2 U
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	ug/l	0.46 U
106-93-4	1,2-DIBROMOETHANE	ug/l	0.41 U
95-50-1	1,2-DICHLOROBENZENE	ug/l	0.45 U
107-06-2	1,2-DICHLOROETHANE	ug/l	0.48 U
XYLMP	M,P-XYLENE (SUM OF ISOMERS)	ug/l	0.95 U
78-87-5	1,2-DICHLOROPROPANE	ug/l	0.46 U
541-73-1	1,3-DICHLOROBENZENE	ug/l	0.43 U
106-46-7	1,4-DICHLOROBENZENE	ug/l	0.32 U
123-91-1	1,4-DIOXANE (P-DIOXANE)	ug/l	50 U
591-78-6	2-HEXANONE	ug/l	1.9 U
67-64-1	ACETONE	ug/l	0.5 U
71-43-2	BENZENE	ug/l	0.32 U
74-97-5	BROMOCHLOROMETHANE	ug/l	0.2 U
75-27-4	BROMODICHLOROMETHANE	ug/l	0.36 U
75-25-2	BROMOFORM	ug/l	0.47 U
74-83-9	BROMOMETHANE	ug/l	0.2 U
75-15-0	CARBON DISULFIDE	ug/l	0.2 U
56-23-5	CARBON TETRACHLORIDE	ug/l	0.2 U
108-90-7	CHLOROBENZENE	ug/l	0.49 U
75-00-3	CHLOROETHANE	ug/l	0.2 U
67-66-3	CHLOROFORM	ug/l	0.34 U
74-87-3	CHLOROMETHANE	ug/l	0.2 U
156-59-2	CIS-1,2-DICHLOROETHYLENE	ug/l	0.35 U
10061-01-5	CIS-1,3-DICHLOROPROPENE	ug/l	0.31 U
110-82-7	CYCLOHEXANE	ug/l	0.2 UJ
124-48-1	DIBROMOCHLOROMETHANE	ug/l	0.2 U
75-71-8	DICHLORODIFLUOROMETHANE	ug/l	0.2 U
100-41-4	ETHYLBENZENE	ug/l	0.2 U
98-82-8	ISOPROPYLBENZENE (CUMENE)	ug/l	0.45 U
79-20-9	METHYL ACETATE	ug/l	0.2 U
78-93-3	METHYL ETHYL KETONE	ug/l	1.3 U
108-10-1	METHYL ISOBUTYL KETONE	ug/l	2.1 U
108-87-2	METHYLCYCLOHEXANE	ug/l	0.2 U
75-09-2	METHYLENE CHLORIDE	ug/l	0.41 U
95-47-6	O-XYLENE (1,2-DIMETHYLBENZENE)	ug/l	0.43 U
100-42-5	STYRENE	ug/l	0.36 U
1634-04-4	TERT-BUTYL METHYL ETHER	ug/l	0.35 U
127-18-4	TETRACHLOROETHYLENE(PCE)	ug/l	0.27 U
108-88-3	TOLUENE	ug/l	0.37 U
156-60-5	TRANS-1,2-DICHLOROETHENE	ug/l	0.41 U
10061-02-6	TRANS-1,3-DICHLOROPROPENE	ug/l	0.29 U
79-01-6	TRICHLOROETHYLENE (TCE)	ug/l	0.28 U
75-69-4	TRICHLOROFUOROMETHANE	ug/l	0.35 U
75-01-4	VINYL CHLORIDE	ug/l	0.34 U

Con Ed - Halleck Street		Location ID:	FIELDQC
Validated Groundwater Analytical Data		Sample ID:	TB112713-20131127
SDG: E4638		Lab Sample Id:	E4638-08
		Source:	CTECH
		SDG:	E4638
		Matrix:	WATER
		Sampled:	11/27/2013 8:00
		Validated:	2/4/2014
CAS NO.	COMPOUND	UNITS:	
	SEMIVOLATILES		
95-94-3	1,2,4,5-TETRACHLOROBENZENE	ug/l	
58-90-2	2,3,4,6-TETRACHLOROPHENOL	ug/l	
95-95-4	2,4,5-TRICHLOROPHENOL	ug/l	
88-06-2	2,4,6-TRICHLOROPHENOL	ug/l	
120-83-2	2,4-DICHLOROPHENOL	ug/l	
105-67-9	2,4-DIMETHYLPHENOL	ug/l	
51-28-5	2,4-DINITROPHENOL	ug/l	
121-14-2	2,4-DINITROTOLUENE	ug/l	
606-20-2	2,6-DINITROTOLUENE	ug/l	
91-58-7	2-CHLORONAPHTHALENE	ug/l	
95-57-8	2-CHLOROPHENOL	ug/l	
91-57-6	2-METHYLNAPHTHALENE	ug/l	
95-48-7	2-METHYLPHENOL (O-CRESOL)	ug/l	
88-74-4	2-NITROANILINE	ug/l	
88-75-5	2-NITROPHENOL	ug/l	
MEPH3MEPH	3- AND 4- METHYLPHENOL (TOTAL)	ug/l	
91-94-1	3,3'-DICHLOROBENZIDINE	ug/l	
99-09-2	3-NITROANILINE	ug/l	
534-52-1	4,6-DINITRO-2-METHYLPHENOL	ug/l	
101-55-3	4-BROMOPHENYL PHENYL ETHER	ug/l	
59-50-7	4-CHLORO-3-METHYLPHENOL	ug/l	
106-47-8	4-CHLOROANILINE	ug/l	
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	ug/l	
100-01-6	4-NITROANILINE	ug/l	
100-02-7	4-NITROPHENOL	ug/l	
83-32-9	ACENAPHTHENE	ug/l	
208-96-8	ACENAPHTHYLENE	ug/l	
98-86-2	ACETOPHENONE	ug/l	
120-12-7	ANTHRACENE	ug/l	
1912-24-9	ATRAZINE	ug/l	
100-52-7	BENZALDEHYDE	ug/l	
56-55-3	BENZO(A)ANTHRACENE	ug/l	
50-32-8	BENZO(A)PYRENE	ug/l	
205-99-2	BENZO(B)FLUORANTHENE	ug/l	
191-24-2	BENZO(G,H,I)PERYLENE	ug/l	
207-08-9	BENZO(K)FLUORANTHENE	ug/l	
85-68-7	BENZYL BUTYL PHTHALATE	ug/l	
92-52-4	BIPHENYL (DIPHENYL)	ug/l	
111-91-1	BIS(2-CHLOROETHOXY) METHANE	ug/l	
111-44-4	BIS(2-CHLOROETHYL) ETHER	ug/l	
108-60-1	BIS(2-CHLOROISOPROPYL) ETHER	ug/l	
117-81-7	BIS(2-ETHYLHEXYL) PHTHALATE	ug/l	
105-60-2	CAPROLACTAM	ug/l	
86-74-8	CARBAZOLE	ug/l	
218-01-9	CHRYSENE	ug/l	
53-70-3	DIBENZ(A,H)ANTHRACENE	ug/l	
132-64-9	DIBENZOFURAN	ug/l	
84-66-2	DIETHYL PHTHALATE	ug/l	
131-11-3	DIMETHYL PHTHALATE	ug/l	
84-74-2	DI-N-BUTYL PHTHALATE	ug/l	
117-84-0	DI-N-OCTYLPHTHALATE	ug/l	

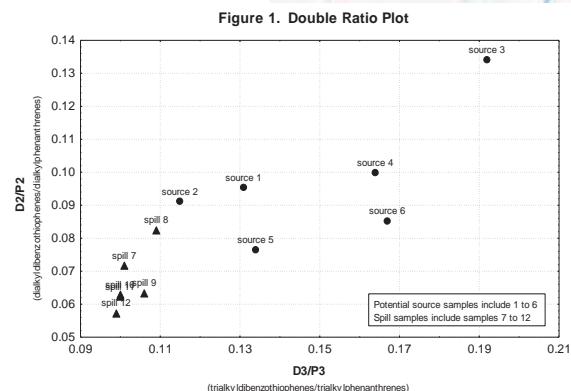
Con Ed - Halleck Street Validated Groundwater Analytical Data SDG: E4638	Location ID: Sample ID: Lab Sample Id: Source: SDG: Matrix: Sampled: Validated:	FIELDQC TB112713-20131127 E4638-08 CTECH E4638 WATER 11/27/2013 8:00 2/4/2014
CAS NO.	COMPOUND	UNITS:
	SEMIVOLATILES	
206-44-0	FLUORANTHENE	ug/l
86-73-7	FLUORENE	ug/l
118-74-1	HEXACHLOROBENZENE	ug/l
87-68-3	HEXACHLOROBUTADIENE	ug/l
77-47-4	HEXACHLOROCYCLOPENTADIENE	ug/l
67-72-1	HEXACHLOROETHANE	ug/l
193-39-5	INDENO(1,2,3-C,D)PYRENE	ug/l
78-59-1	ISOPHORONE	ug/l
91-20-3	NAPHTHALENE	ug/l
98-95-3	NITROBENZENE	ug/l
621-64-7	N-NITROSODI-N-PROPYLAMINE	ug/l
86-30-6	N-NITROSODIPHENYLAMINE	ug/l
87-86-5	PENTACHLOROPHENOL	ug/l
85-01-8	PHENANTHRENE	ug/l
108-95-2	PHENOL	ug/l
129-00-0	PYRENE	ug/l
	INORGANICS	
7429-90-5	ALUMINUM	ug/l
7440-36-0	ANTIMONY	ug/l
7440-38-2	ARSENIC	ug/l
7440-39-3	BARIUM	ug/l
7440-41-7	BERYLLIUM	ug/l
7440-43-9	CADMIUM	ug/l
7440-70-2	CALCIUM	ug/l
7440-47-3	CHROMIUM, TOTAL	ug/l
7440-48-4	COBALT	ug/l
7440-50-8	COPPER	ug/l
7439-89-6	IRON	ug/l
7439-92-1	LEAD	ug/l
7439-95-4	MAGNESIUM	ug/l
7439-96-5	MANGANESE	ug/l
7439-97-6	MERCURY	ug/l
7440-02-0	NICKEL	ug/l
7440-09-7	POTASSIUM	ug/l
7782-49-2	SELENIUM	ug/l
7440-22-4	SILVER	ug/l
7440-23-5	SODIUM	ug/l
7440-28-0	THALLIUM	ug/l
7440-62-2	VANADIUM	ug/l
7440-66-6	ZINC	ug/l
57-12-5	CYANIDE	ug/l

## **APPENDIX D**

### **HYDROCARBON FINGERPRINT RESULTS**

# Environmental Forensic Report

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**Halleck Street**

SDG: PA131108 (MC26103)

*Report To:*

**Parsons  
200 Cottontail Lane  
Somerset, NJ**

*Report By:*

**META Environmental, Inc.  
115 Dean Avenue, Suite 300  
Box 13  
Franklin MA 02038**

**November 29, 2013**

***Identifying and allocating sources of pollutants in complex environments.***

---

## Certification

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This certifies that this package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed herein. The results included in this data report relate only to the samples as received and analyzed by the laboratory.

This report shall not be reproduced except in full, without the written approval of META Environmental, Inc.

Release of the data contained in this hardcopy or electronic copy data package has been authorized by the following signature(s).



David M. Mauro  
President, Senior Scientist

November 29, 2013

Date

META Environmental, Inc.  
115 Dean Avenue, Suite 300  
Franklin MA 02038  
Phone: 508-541-9146  
E-Mail [dmauro@metaenv.com](mailto:dmauro@metaenv.com)

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## Sample Delivery Group

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Project: Halleck Street  
 Client Parsons  
 200 Cottontail Lane  
 Somerset, NJ Report  
 Contact: Eric Gaulin  
 Dates of Receipt: 11/8/2013  
 Sample Summary: The samples received for this project are summarized in the attached sample login forms in Appendix A.  
 META Project Number: P06022  
 SDG No.: PA131108 (MC26103)

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## Chain of Custody

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The sample was received by Accutest Laboratories of New England for META Environmental, Inc. in good condition. The internal temperature of the shipping container was within the recommended 0-6°C range and was as follows:

Samples received: 11/08/2013      0.5°C      Ice present

Internal chain of custody procedures were followed after sample receipt. Samples were stored in a locked refrigerator. A sample custody logbook contains the record of sample removal from the secure sample storage area to the sample preparation laboratory. The custody record for the sample extracts is present on the sample extraction logbook page.

The disposal of samples and extracts will be authorized one month after the release of this data report. Sample disposal will be documented.

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## Methods

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The NAPL sample was prepared by solvent dilution (EPA 3580) using dichloromethane (DCM). The extracts were spiked with internal standard and analyzed by GC/FID (EPA 8015M) for fingerprinting and TPH and by GC/MS/SIM (EPA 8270M) for mono- and polycyclic aromatic hydrocarbons (MAHs and PAHs), alkyl PAH homologues and other selected compounds.

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## Results

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Sample results are presented in several appendices which follow this narrative.

- Appendix B: GC/FID Fingerprints
- Appendix C: Data Summary Sheets
- Appendix D: Extracted Ion Current Profiles (EICPs)

## Appendix E: Accutest Laboratory Report

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### Quality Control

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The quality control measures, criteria, and results are included in the Accutest laboratory report (Appendix E).

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### Interpretation

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#### Introduction

One sample of non-aqueous phase liquid (NAPL) (SB-4(15-17)NAPL) was received by Accutest Laboratories in Marlborough MA for META from the Halleck Street site on 11/8/2013. The sample was analyzed for hydrocarbon fingerprint and an expanded list of MAHs and PAHs.

This report summarizes the findings and compares the samples.

#### Sample-Specific Observations

##### SB-4(15-17)NAPL

Sample SB-4(15-17)NAPL contained pyrogenic material (see definitions). The pyrogenic material was indicated by the wide range distribution of unsubstituted mono- and polycyclic aromatic hydrocarbons (MAHs & PAHs), with the 2 and 3 ring PAHs most abundant. The abundance of naphthalene relative to the higher molecular weight PAHs suggested that the pyrogenic material in the sample had not experienced substantial weathering.

The ratio of fluoranthene to pyrene (Fl/Py – Table 1) as well as the double ratio plot of benzofluorenes/methylpyrenes (BF/MP) to Fl/Py (Figure 1) shows that sample SB-4(15-17)NAPL was very similar to tars in META's reference library that were formed from manufactured gas plants (MGPs) utilizing carbureted water gas processes.

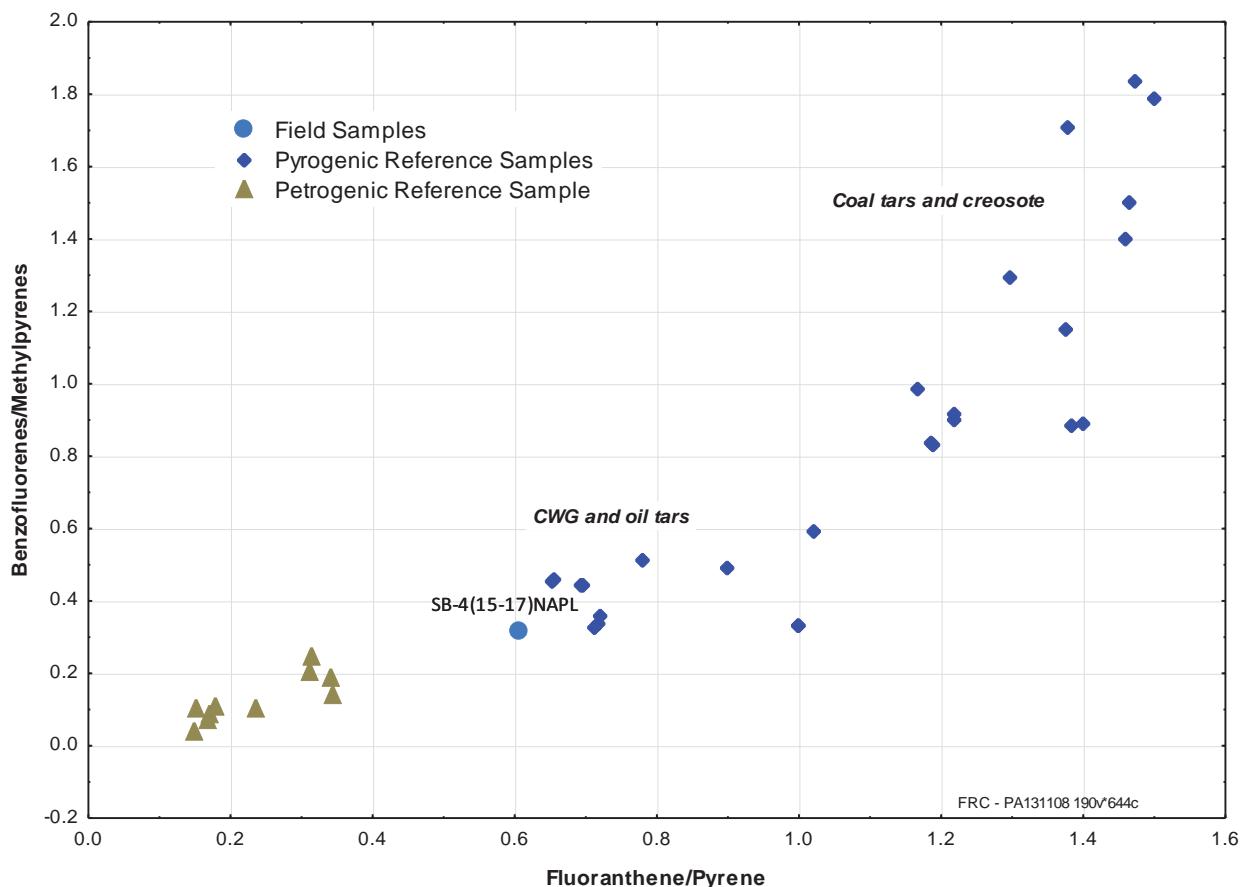
No other pyrogenic or petrogenic substances were detected in the sample.

**Table 1. Selected Source and Weathering Ratios**

	Fl/Py	D/F	C3D/C3PA	C2D/C2PA	BF/MP
SB-4(15-17)NAPL	0.606	0.194	0.702	0.475	0.318

#### Ratios:

Fl/Py	fluoranthene/pyrene
D/F	dibenzofuran/fluorene
C3D/C3PA	trikylbibenzothiophenes/trikylphenanthrenes/anthracenes
C2D/C2PA	dialkyldibenzothiophenes/dialkylphenanthrenes/anthracenes
BF/MP	benzofluorenes/methylpyrenes
NC	Not calculable

**Figure 1. Selected Diagnostic Ratios**


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## Definitions

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Pyrogenic substances are complex mixtures of primarily hydrocarbons produced from organic matter subjected to high temperatures but with insufficient oxygen for complete combustion. Pyrogenic materials are produced by fires, internal combustion engines, and furnaces. They also are formed when coke or gas are produced from coal or oil. Coal-tar based products, such as roofing, pavement sealers, waterproofing, pesticides, and some shampoos contain pyrogenic materials.

Petrogenic substances include crude oil and crude oil derivatives such as gasoline, heating oil, and asphalt.

Pitch is the semi-solid or solid material consisting of high molecular weight hydrocarbons that remain following coal tar distillation.

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## References

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“Chemical Fingerprinting of Hydrocarbons,” in: Introduction to Environmental Forensics. B.L. PA131108 MC26103 Halleck St Report  
November 29, 2013

Murphy and R.D. Morrison editors, Academic Press, San Diego, CA 2002.

Mauro, D.M., "Chemical Source Attribution at former MGP Sites," EPRI Report 1000728, December 2000.

## **Appendix A**

## **Chains of Custody**

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META Environmental, Inc.						
Sample Receipt Log						
Lab ID	Field ID	Matrix	Date Sampled	Date Received	Project #	Container
PA 3108-01	SB-1 (5.17)NAPL	soil	10/31/2013	11/8/2013	Proj022-60	MC26103

Logged By: \_\_\_\_\_

Date: \_\_\_\_\_

Reviewed By: \_\_\_\_\_

Date: \_\_\_\_\_

## **Appendix B**

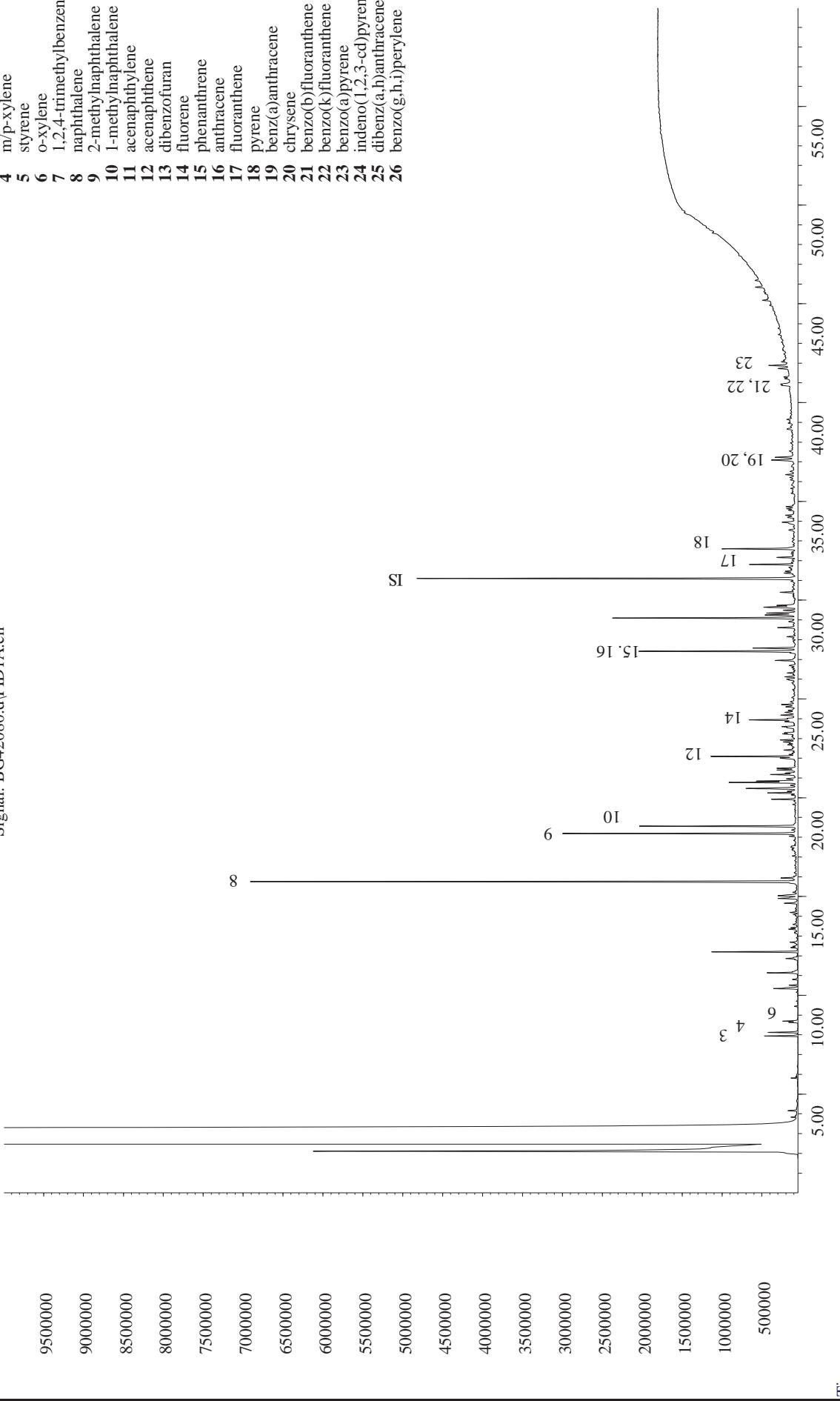
## **GC/FID Fingerprints**

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# GC/FID Fingerprint

10 of 133

Signal: BG42080.d\FID1A.ch



Analysis Date: 11/20/2013

IS -  $\delta$ -androstane  
SS2 - o-terphenyl

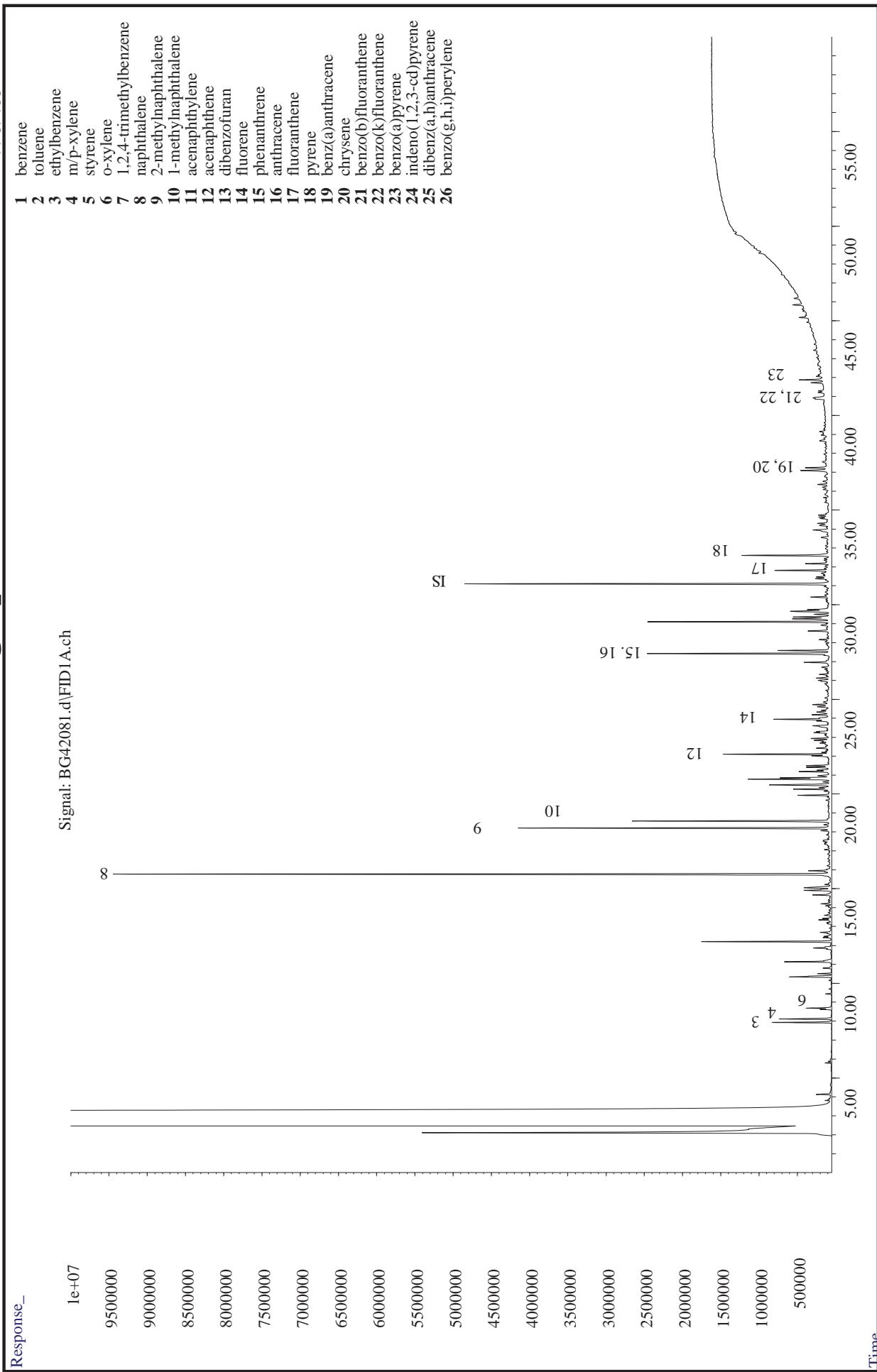
**Field ID:** SB-4(15-17)NAPL

Laboratory ID: MC26103-1

Method: EPA 8015M

# GC/FID Fingerprint

11 of 133



Analysis Date: 11/20/2013

IS – 5 $\alpha$ -androstane  
SS2 – o-terphenyl

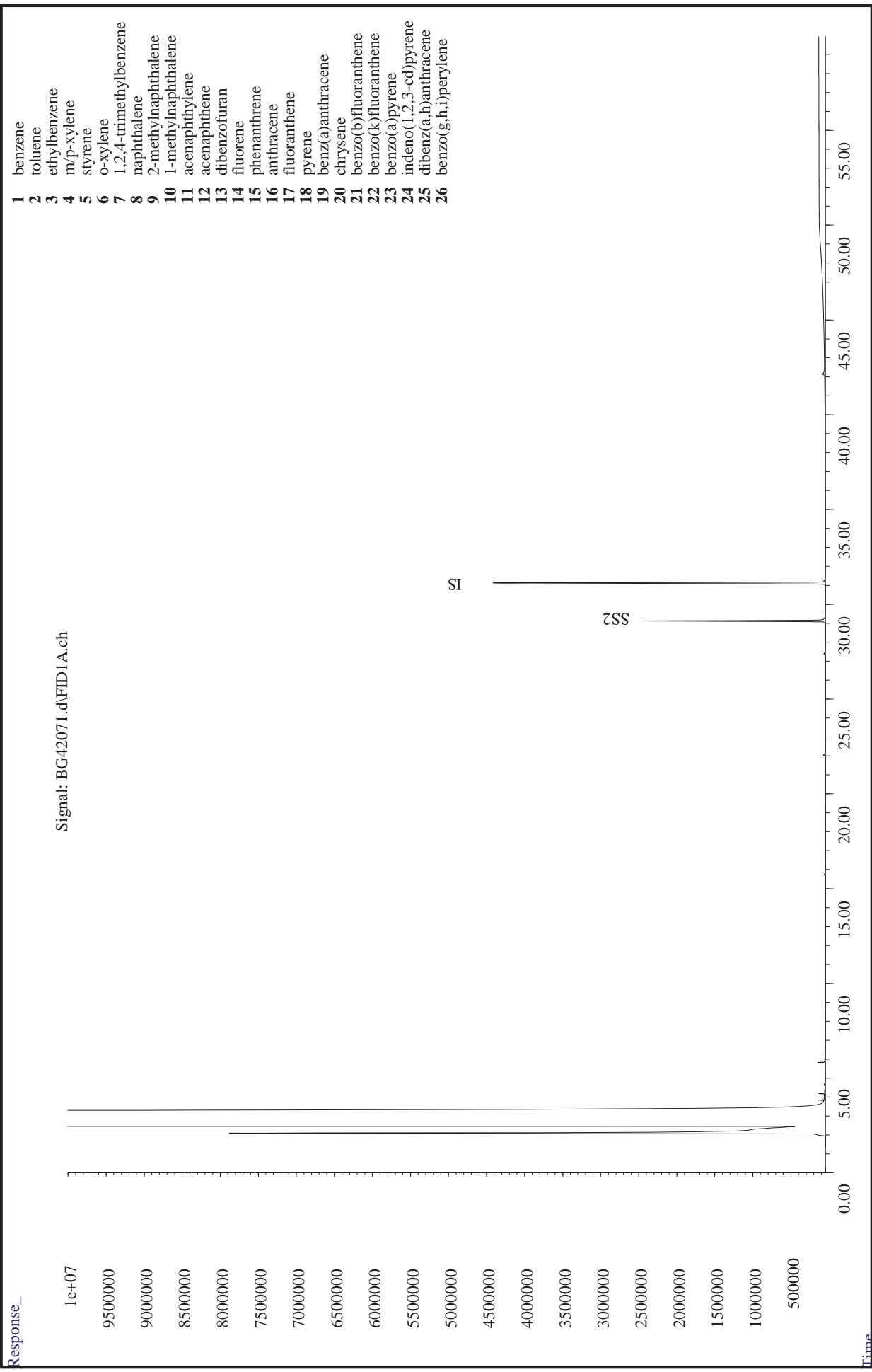
Field ID: SB-4(15-17)NAPL dup

Laboratory ID: OP35833-dup1

Method: EPA 8015M

# GC/FID Fingerprint

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**Analysis Date:** 11/20/2013

*IS* – 5 $\alpha$ -androstane  
*SS2* – *o*-terphenyl

**Field ID:** Method Blank  
**Laboratory ID:** op35833-MB  
**Method:** EPA 8015M

## **Appendix C**

## **Data Summary Sheets**

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Accutest Labs of New England, Inc. 11/29/2013  
 Job Number: MC26103  
 Account: META Environmental, Inc.  
 Project: Parsons, Halleck Street, NJ  
 Project Number:

Client Sample ID: SB-4(15-17)NAPL  
 Lab Sample ID: MC26103-1  
 Date Sampled: 10/31/2013  
 Matrix: Oil

GC/MS Semi-volatiles (D5739-06/8270C SIM)

Benzene	mg/kg	10900
C1-Benzene	mg/kg	1580
C2-Benzenes	mg/kg	72300
C3-Benzenes	mg/kg	82600
C4-Benzenes	mg/kg	43300
C5-Benzenes	mg/kg	12800
Toluene	mg/kg	2190
Ethylbenzene	mg/kg	55300
m,p-Xylene	mg/kg	52800
Styrene	mg/kg	11300
o-Xylene	mg/kg	23800
Isopropylbenzene	mg/kg	6660
n-Propylbenzene	mg/kg	2440
1,3,5-Trimethylbenzene	mg/kg	13200
1,2,3-Trimethylbenzene	mg/kg	13500
1,2,4-Trimethylbenzene	mg/kg	40900
t-Butylbenzene	mg/kg	5370
sec-Butylbenzene	mg/kg	ND
p-Isopropyltoluene	mg/kg	6740
n-Butylbenzene	mg/kg	1480
trans-Decalin	mg/kg	ND
cis-Decalin	mg/kg	ND
Benzo(b)thiophene	mg/kg	26600
Naphthalene	mg/kg	744000
2-Methylnaphthalene	mg/kg	310000
1-Methylnaphthalene	mg/kg	195000
C1-Naphthalenes	mg/kg	331000
C2-Naphthalenes	mg/kg	155000
C3-Naphthalenes	mg/kg	51800
C4-Naphthalenes	mg/kg	14800
Biphenyl	mg/kg	27300
Acenaphthylene	mg/kg	25200
Acenaphthene	mg/kg	102000
Dibenzofuran	mg/kg	9390
Fluorene	mg/kg	48300
C1-Fluorenes	mg/kg	23600
C2-Fluorenes	mg/kg	13900
C3-Fluorenes	mg/kg	6710
Dibenzothiophene	mg/kg	28500

Client Sample ID: **SB-4(15-17)NAPL**  
 Lab Sample ID: MC26103-1  
 Date Sampled: 10/31/2013  
 Matrix: Oil

**GC/MS Semi-volatiles (D5739-06/8270C SIM)**

C1-Dibenzothiophenes	mg/kg	26300
C2-Dibenzothiophenes	mg/kg	16100
C3-Dibenzothiophenes	mg/kg	6760
C4-Dibenzothiophenes	mg/kg	1670
Phenanthrene	mg/kg	179000
Anthracene	mg/kg	53200
C1-Phenanthrenes/Anthracenes	mg/kg	90900
C2-Phenanthrenes/Anthracenes	mg/kg	33900
C3-Phenanthrenes/Anthracenes	mg/kg	9630
C4-Phenanthrenes/Anthracenes	mg/kg	2610
Retene	mg/kg	1020
Benzo(b)naphtho(2,1-d)thiophene	mg/kg	4680
Fluoranthene	mg/kg	53200
Pyrene	mg/kg	87800
C1-Fluoranthenes/Pyrenes	mg/kg	45700
C2-Fluoranthenes/Pyrenes	mg/kg	11500
C3-Fluoranthenes/Pyrenes	mg/kg	2630
Benzo(b)fluorene	mg/kg	3540
Benzo(c)fluorene	mg/kg	2140
2-Methylpyrene	mg/kg	6480
4-Methylpyrene	mg/kg	5760
1-Methylpyrene	mg/kg	5610
Benzo(a)anthracene	mg/kg	22200
Chrysene	mg/kg	22000
C1-Benzo(a)anthracenes/Chrysenes	mg/kg	10300
C2-Benzo(a)anthracenes/Chrysenes	mg/kg	4160
C3-Benzo(a)anthracenes/Chrysenes	mg/kg	ND
C4-Benzo(a)anthracenes/Chrysenes	mg/kg	ND
Benzo(b)fluoranthene	mg/kg	8440
Benzo(k)fluoranthene	mg/kg	12200
Benzo(e)pyrene	mg/kg	11700
Benzo(a)pyrene	mg/kg	21800
Perylene	mg/kg	2850
Indeno(1,2,3-cd)pyrene	mg/kg	7000
Dibenzo(a,h)anthracene	mg/kg	2040
Benzo(g,h,i)perylene	mg/kg	9150
Coronene	mg/kg	2100

**GC Semi-volatiles (ASTM D3328-06)**

TPH (C8-C40)	mg/kg	2510000
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## **Appendix D**

# **Extracted Ion Current Profiles (EICPs)**

---

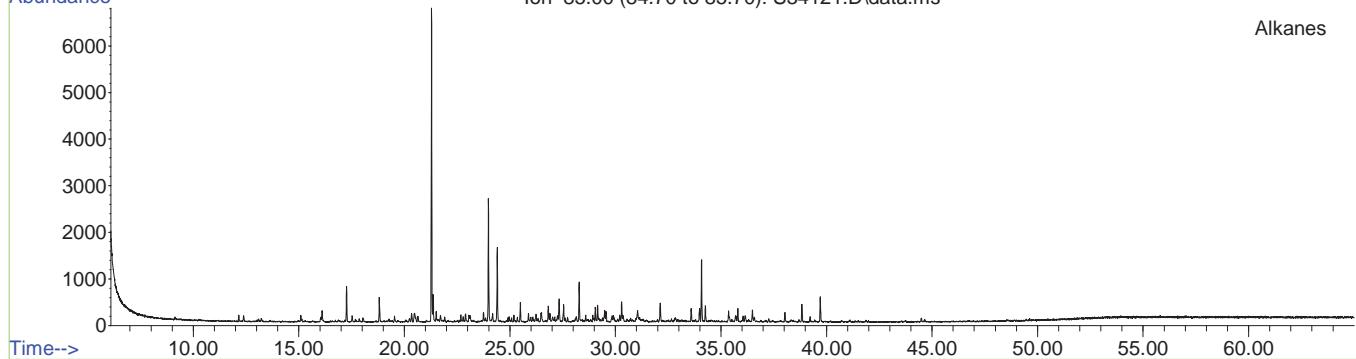
File: Z:\1\data\S131120\S34121.D

Date Acquired: 20 Nov 2013 7:05 pm

Sample Name: mc26103-1

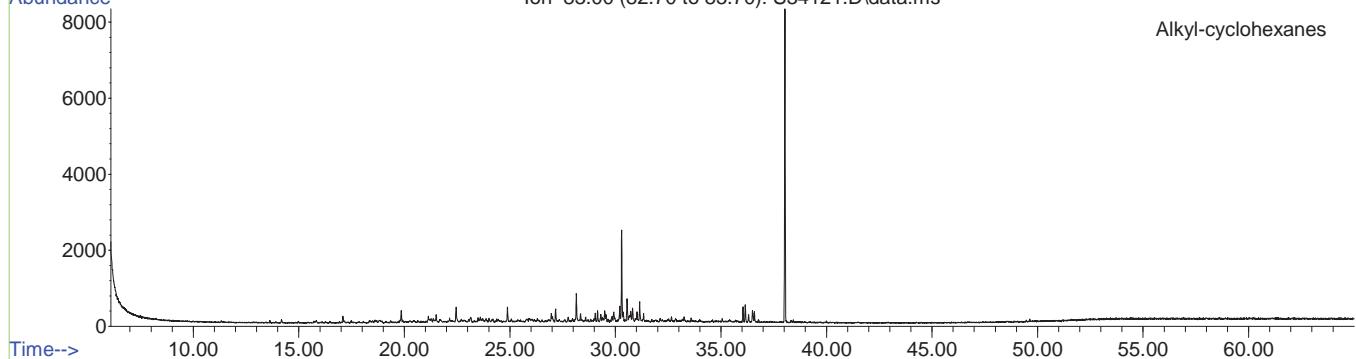
Misc Info: op35834,mss1523,0.0011,,,10,10

Abundance Ion 85.00 (84.70 to 85.70): S34121.D\data.ms



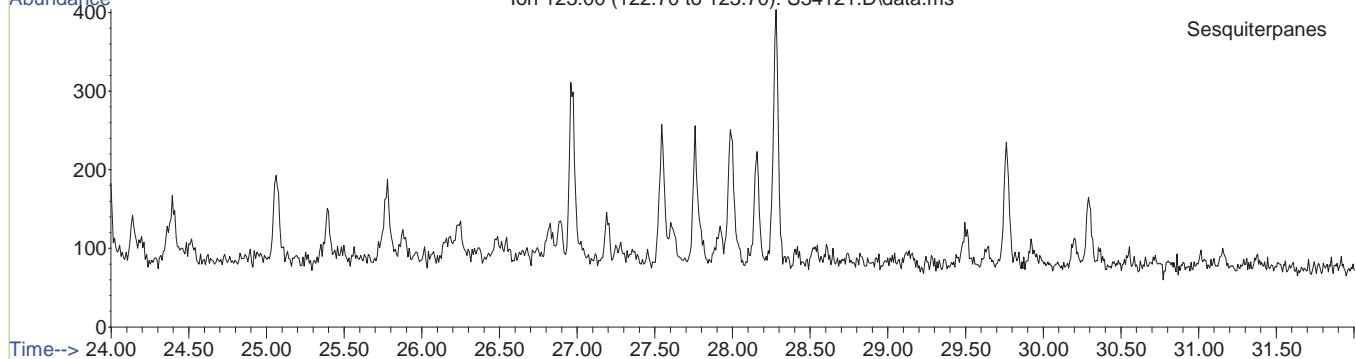
Alkanes

Abundance Ion 83.00 (82.70 to 83.70): S34121.D\data.ms



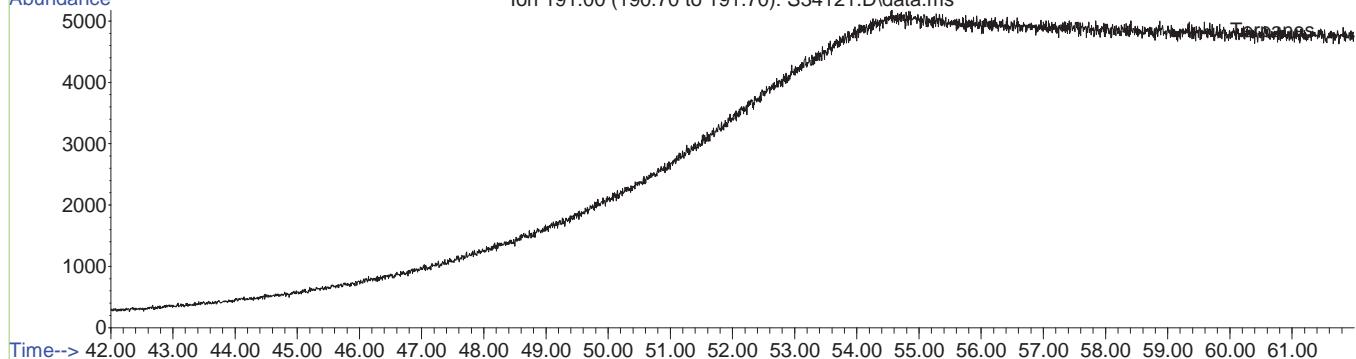
Alkyl-cyclohexanes

Abundance Ion 123.00 (122.70 to 123.70): S34121.D\data.ms



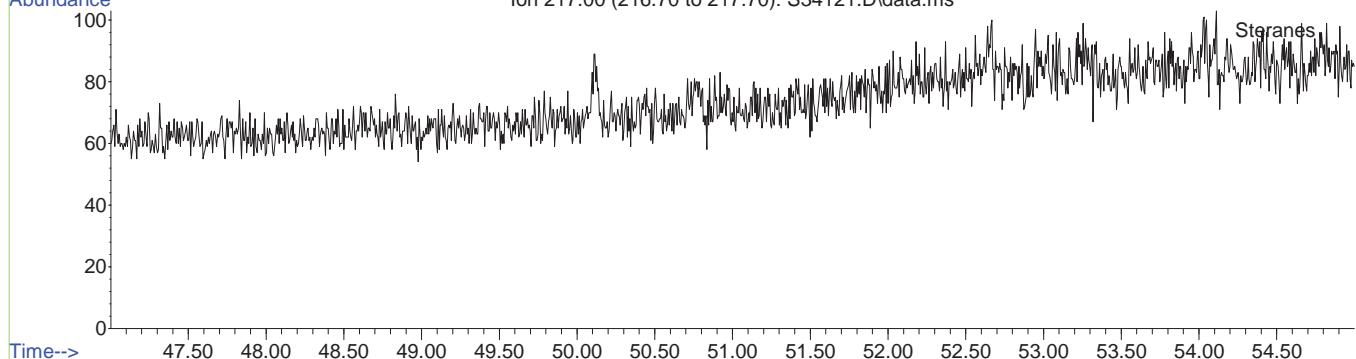
Sesquiterpanes

Abundance Ion 191.00 (190.70 to 191.70): S34121.D\data.ms



Terpanes

Abundance Ion 217.00 (216.70 to 217.70): S34121.D\data.ms



Steranes

Time-->

47.50 48.00 48.50 49.00 49.50 50.00 50.50 51.00 51.50 52.00 52.50 53.00 53.50 54.00 54.50

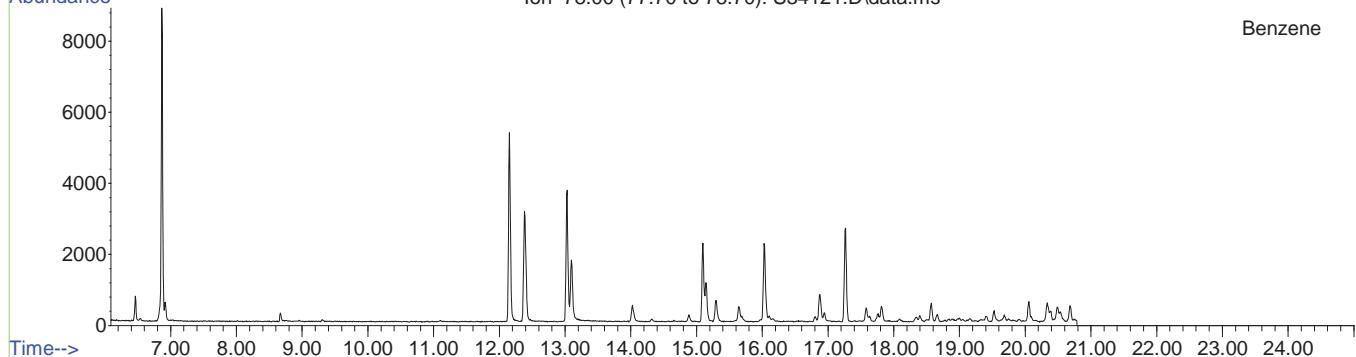
File: Z:\1\data\S131120\S34121.D

Date Acquired: 20 Nov 2013 7:05 pm

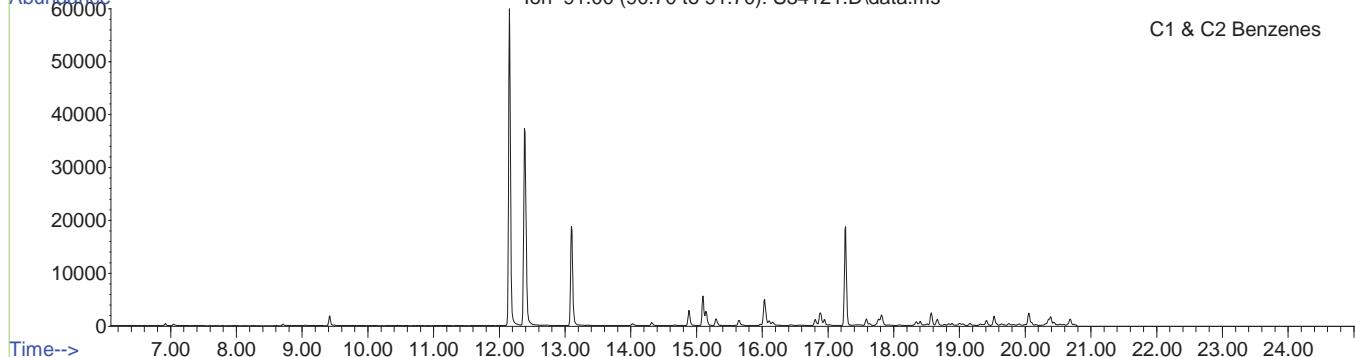
Sample Name: mc26103-1

Misc Info: op35834,mss1523,0.0011,,,10,10

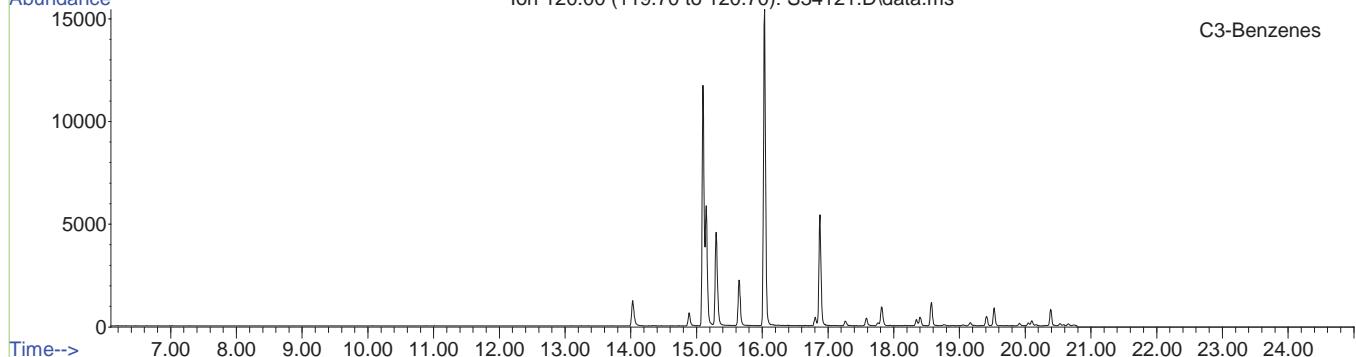
Abundance Ion 78.00 (77.70 to 78.70): S34121.D\data.ms



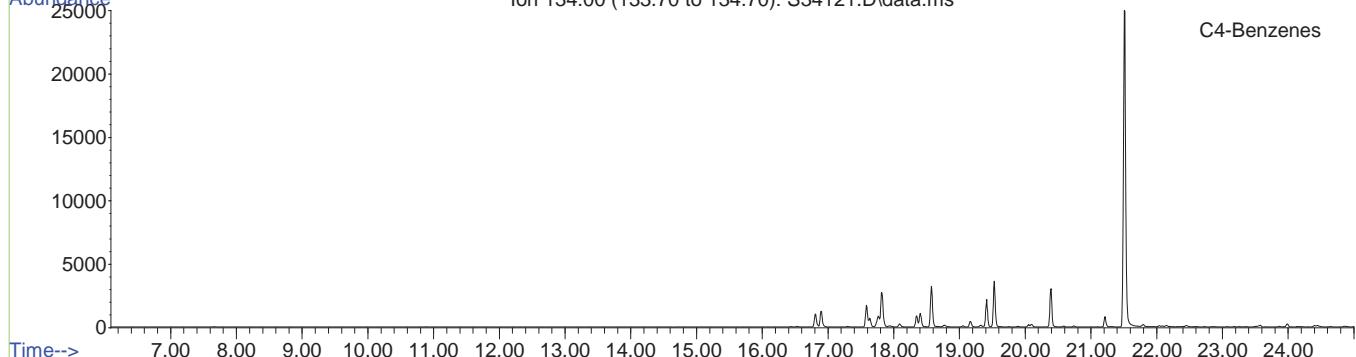
Time--> Abundance Ion 91.00 (90.70 to 91.70): S34121.D\data.ms



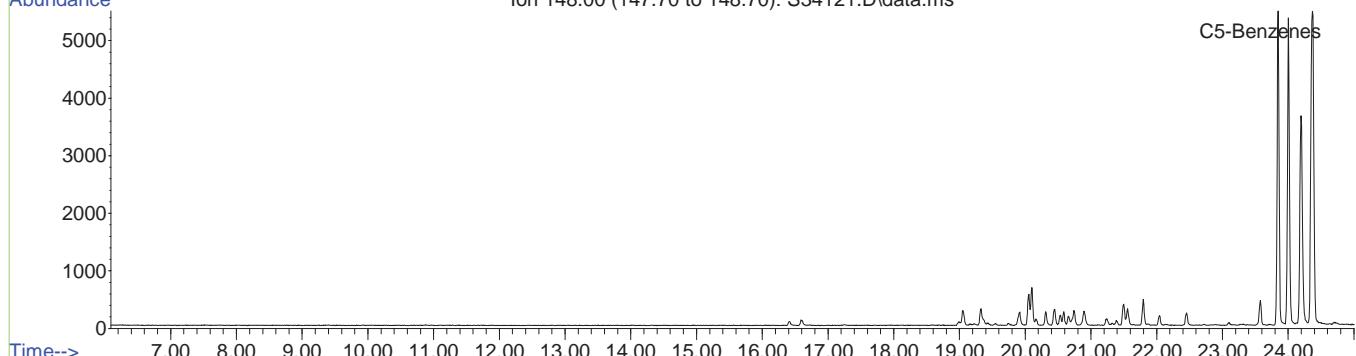
Time--> Abundance Ion 120.00 (119.70 to 120.70): S34121.D\data.ms



Time--> Abundance Ion 134.00 (133.70 to 134.70): S34121.D\data.ms



Time--> Abundance Ion 148.00 (147.70 to 148.70): S34121.D\data.ms



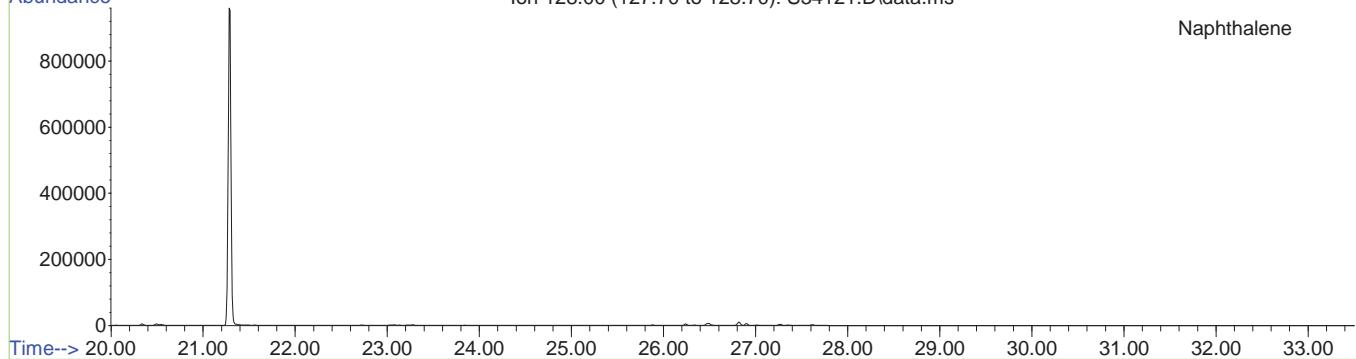
File: Z:\1\data\S131120\S34121.D

Date Acquired: 20 Nov 2013 7:05 pm

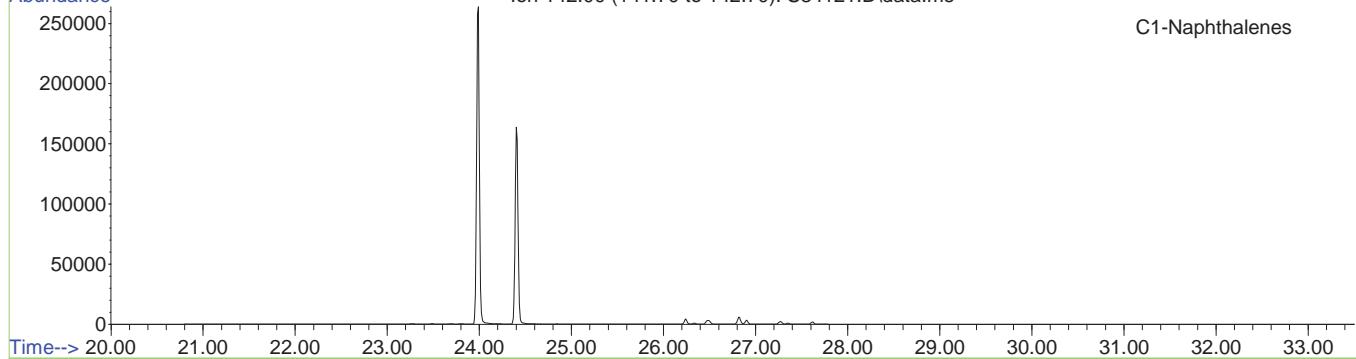
Sample Name: mc26103-1

Misc Info: op35834,mss1523,0.0011,,,10,10

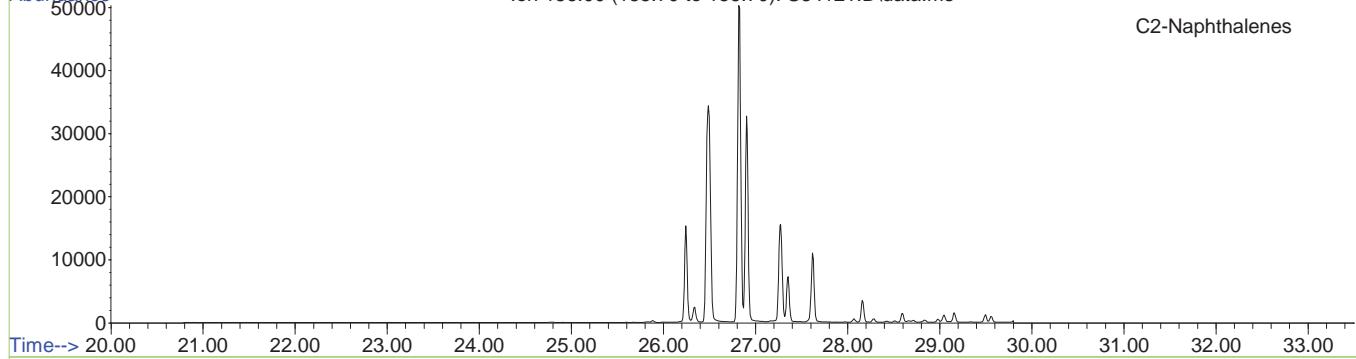
Abundance Ion 128.00 (127.70 to 128.70): S34121.D\data.ms



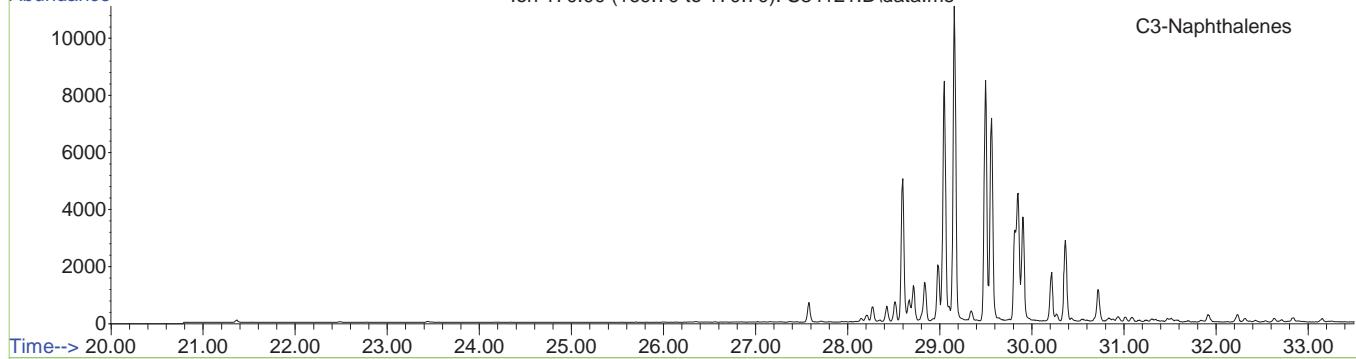
Abundance Ion 142.00 (141.70 to 142.70): S34121.D\data.ms



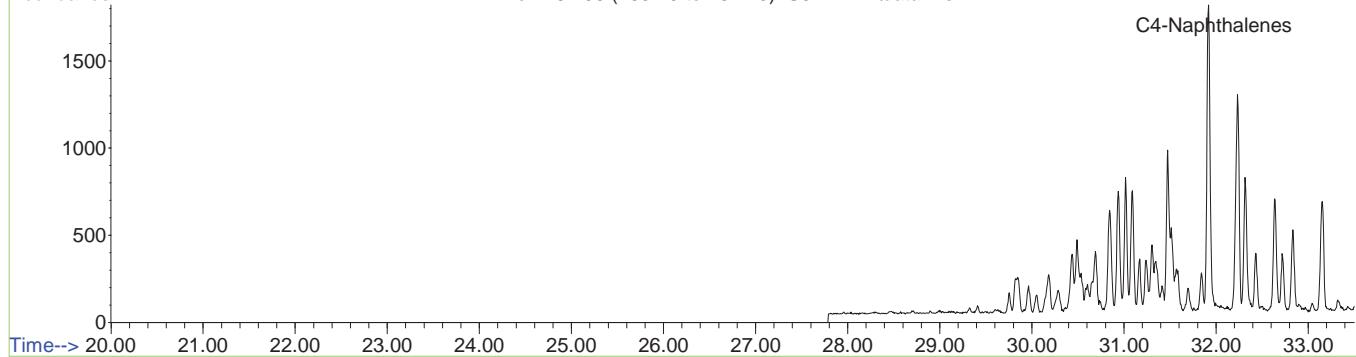
Abundance Ion 156.00 (155.70 to 156.70): S34121.D\data.ms



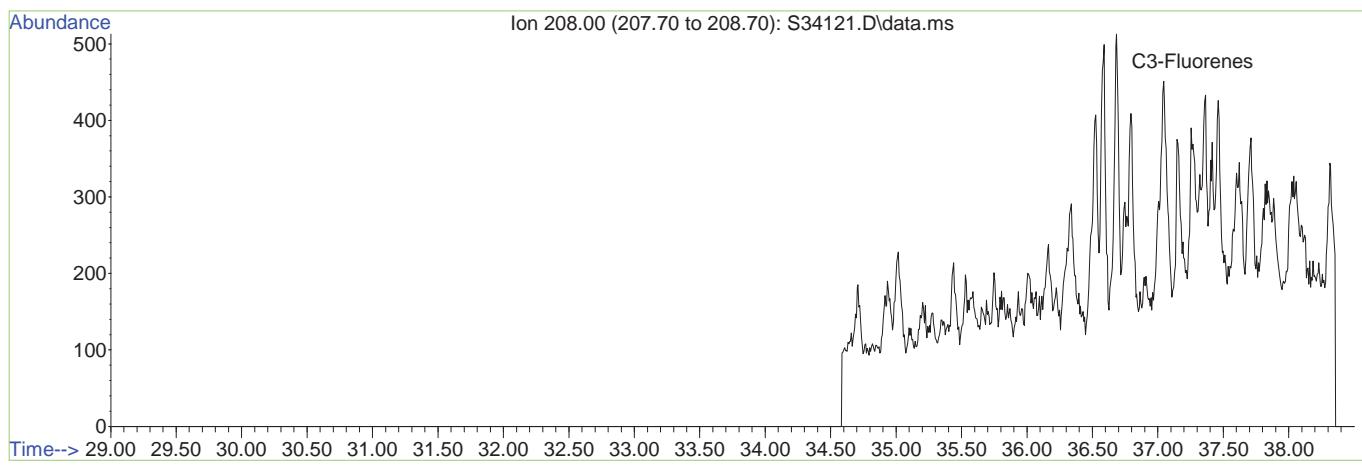
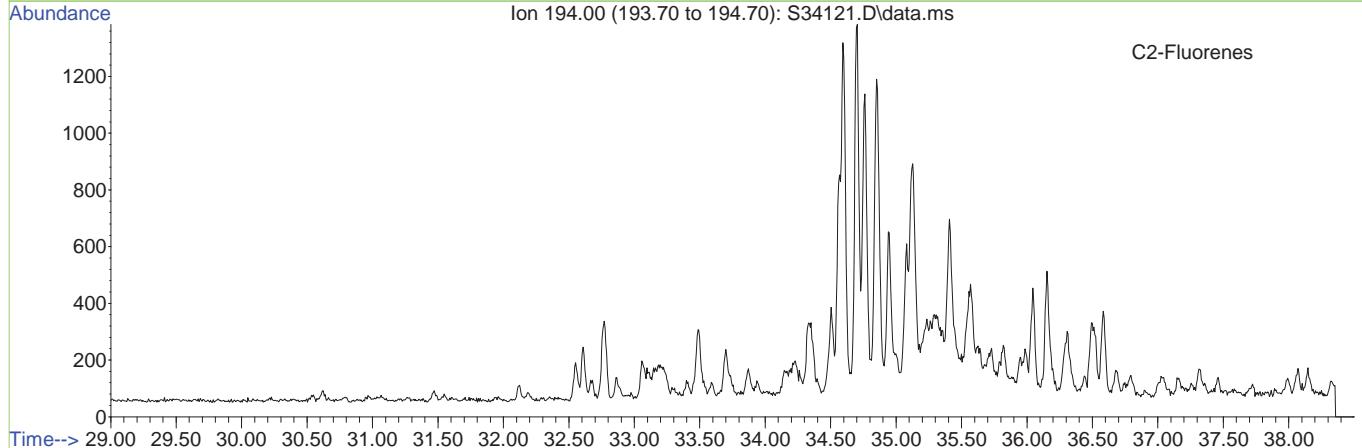
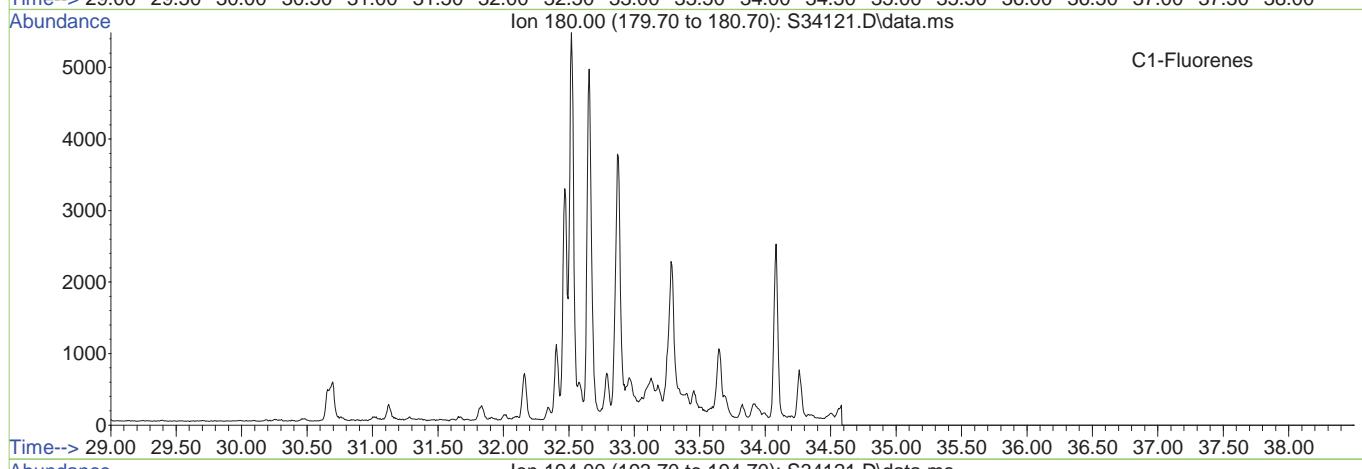
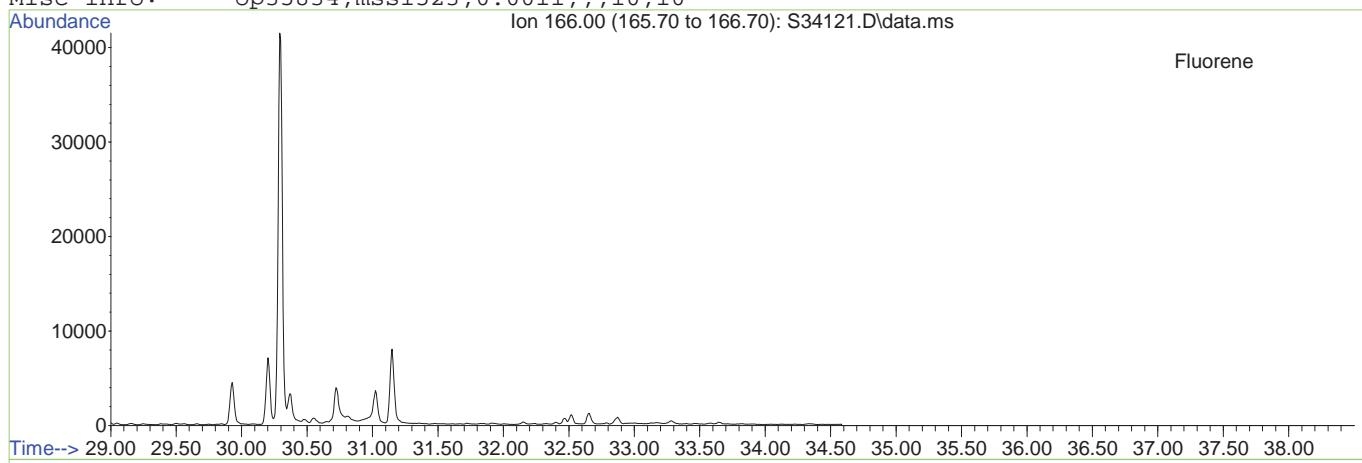
Abundance Ion 170.00 (169.70 to 170.70): S34121.D\data.ms



Abundance Ion 184.00 (183.70 to 184.70): S34121.D\data.ms



File: Z:\1\data\S131120\S34121.D  
Date Acquired: 20 Nov 2013 7:05 pm  
Method File: SAPAHSIM-MTBE.M  
Sample Name: mc26103-1  
Misc Info: op35834,mss1523,0.0011,,,10,10



File: Z:\1\data\S131120\S34121.D

Date Acquired: 20 Nov 2013 7:05 pm

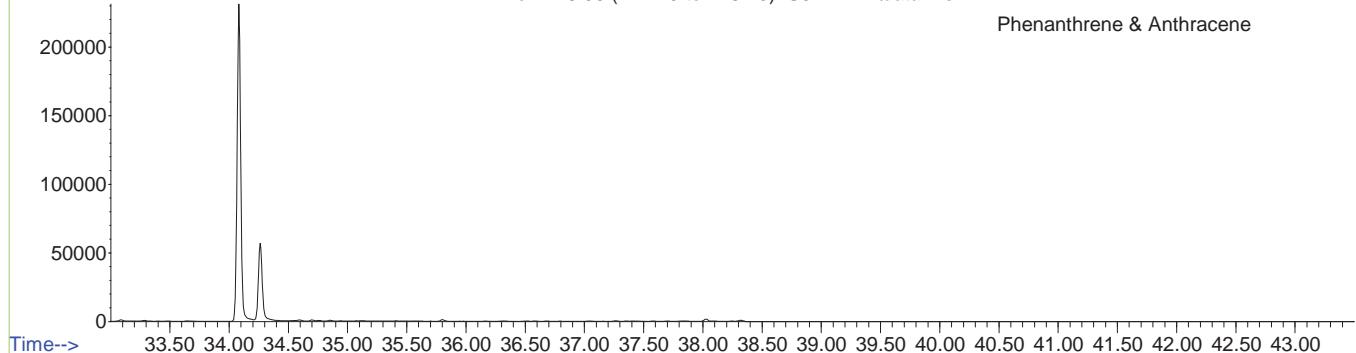
Sample Name: mc26103-1

Misc Info: op35834,mss1523,0.0011,,,10,10

Abundance

Ion 178.00 (177.70 to 178.70): S34121.D\data.ms

Phenanthrene &amp; Anthracene



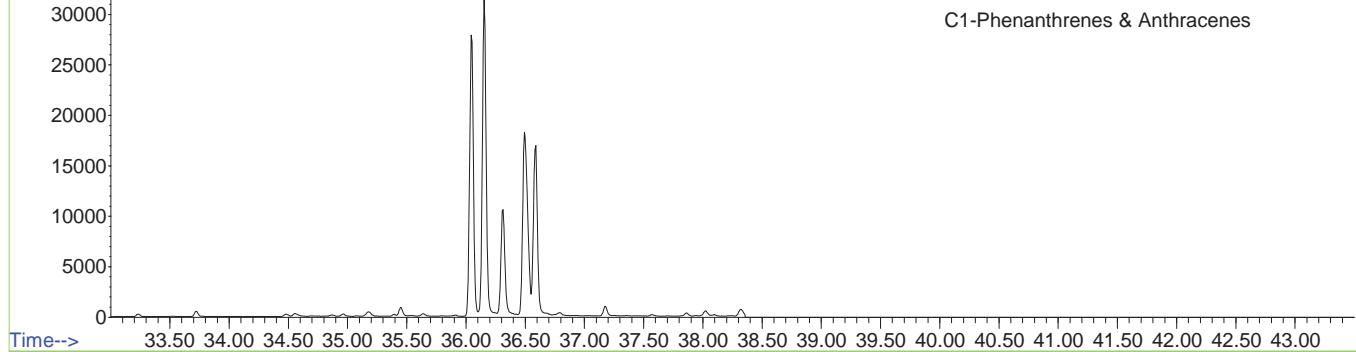
Time--&gt;

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00

Abundance

Ion 192.00 (191.70 to 192.70): S34121.D\data.ms

C1-Phenanthrenes &amp; Anthracenes



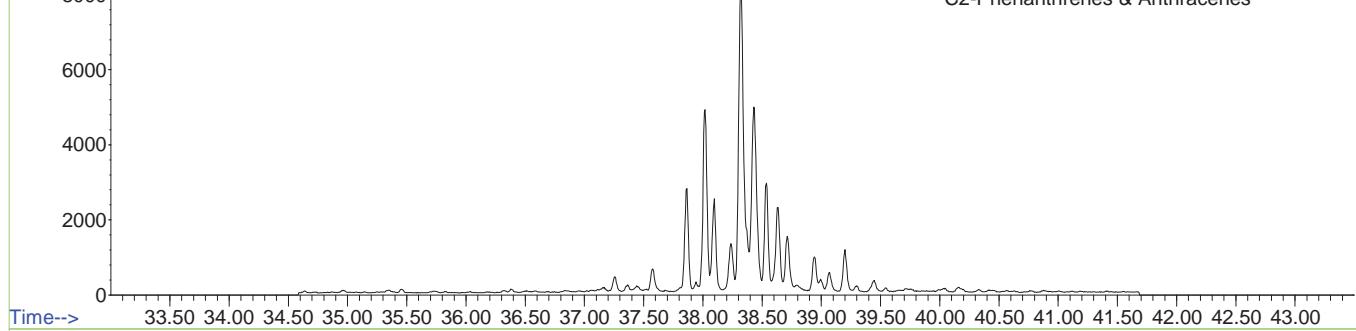
Time--&gt;

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00

Abundance

Ion 206.00 (205.70 to 206.70): S34121.D\data.ms

C2-Phenanthrenes &amp; Anthracenes



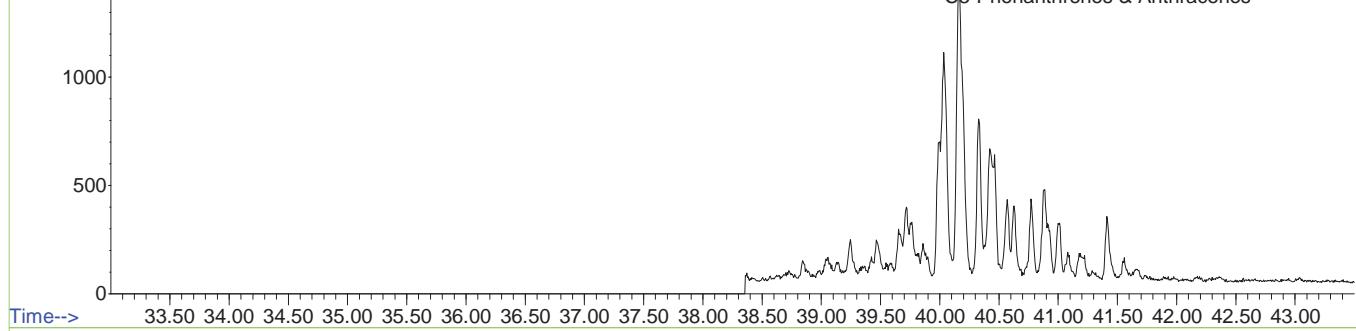
Time--&gt;

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00

Abundance

Ion 220.00 (219.70 to 220.70): S34121.D\data.ms

C3-Phenanthrenes &amp; Anthracenes



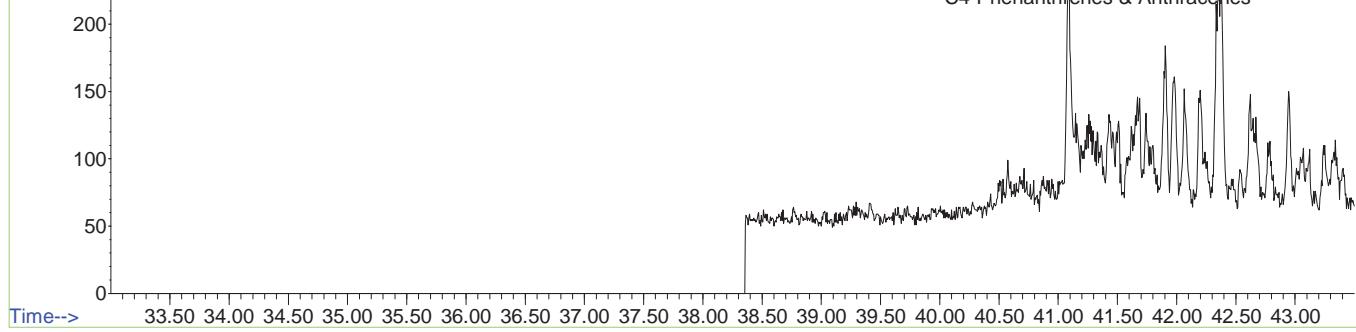
Time--&gt;

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00

Abundance

Ion 234.00 (233.70 to 234.70): S34121.D\data.ms

C4-Phenanthrenes &amp; Anthracenes



Time--&gt;

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00

File: Z:\1\data\S131120\S34121.D

Date Acquired: 20 Nov 2013 7:05 pm

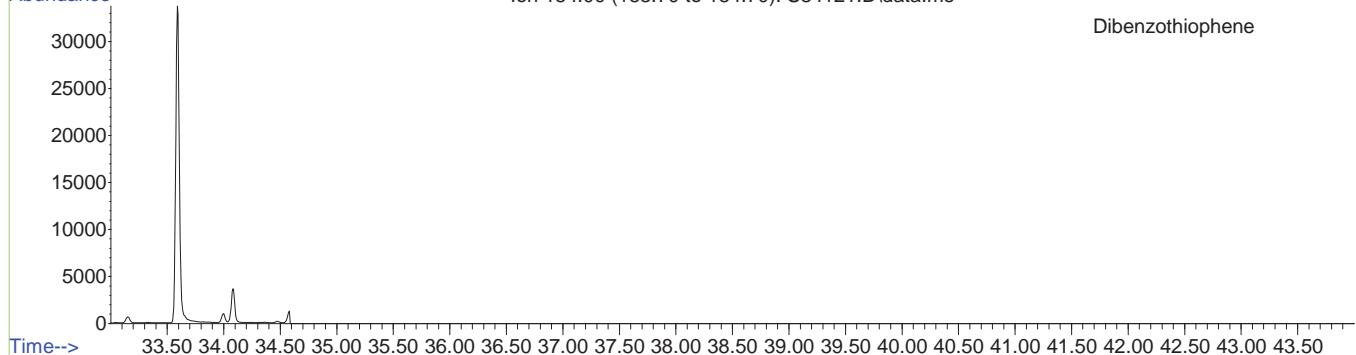
Sample Name: mc26103-1

Misc Info: op35834,mss1523,0.0011,,,10,10

Abundance

Ion 184.00 (183.70 to 184.70): S34121.D\data.ms

Dibenzothiophene



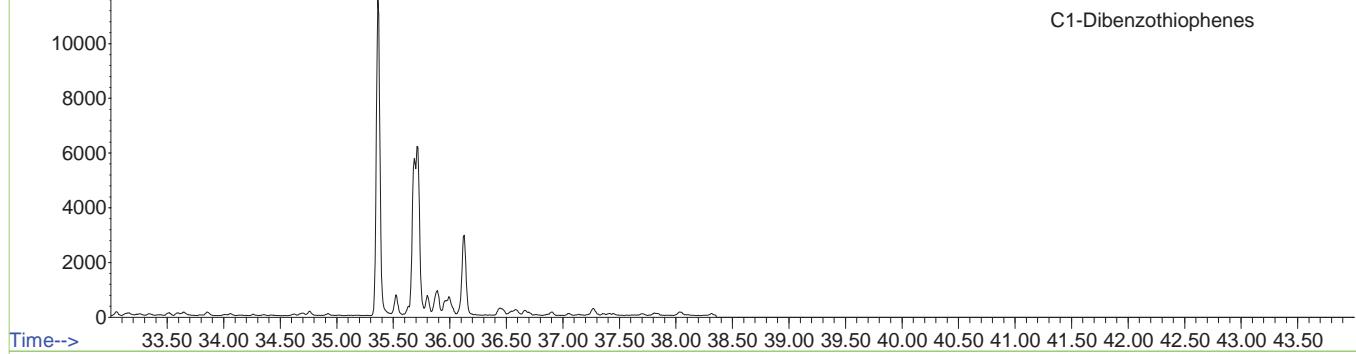
Time--&gt;

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50

Abundance

Ion 198.00 (197.70 to 198.70): S34121.D\data.ms

C1-Dibenzothiophenes



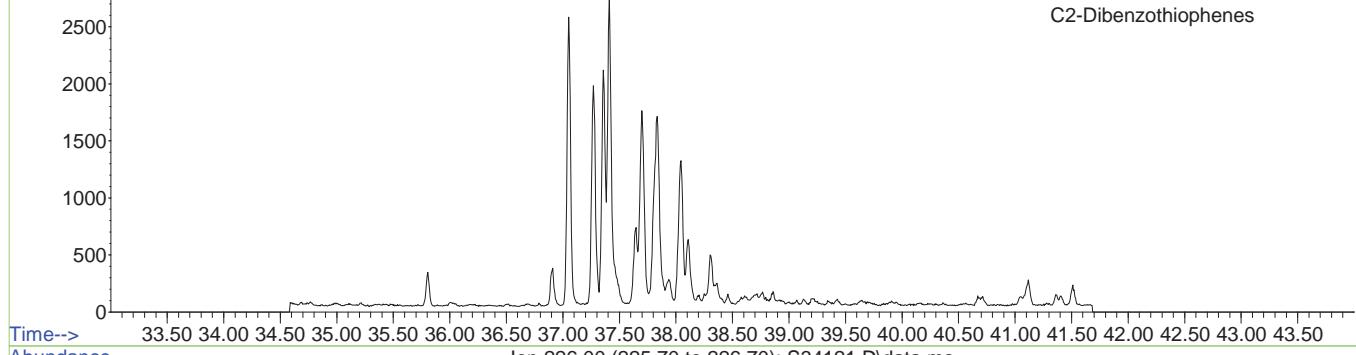
Time--&gt;

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50

Abundance

Ion 212.00 (211.70 to 212.70): S34121.D\data.ms

C2-Dibenzothiophenes



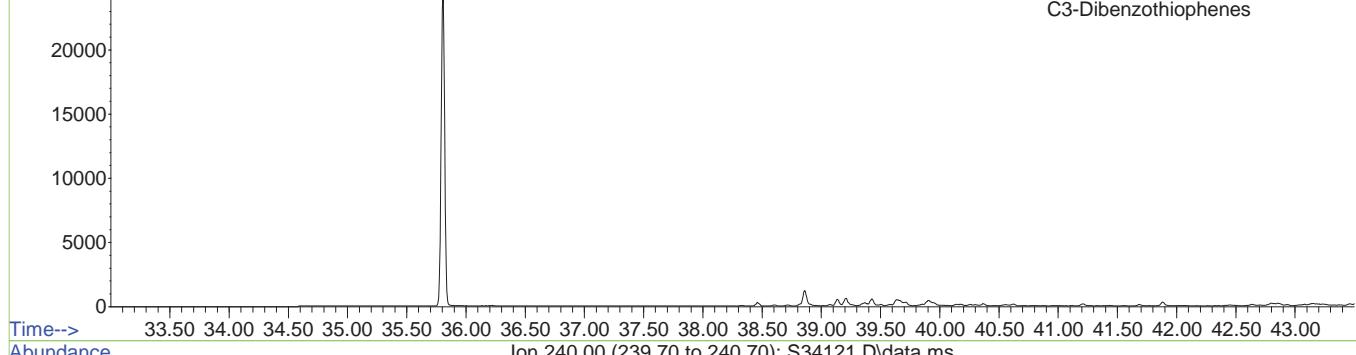
Time--&gt;

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50

Abundance

Ion 226.00 (225.70 to 226.70): S34121.D\data.ms

C3-Dibenzothiophenes



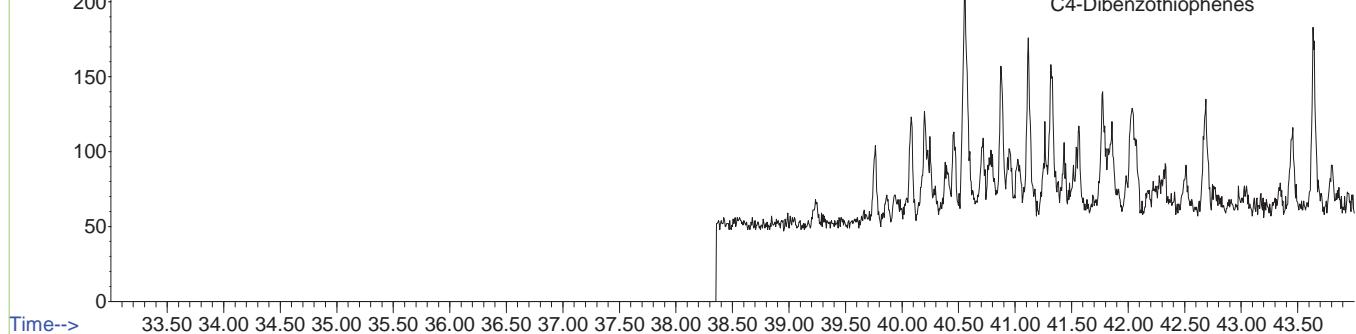
Time--&gt;

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50

Abundance

Ion 240.00 (239.70 to 240.70): S34121.D\data.ms

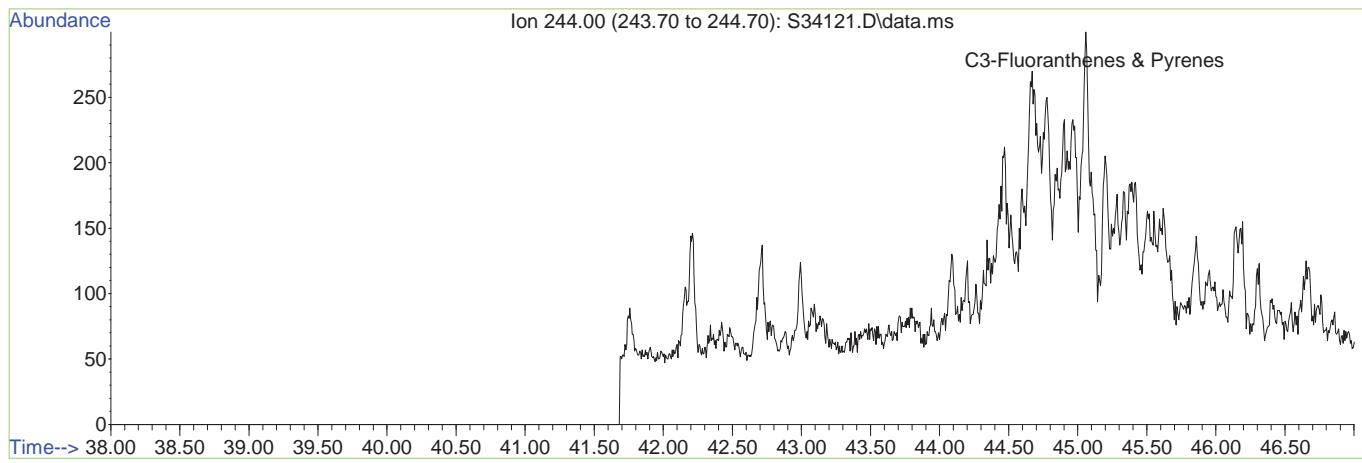
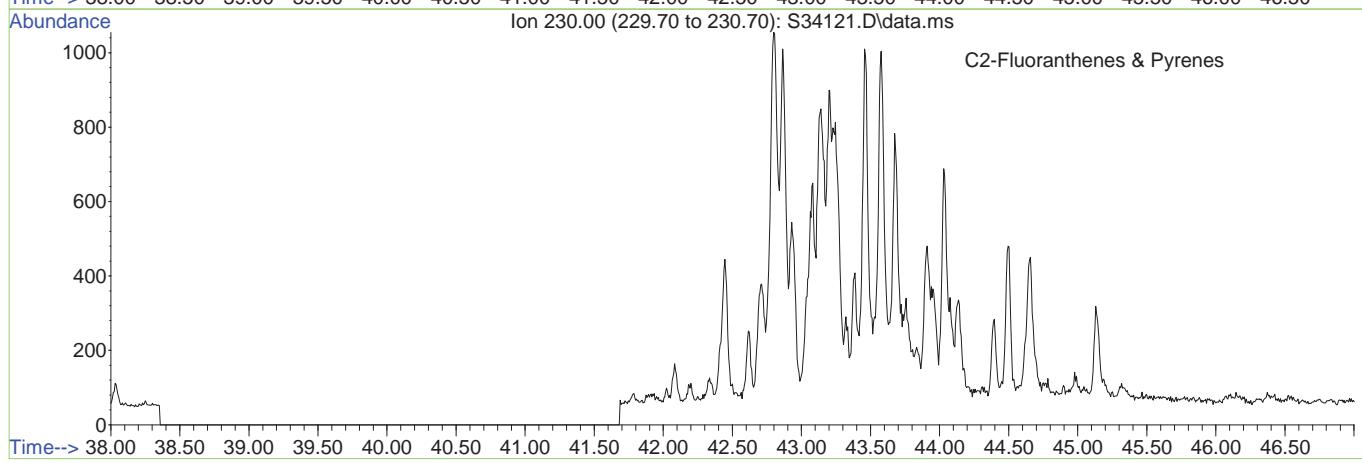
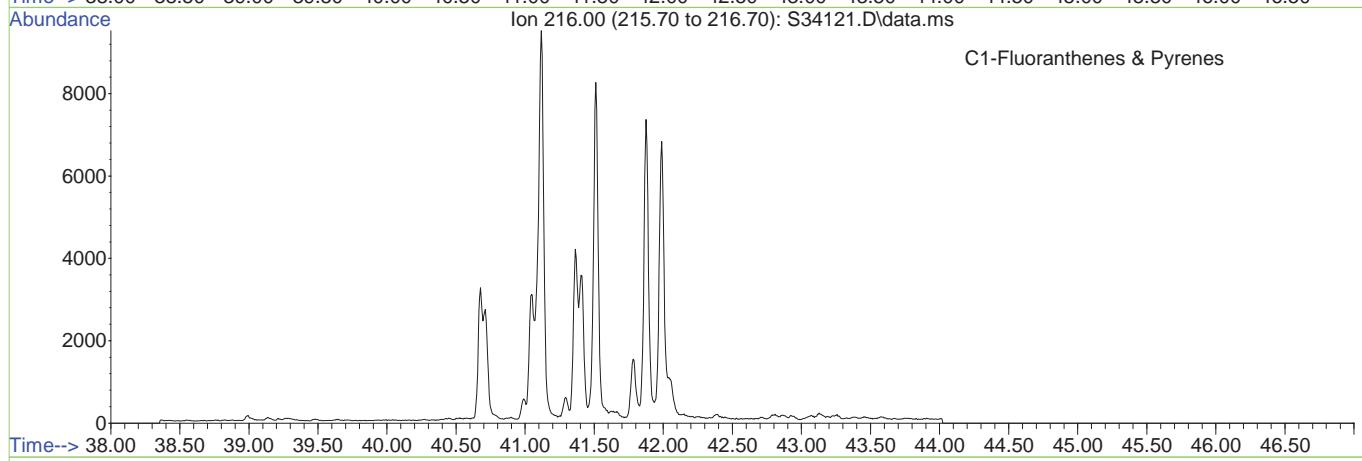
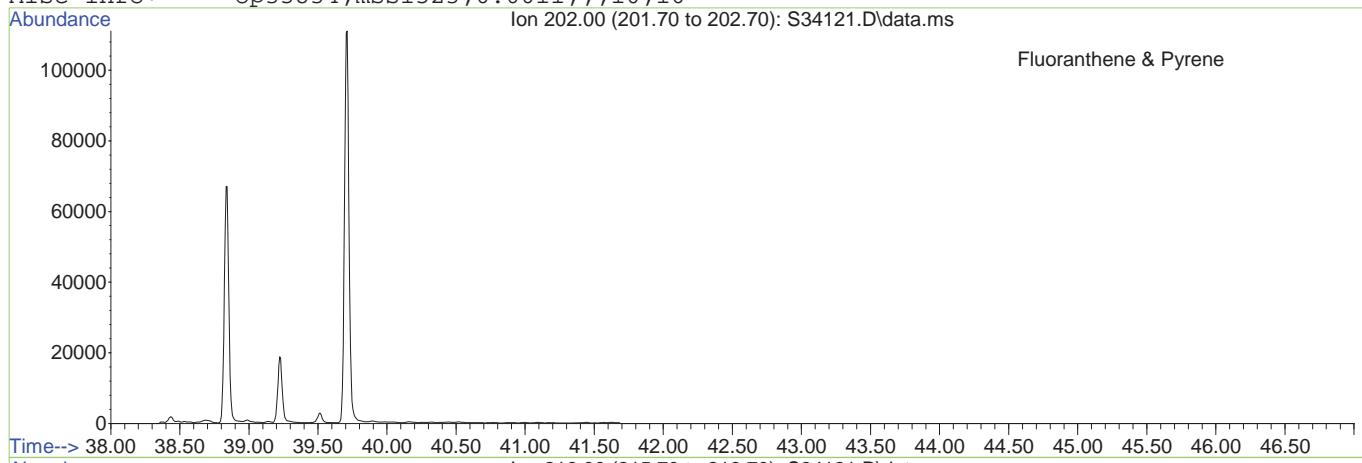
C4-Dibenzothiophenes



Time--&gt;

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50

File: Z:\1\data\S131120\S34121.D  
Date Acquired: 20 Nov 2013 7:05 pm  
Method File: SAPAHSIM-MTBE.M  
Sample Name: mc26103-1  
Misc Info: op35834,mss1523,0.0011,,,10,10



File: Z:\1\data\S131120\S34121.D

Date Acquired: 20 Nov 2013 7:05 pm

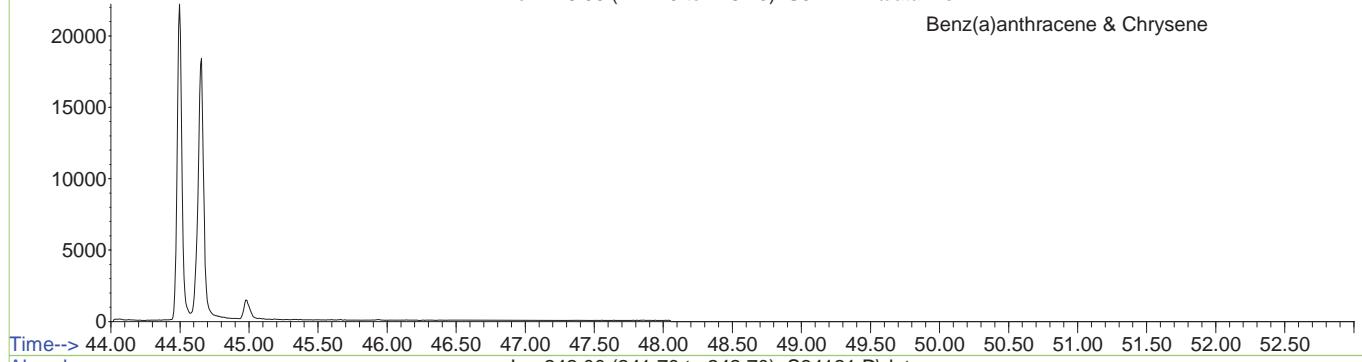
Sample Name: mc26103-1

Misc Info: op35834,mss1523,0.0011,,,10,10

Abundance

Ion 228.00 (227.70 to 228.70): S34121.D\data.ms

Benz(a)anthracene & Chrysene

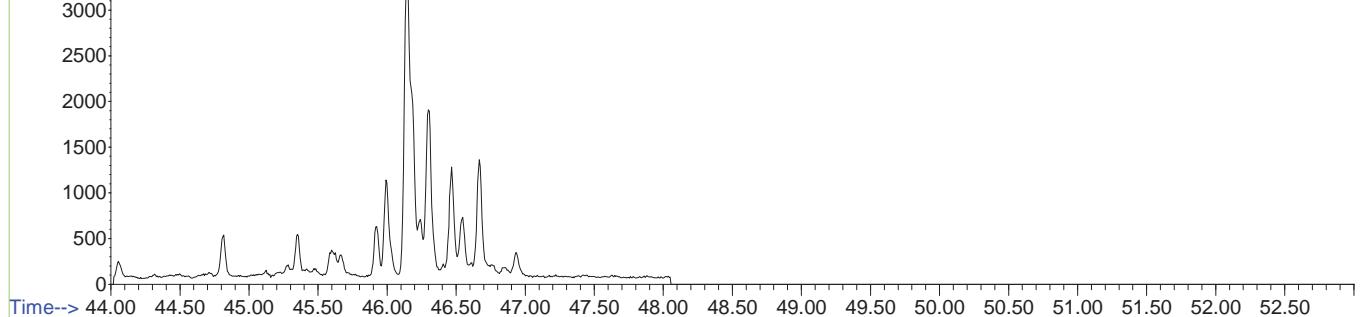


Time--> 44.00 44.50 45.00 45.50 46.00 46.50 47.00 47.50 48.00 48.50 49.00 49.50 50.00 50.50 51.00 51.50 52.00 52.50

Abundance

Ion 242.00 (241.70 to 242.70): S34121.D\data.ms

C1-Benz(a)anthracenes & Chrysenes

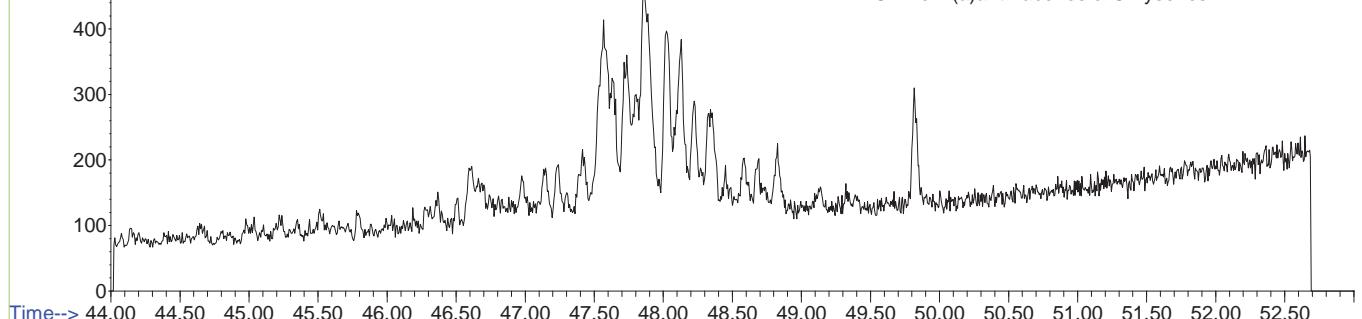


Time--> 44.00 44.50 45.00 45.50 46.00 46.50 47.00 47.50 48.00 48.50 49.00 49.50 50.00 50.50 51.00 51.50 52.00 52.50

Abundance

Ion 256.00 (255.70 to 256.70): S34121.D\data.ms

C2-Benz(a)anthracenes & Chrysenes

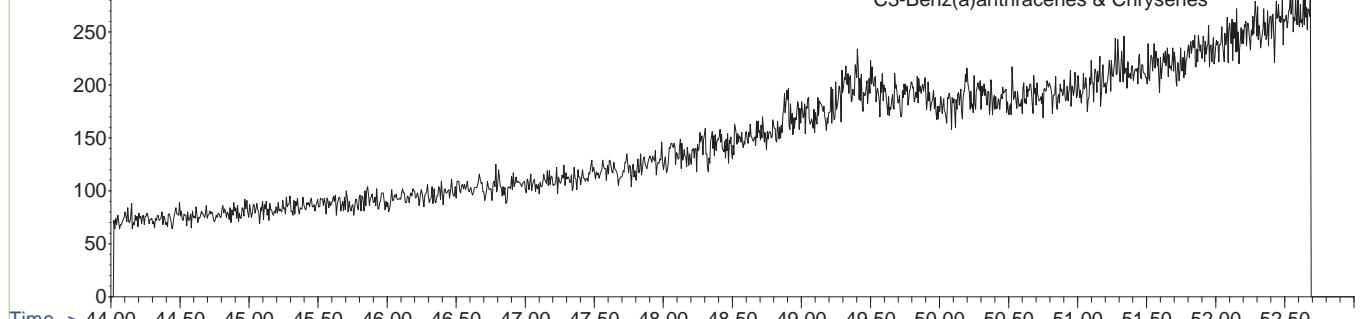


Time--> 44.00 44.50 45.00 45.50 46.00 46.50 47.00 47.50 48.00 48.50 49.00 49.50 50.00 50.50 51.00 51.50 52.00 52.50

Abundance

Ion 270.00 (269.70 to 270.70): S34121.D\data.ms

C3-Benz(a)anthracenes & Chrysenes

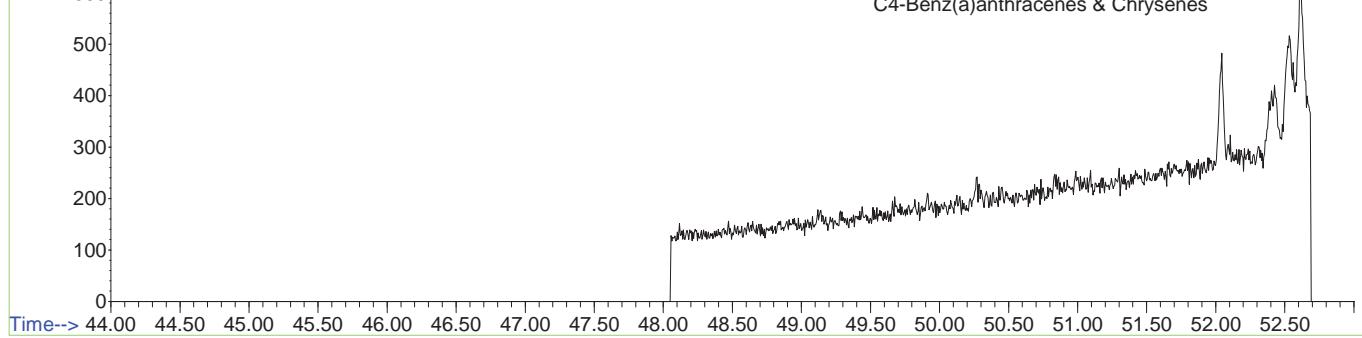


Time--> 44.00 44.50 45.00 45.50 46.00 46.50 47.00 47.50 48.00 48.50 49.00 49.50 50.00 50.50 51.00 51.50 52.00 52.50

Abundance

Ion 284.00 (283.70 to 284.70): S34121.D\data.ms

C4-Benz(a)anthracenes & Chrysenes

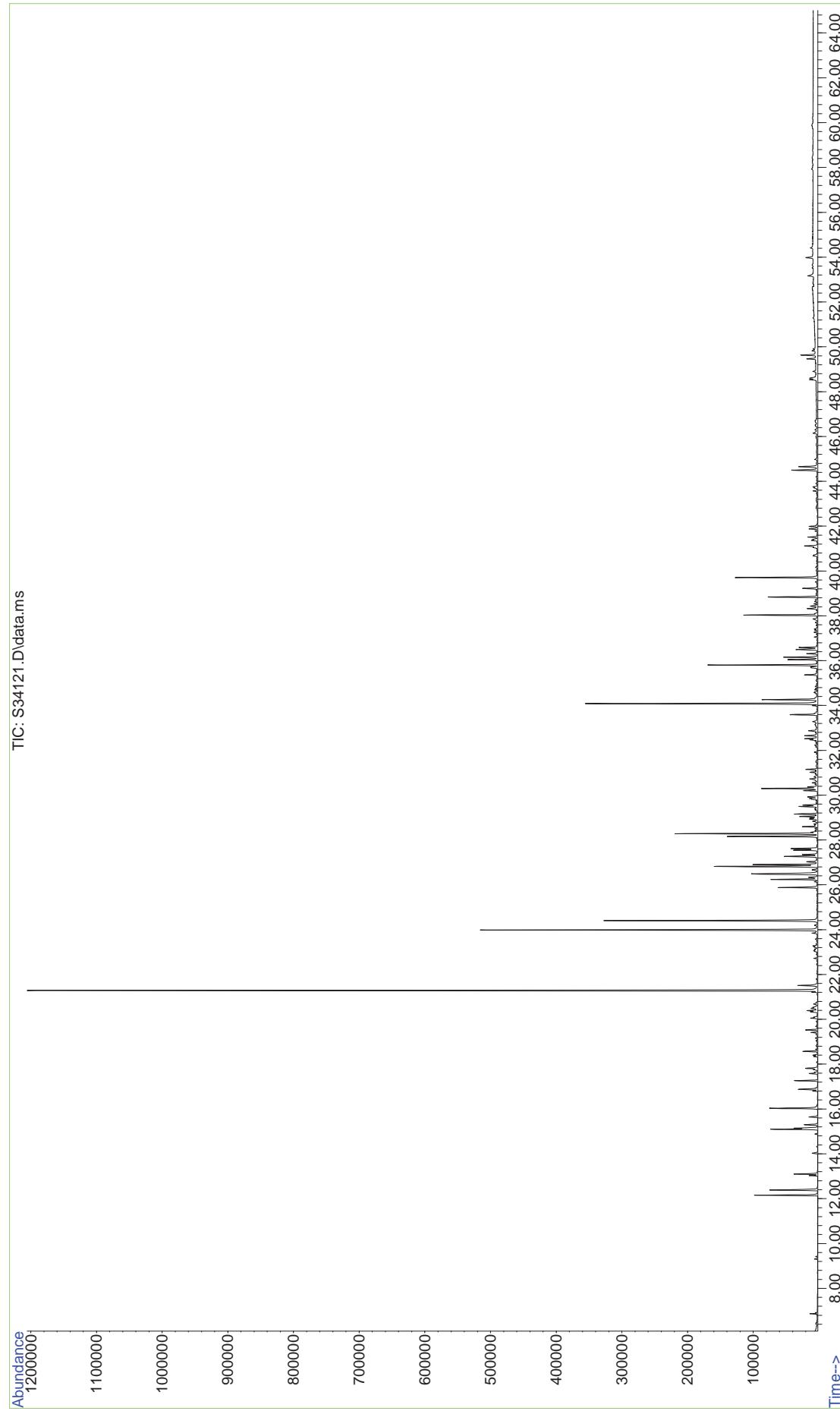


Time--> 44.00 44.50 45.00 45.50 46.00 46.50 47.00 47.50 48.00 48.50 49.00 49.50 50.00 50.50 51.00 51.50 52.00 52.50

## ACCUTEST

## GC/MS TOTAL ION CHROMATOGRAM

File: Z:\1\data\S131120\S34121.D  
Date Acquired: 20 Nov 2013 7:05 pm  
Method File: SAPAHSIM-MTBE.M  
Sample Name: mc26103-1  
Misc Info: op35834,msss1523,0.0011,,10,10



File: Z:\1\data\S131120\S34120.D

Date Acquired: 20 Nov 2013 5:53 pm

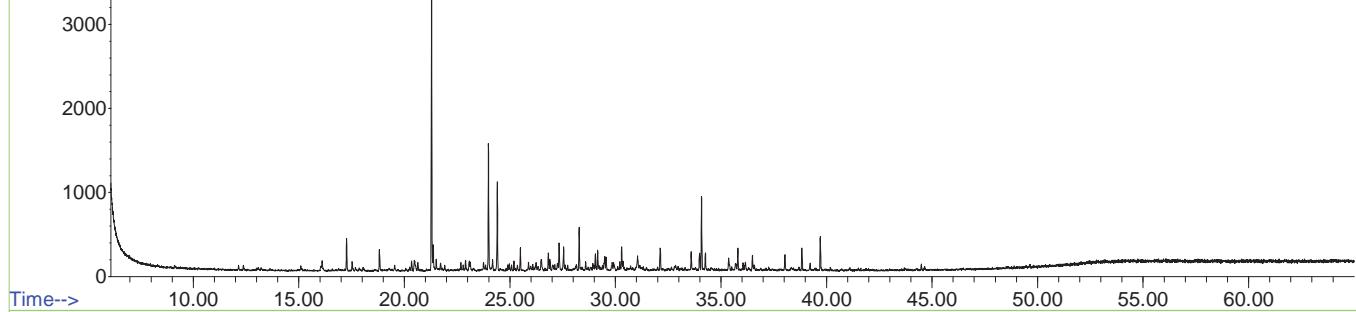
Sample Name: op35834-dup1

Misc Info: op35834,mss1523,0.0015,,,10,10

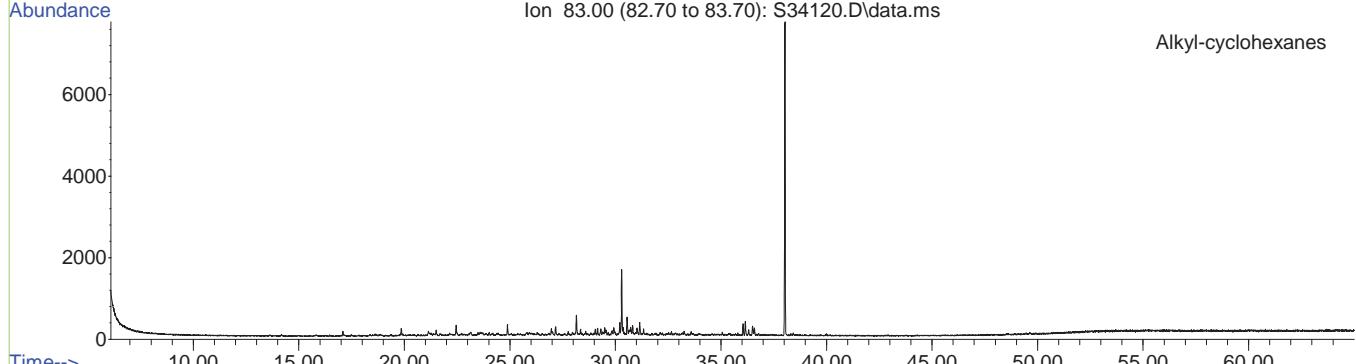
Abundance

Ion 85.00 (84.70 to 85.70): S34120.D\data.ms

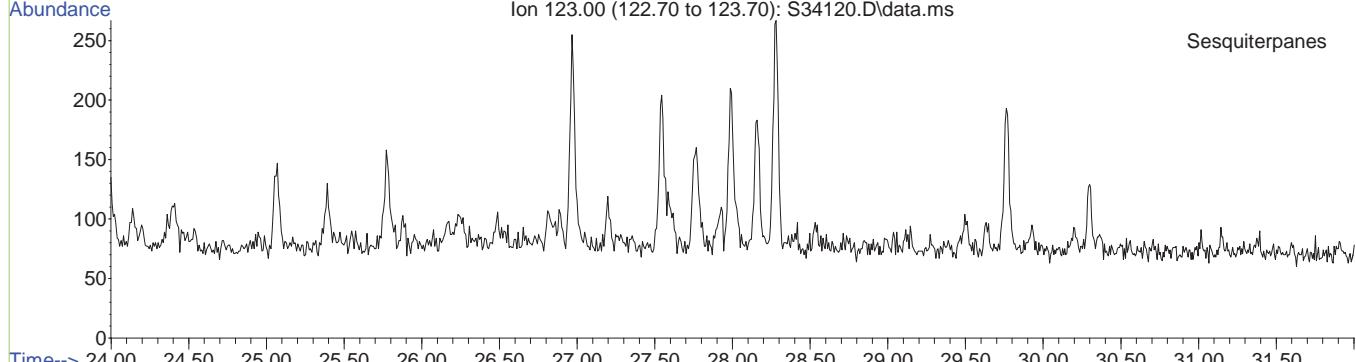
Alkanes



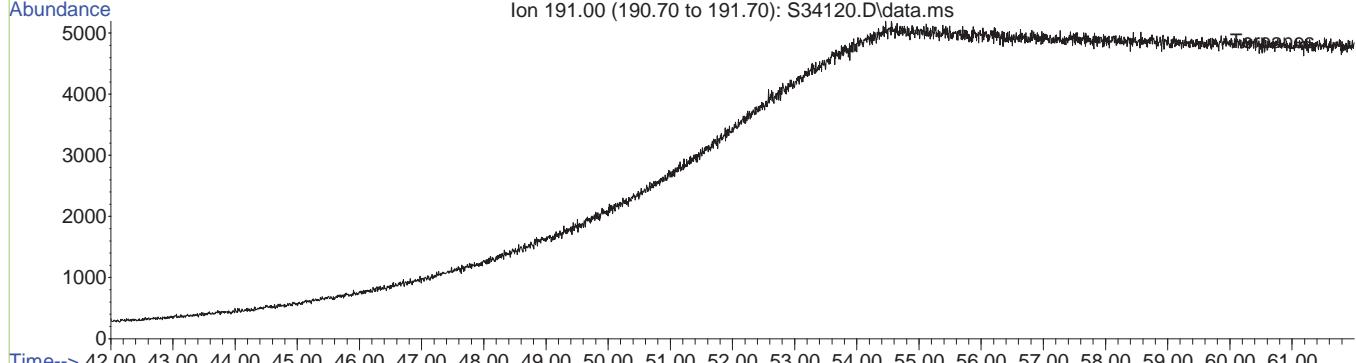
Time--&gt;



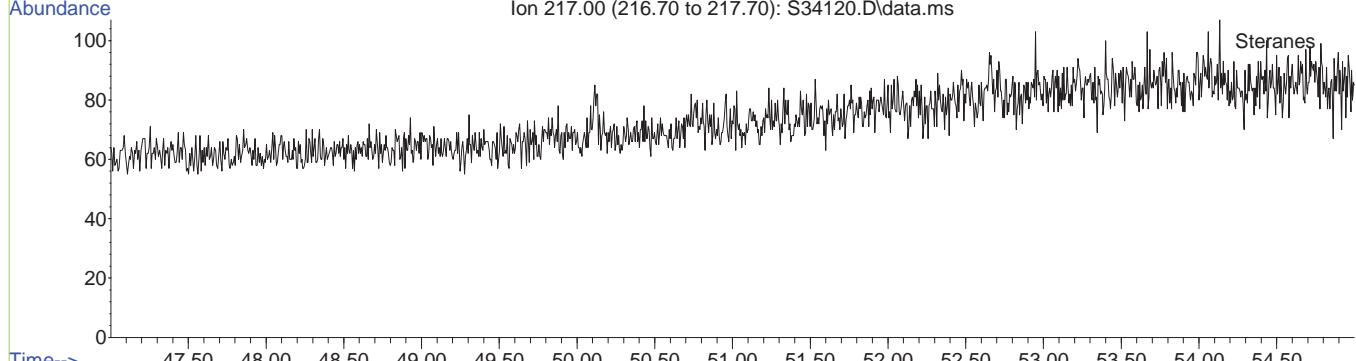
Time--&gt;



Time--&gt;



Time--&gt;



Time--&gt;

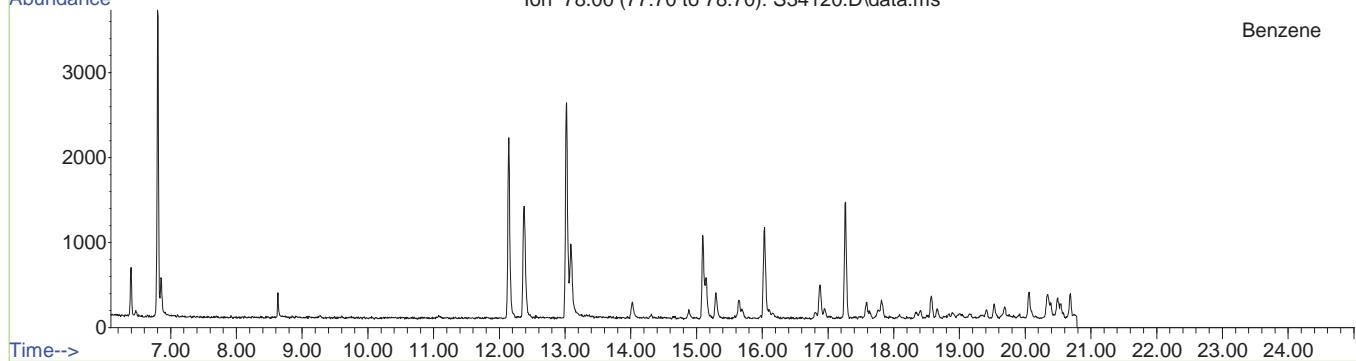
File: Z:\1\data\S131120\S34120.D

Date Acquired: 20 Nov 2013 5:53 pm

Sample Name: op35834-dup1

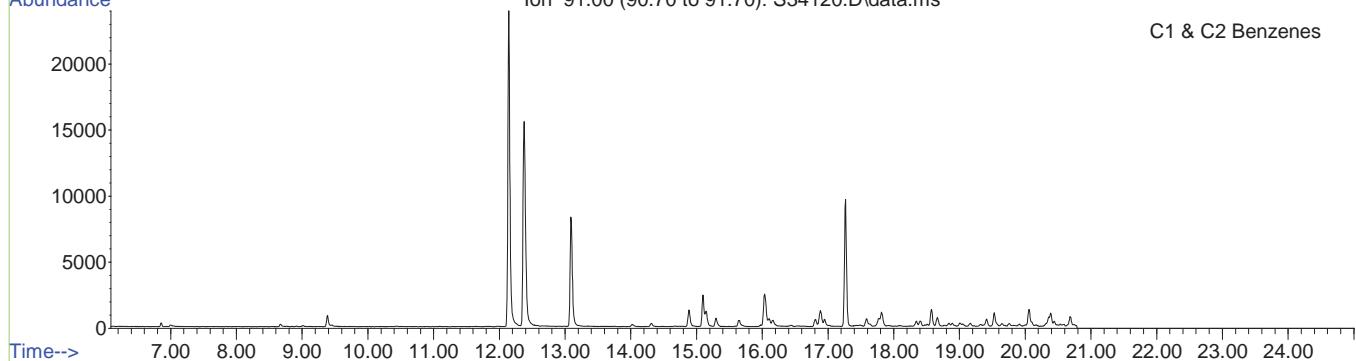
Misc Info: op35834,mss1523,0.0015,,,10,10

Abundance Ion 78.00 (77.70 to 78.70): S34120.D\data.ms



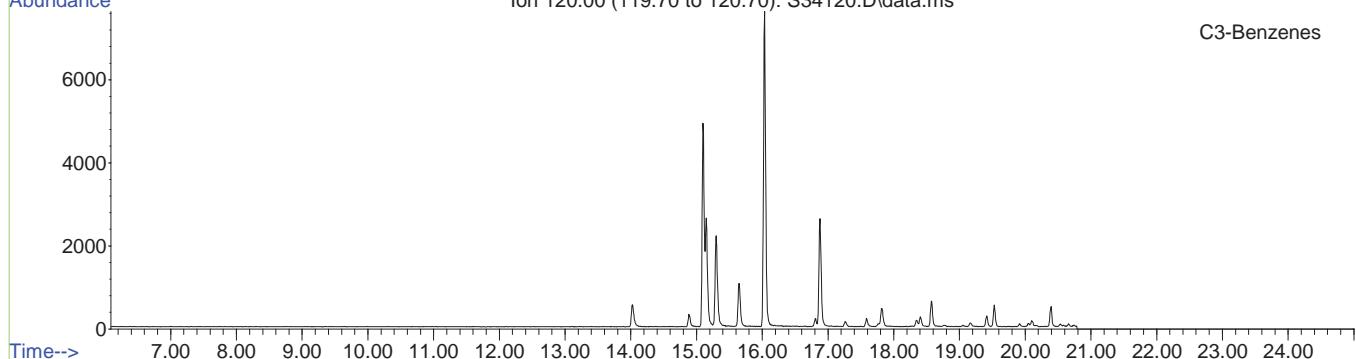
Benzene

Abundance Ion 91.00 (90.70 to 91.70): S34120.D\data.ms



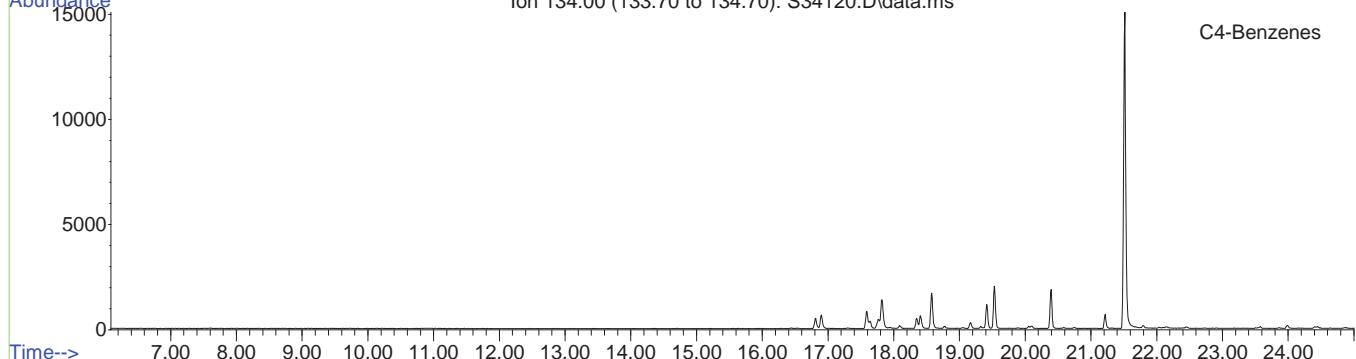
C1 &amp; C2 Benzenes

Abundance Ion 120.00 (119.70 to 120.70): S34120.D\data.ms



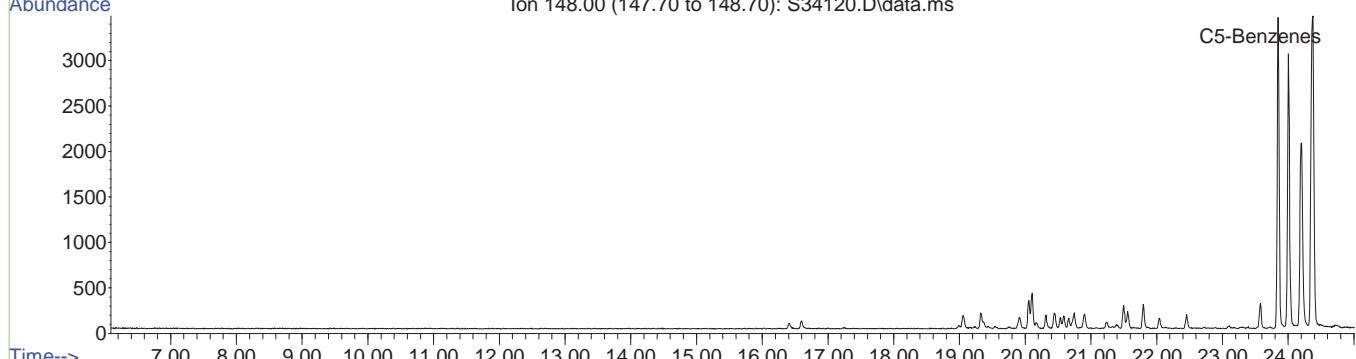
C3-Benzene

Abundance Ion 134.00 (133.70 to 134.70): S34120.D\data.ms



C4-Benzene

Abundance Ion 148.00 (147.70 to 148.70): S34120.D\data.ms



C5-Benzene

File: Z:\1\data\S131120\S34120.D

Date Acquired: 20 Nov 2013 5:53 pm

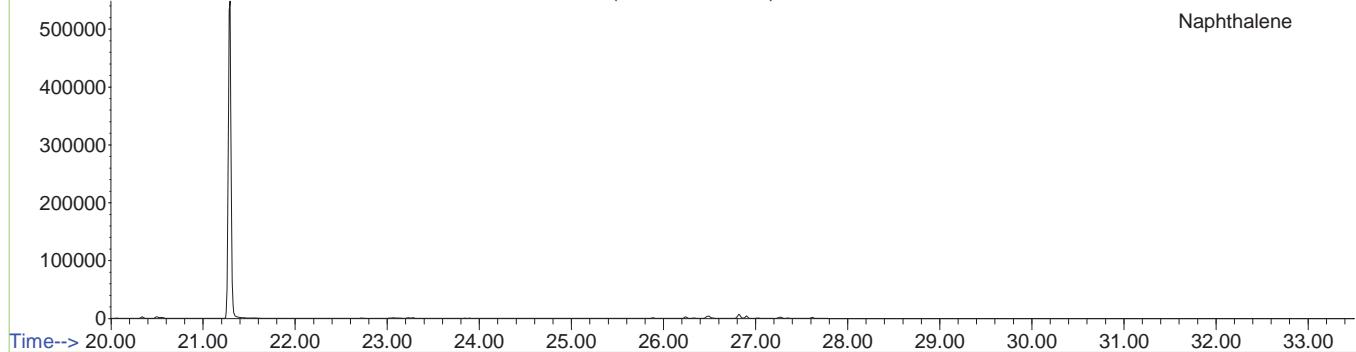
Sample Name: op35834-dup1

Misc Info: op35834,mss1523,0.0015,,,10,10

Abundance

Ion 128.00 (127.70 to 128.70): S34120.D\data.ms

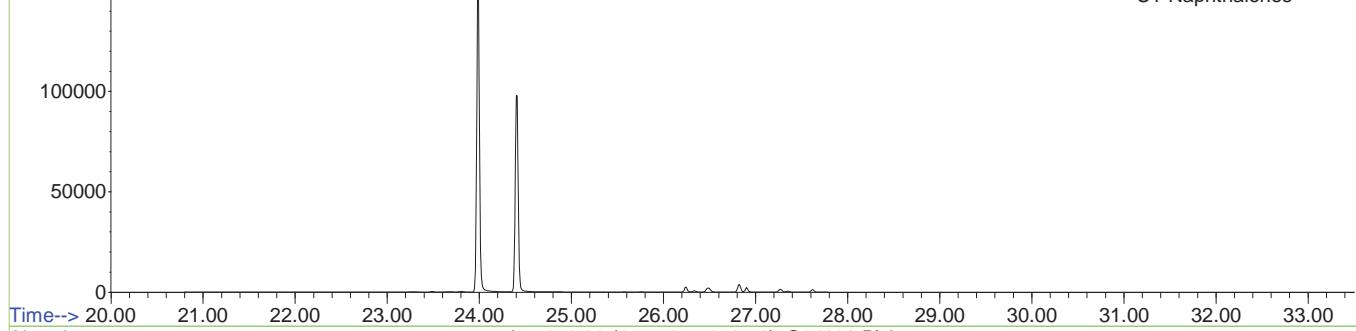
Naphthalene



Abundance

Ion 142.00 (141.70 to 142.70): S34120.D\data.ms

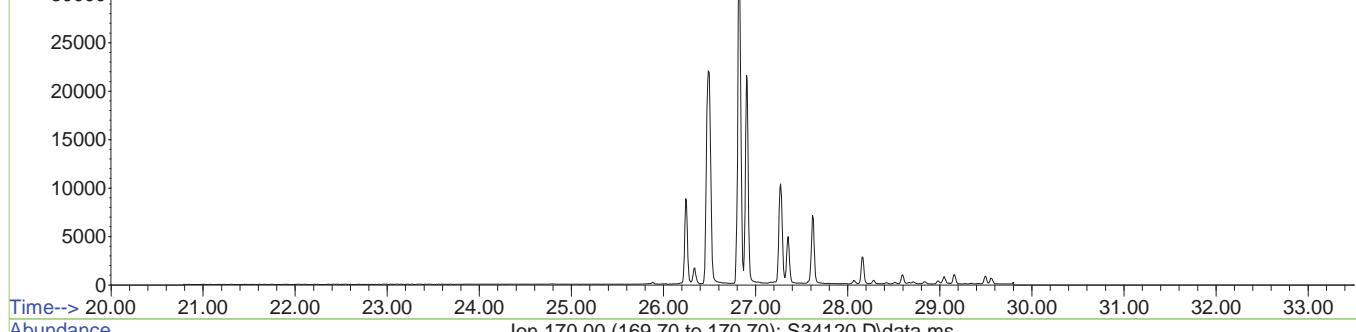
C1-Naphthalenes



Abundance

Ion 156.00 (155.70 to 156.70): S34120.D\data.ms

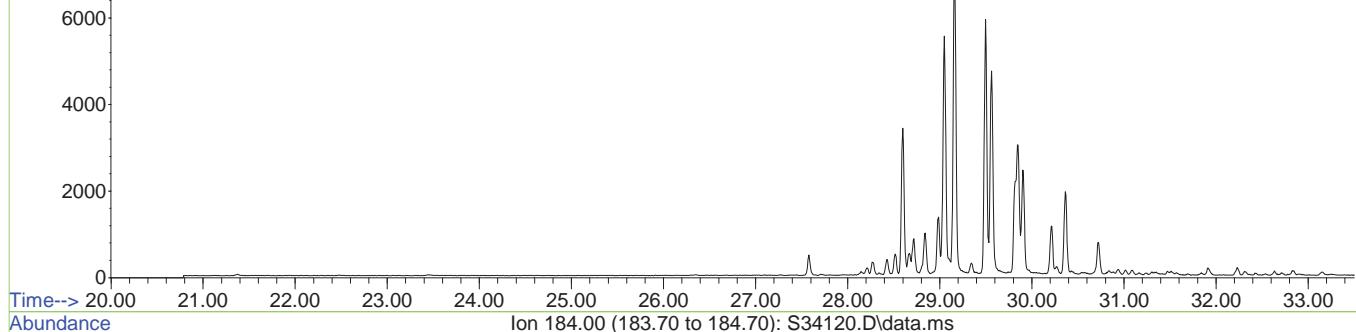
C2-Naphthalenes



Abundance

Ion 170.00 (169.70 to 170.70): S34120.D\data.ms

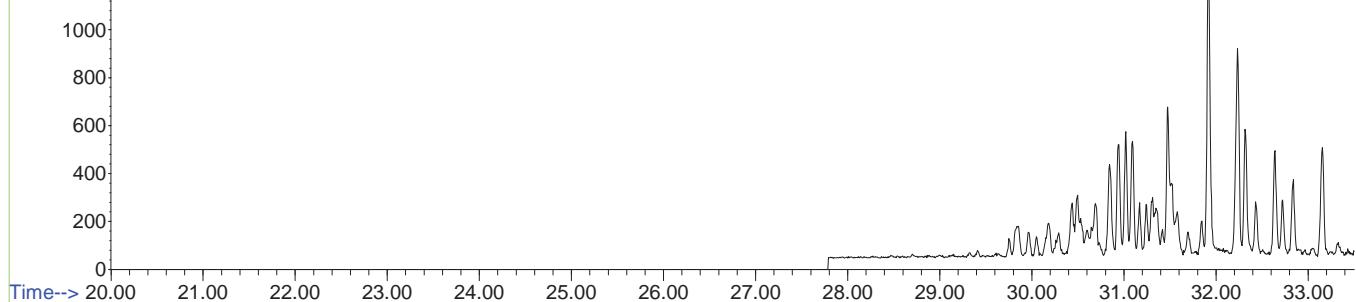
C3-Naphthalenes



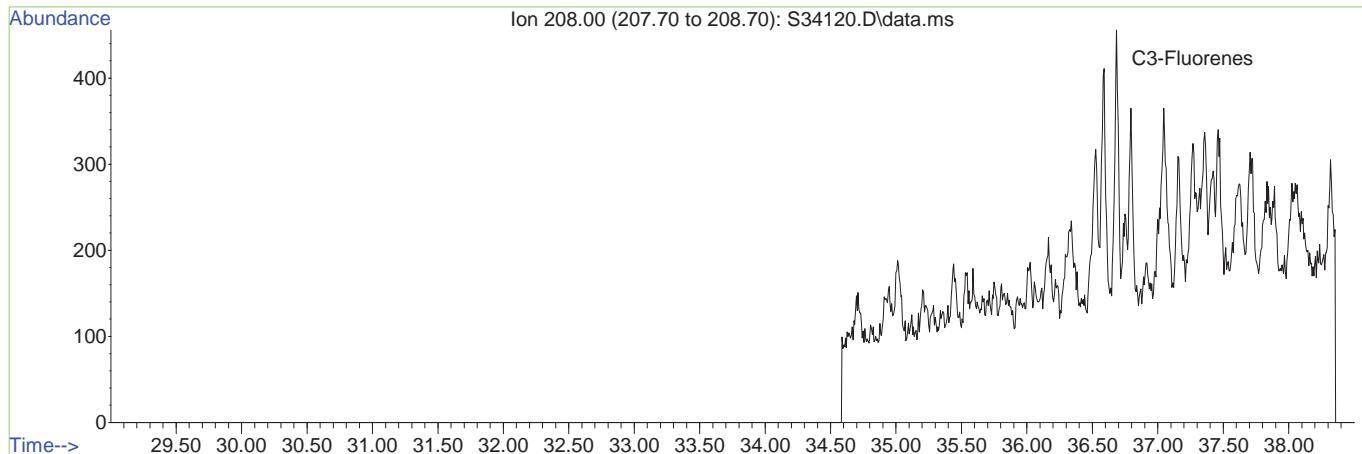
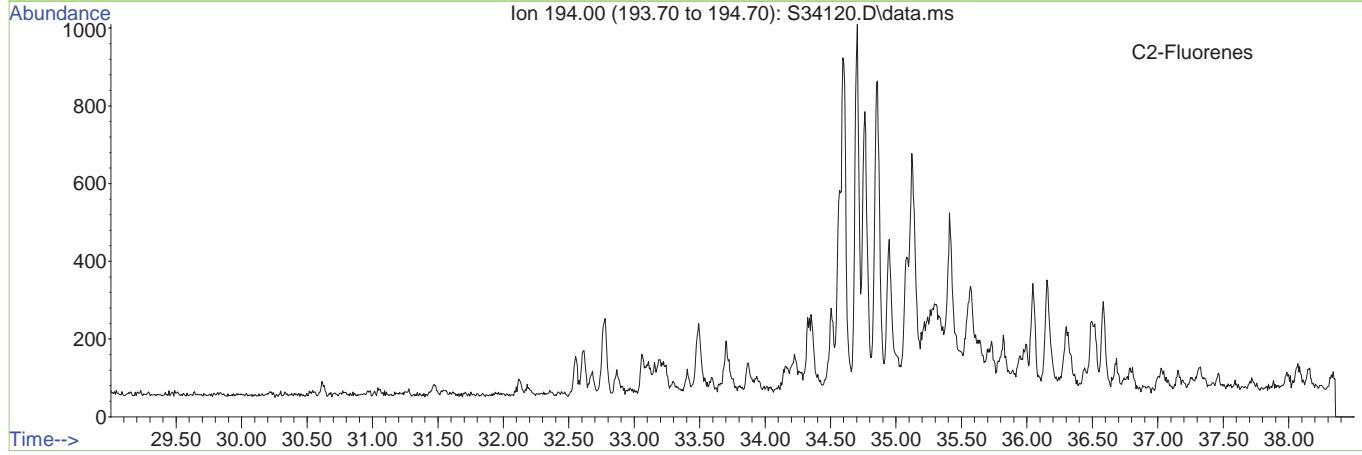
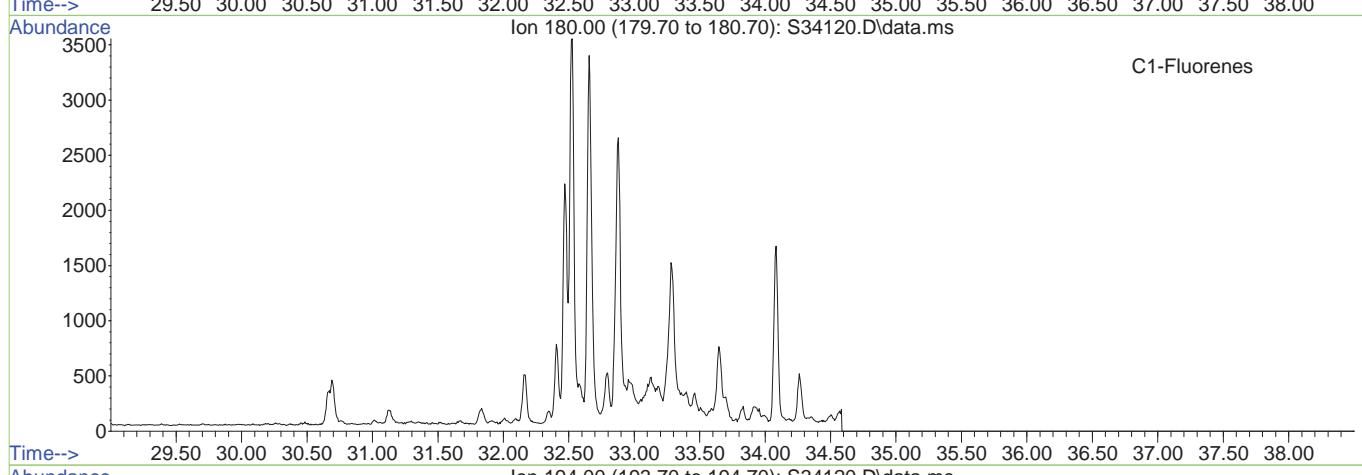
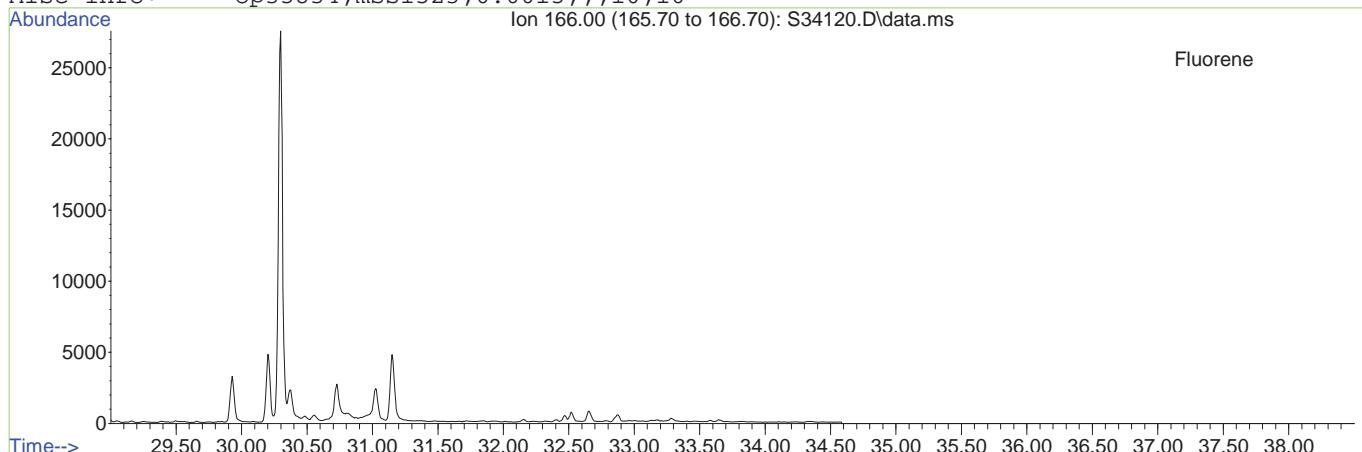
Abundance

Ion 184.00 (183.70 to 184.70): S34120.D\data.ms

C4-Naphthalenes



File: Z:\1\data\S131120\S34120.D  
Date Acquired: 20 Nov 2013 5:53 pm  
Method File: SAPAHSIM-MTBE.M  
Sample Name: op35834-dup1  
Misc Info: op35834,mss1523,0.0015,,,10,10



File: Z:\1\data\S131120\S34120.D

Date Acquired: 20 Nov 2013 5:53 pm

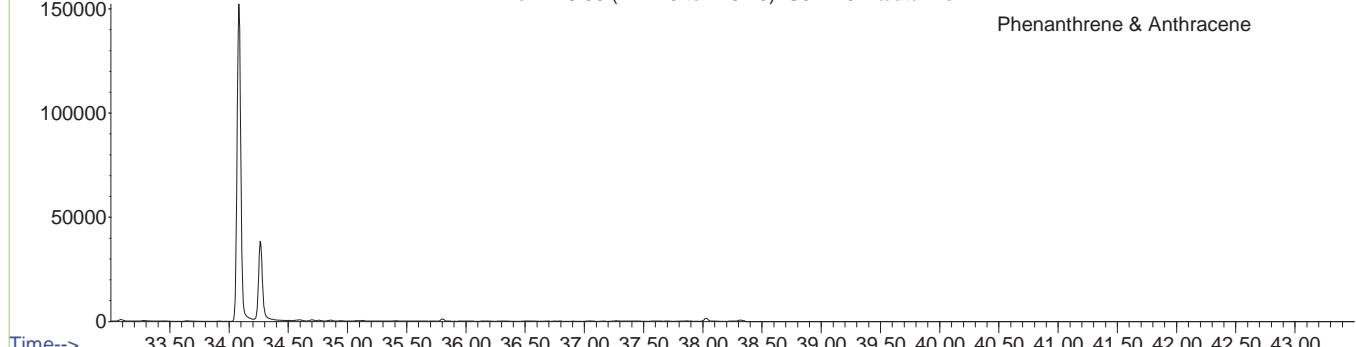
Sample Name: op35834-dup1

Misc Info: op35834,mss1523,0.0015,,,10,10

Abundance

Ion 178.00 (177.70 to 178.70): S34120.D\data.ms

Phenanthrene &amp; Anthracene

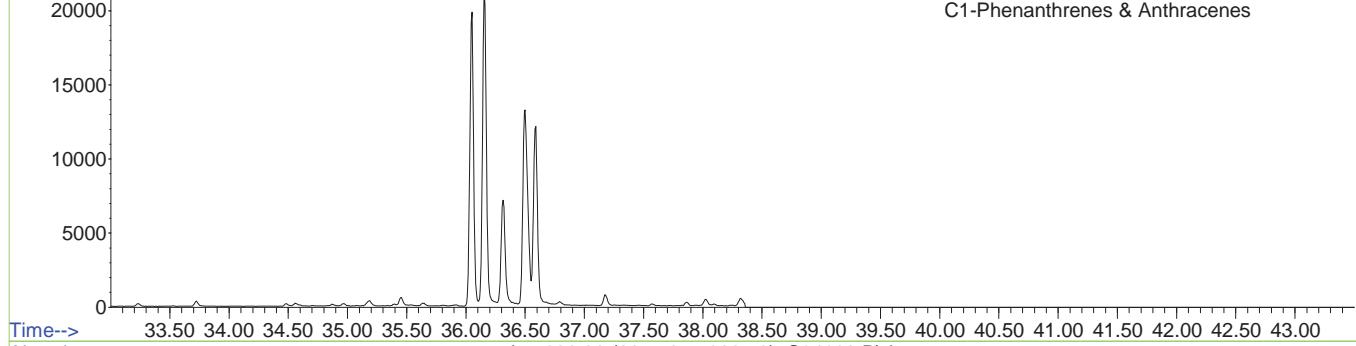


Time--&gt;

Abundance

Ion 192.00 (191.70 to 192.70): S34120.D\data.ms

C1-Phenanthrenes &amp; Anthracenes

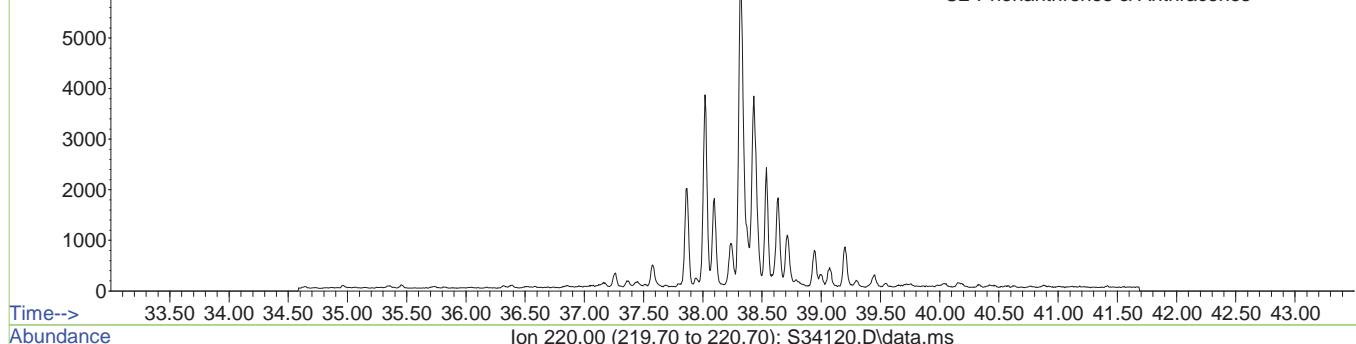


Time--&gt;

Abundance

Ion 206.00 (205.70 to 206.70): S34120.D\data.ms

C2-Phenanthrenes &amp; Anthracenes

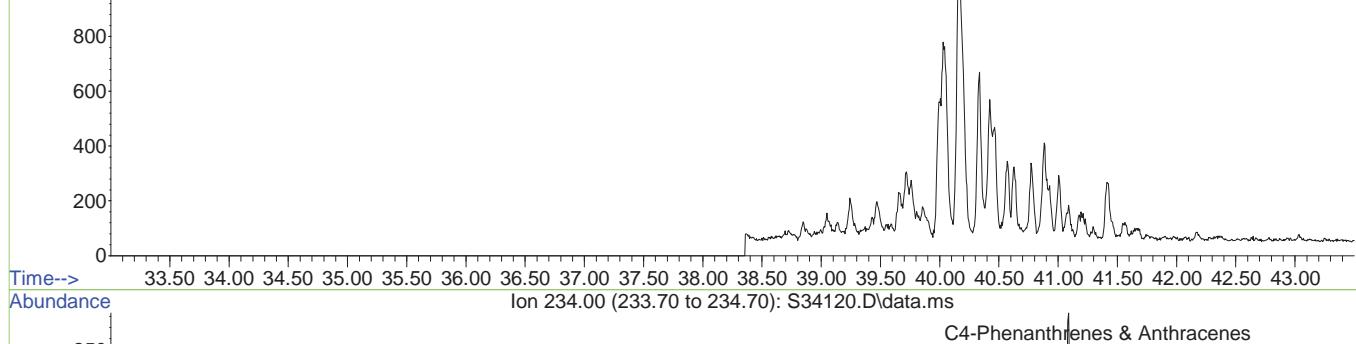


Time--&gt;

Abundance

Ion 220.00 (219.70 to 220.70): S34120.D\data.ms

C3-Phenanthrenes &amp; Anthracenes

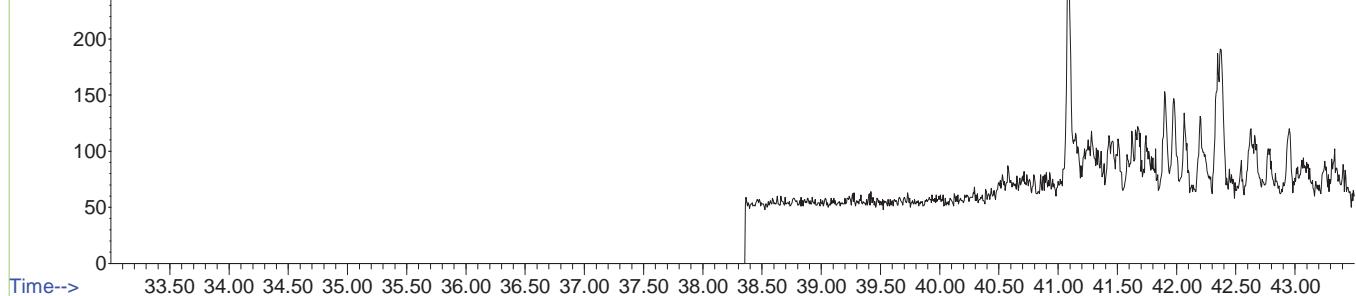


Time--&gt;

Abundance

Ion 234.00 (233.70 to 234.70): S34120.D\data.ms

C4-Phenanthrenes &amp; Anthracenes



Time--&gt;

File: Z:\1\data\S131120\S34120.D

Date Acquired: 20 Nov 2013 5:53 pm

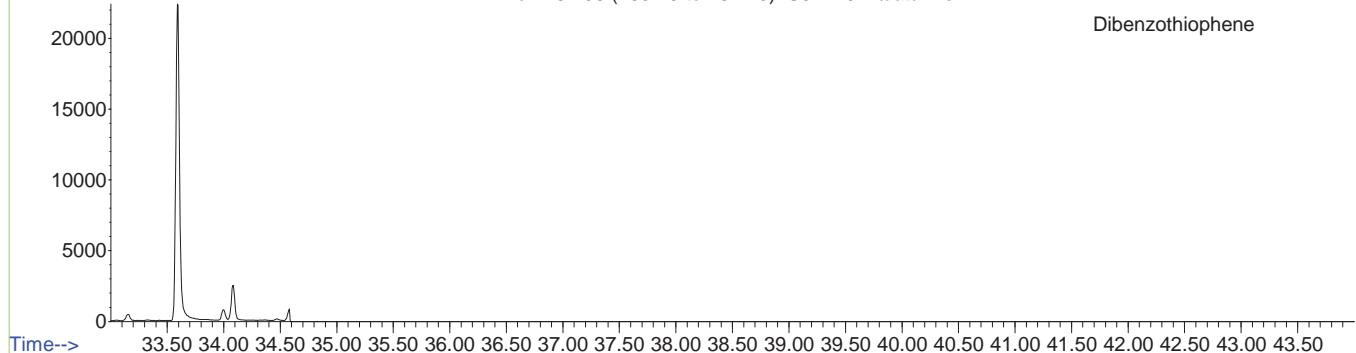
Sample Name: op35834-dup1

Misc Info: op35834,mss1523,0.0015,,,10,10

Abundance

Ion 184.00 (183.70 to 184.70): S34120.D\data.ms

Dibenzothiophene



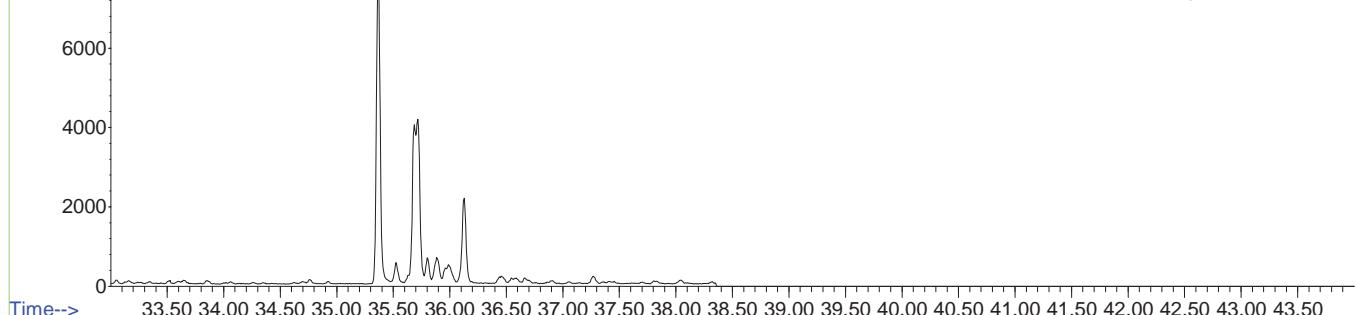
Time-->

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50

Abundance

Ion 198.00 (197.70 to 198.70): S34120.D\data.ms

C1-Dibenzothiophenes



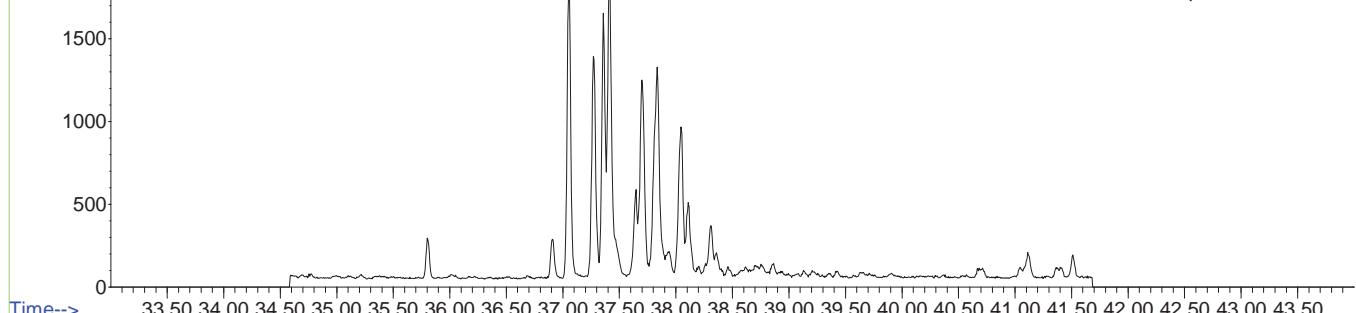
Time-->

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50

Abundance

Ion 212.00 (211.70 to 212.70): S34120.D\data.ms

C2-Dibenzothiophenes



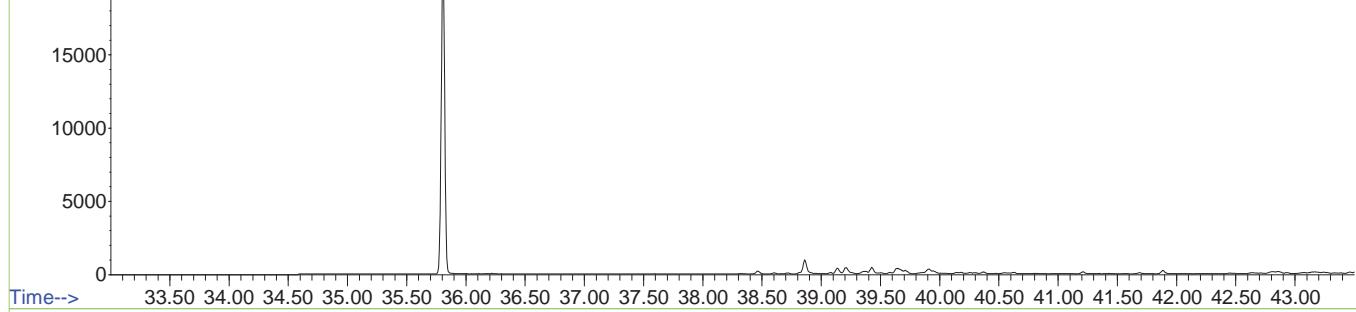
Time-->

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50

Abundance

Ion 226.00 (225.70 to 226.70): S34120.D\data.ms

C3-Dibenzothiophenes



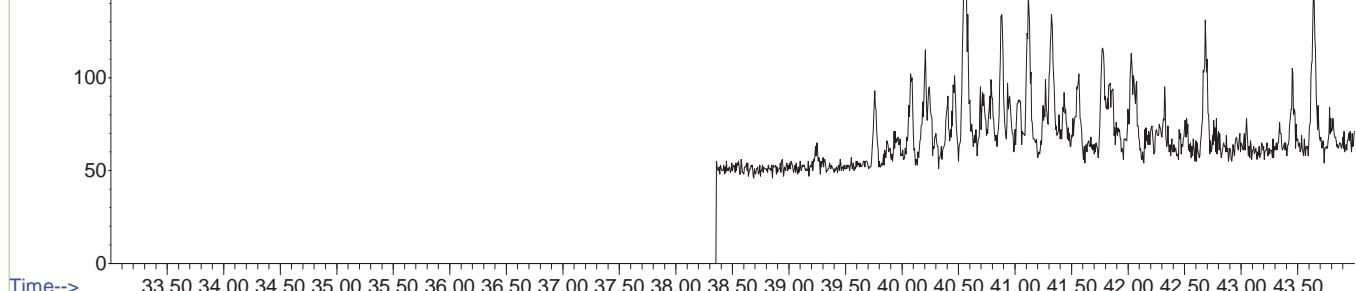
Time-->

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50

Abundance

Ion 240.00 (239.70 to 240.70): S34120.D\data.ms

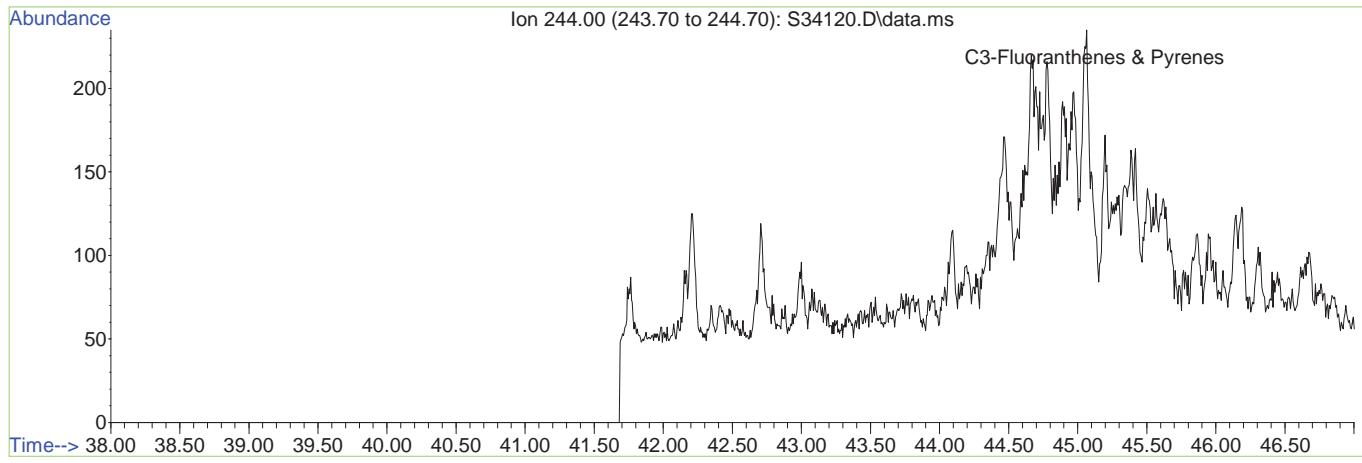
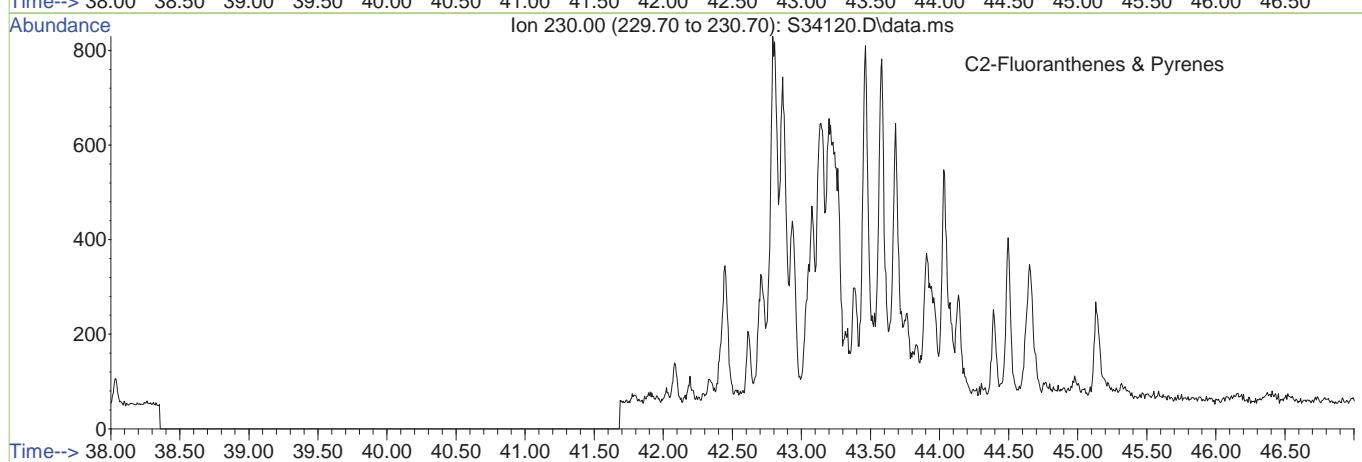
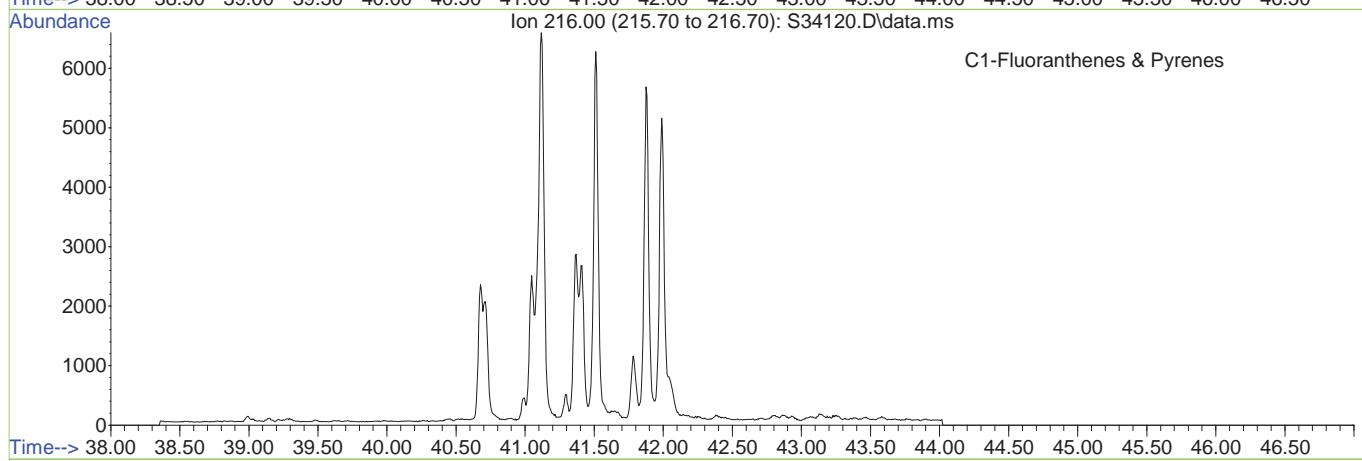
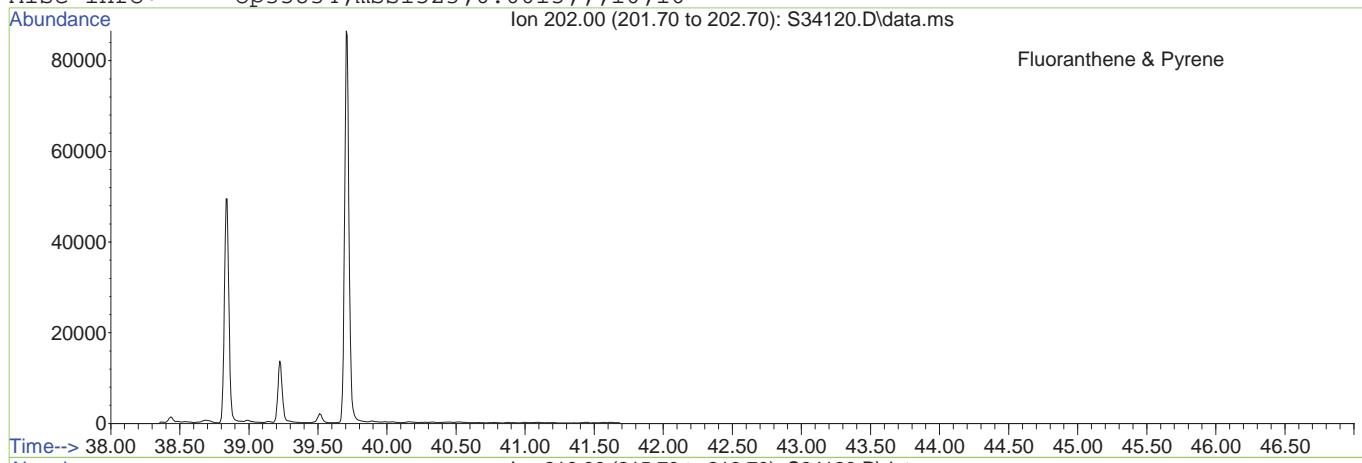
C4-Dibenzothiophenes



Time-->

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50

File: Z:\1\data\S131120\S34120.D  
Date Acquired: 20 Nov 2013 5:53 pm  
Method File: SAPAHSIM-MTBE.M  
Sample Name: op35834-dup1  
Misc Info: op35834,mss1523,0.0015,,,10,10



File: Z:\1\data\S131120\S34120.D

Date Acquired: 20 Nov 2013 5:53 pm

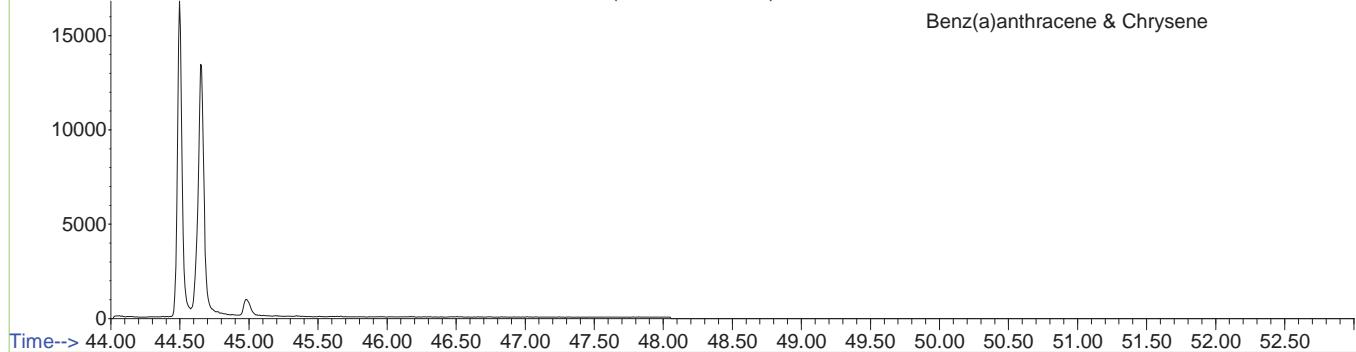
Sample Name: op35834-dup1

Misc Info: op35834,mss1523,0.0015,,,10,10

Abundance

Ion 228.00 (227.70 to 228.70): S34120.D\data.ms

Benz(a)anthracene &amp; Chrysene

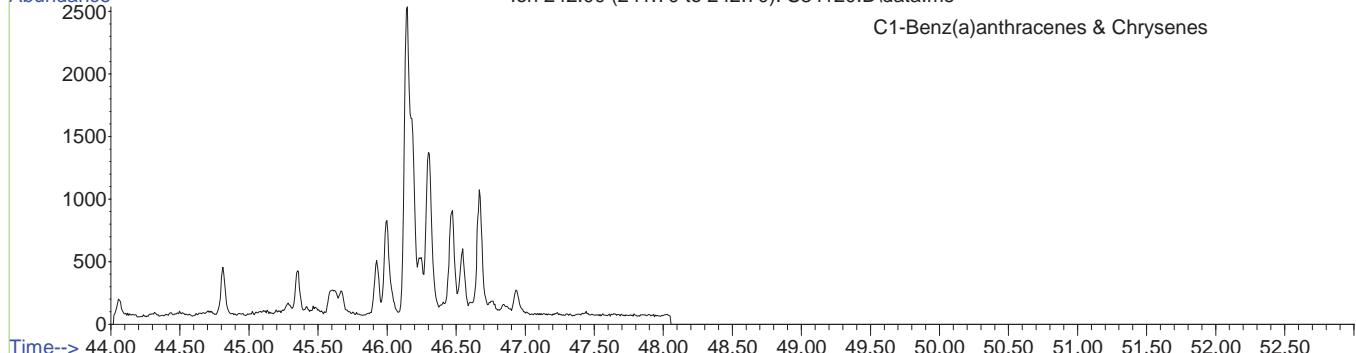


Time--&gt; 44.00 44.50 45.00 45.50 46.00 46.50 47.00 47.50 48.00 48.50 49.00 49.50 50.00 50.50 51.00 51.50 52.00 52.50

Abundance

Ion 242.00 (241.70 to 242.70): S34120.D\data.ms

C1-Benz(a)anthracenes &amp; Chrysenes

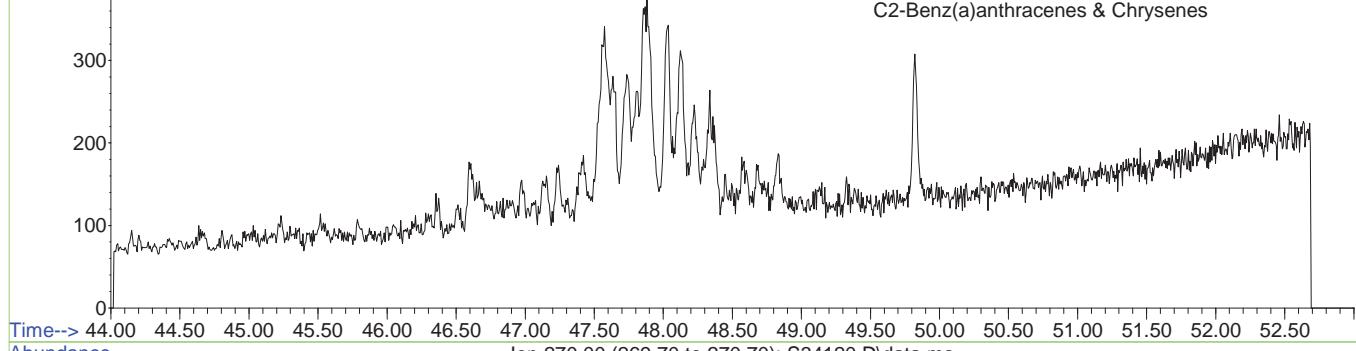


Time--&gt; 44.00 44.50 45.00 45.50 46.00 46.50 47.00 47.50 48.00 48.50 49.00 49.50 50.00 50.50 51.00 51.50 52.00 52.50

Abundance

Ion 256.00 (255.70 to 256.70): S34120.D\data.ms

C2-Benz(a)anthracenes &amp; Chrysenes

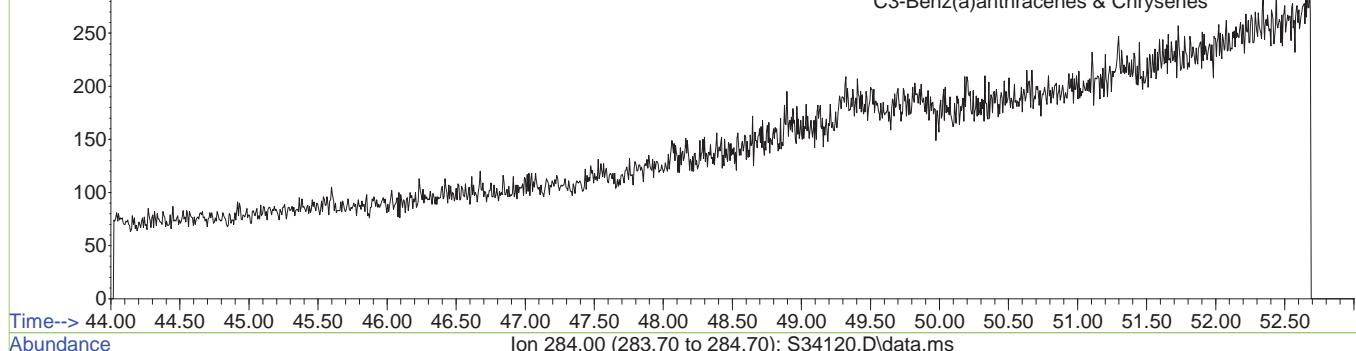


Time--&gt; 44.00 44.50 45.00 45.50 46.00 46.50 47.00 47.50 48.00 48.50 49.00 49.50 50.00 50.50 51.00 51.50 52.00 52.50

Abundance

Ion 270.00 (269.70 to 270.70): S34120.D\data.ms

C3-Benz(a)anthracenes &amp; Chrysenes

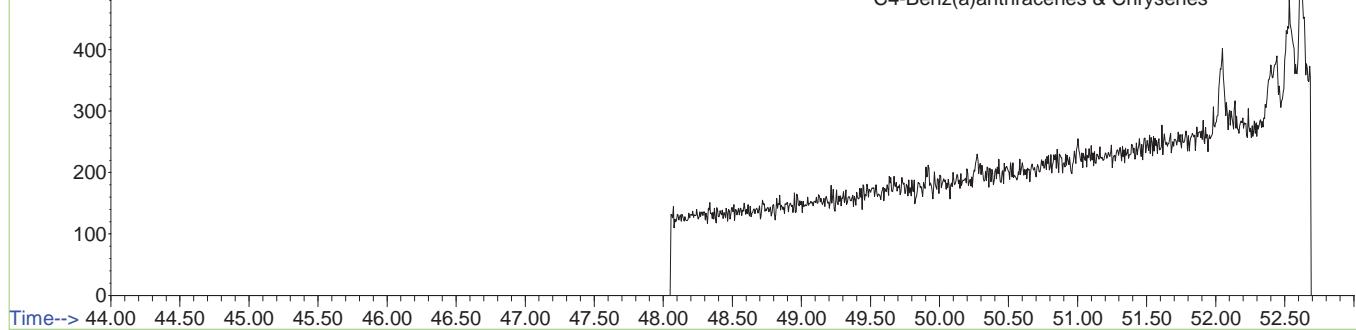


Time--&gt; 44.00 44.50 45.00 45.50 46.00 46.50 47.00 47.50 48.00 48.50 49.00 49.50 50.00 50.50 51.00 51.50 52.00 52.50

Abundance

Ion 284.00 (283.70 to 284.70): S34120.D\data.ms

C4-Benz(a)anthracenes &amp; Chrysenes



Time--&gt; 44.00 44.50 45.00 45.50 46.00 46.50 47.00 47.50 48.00 48.50 49.00 49.50 50.00 50.50 51.00 51.50 52.00 52.50

## ACCUTEST

## GC/MS TOTAL ION CHROMATOGRAM

File: Z:\1\data\S131120\S34120.D

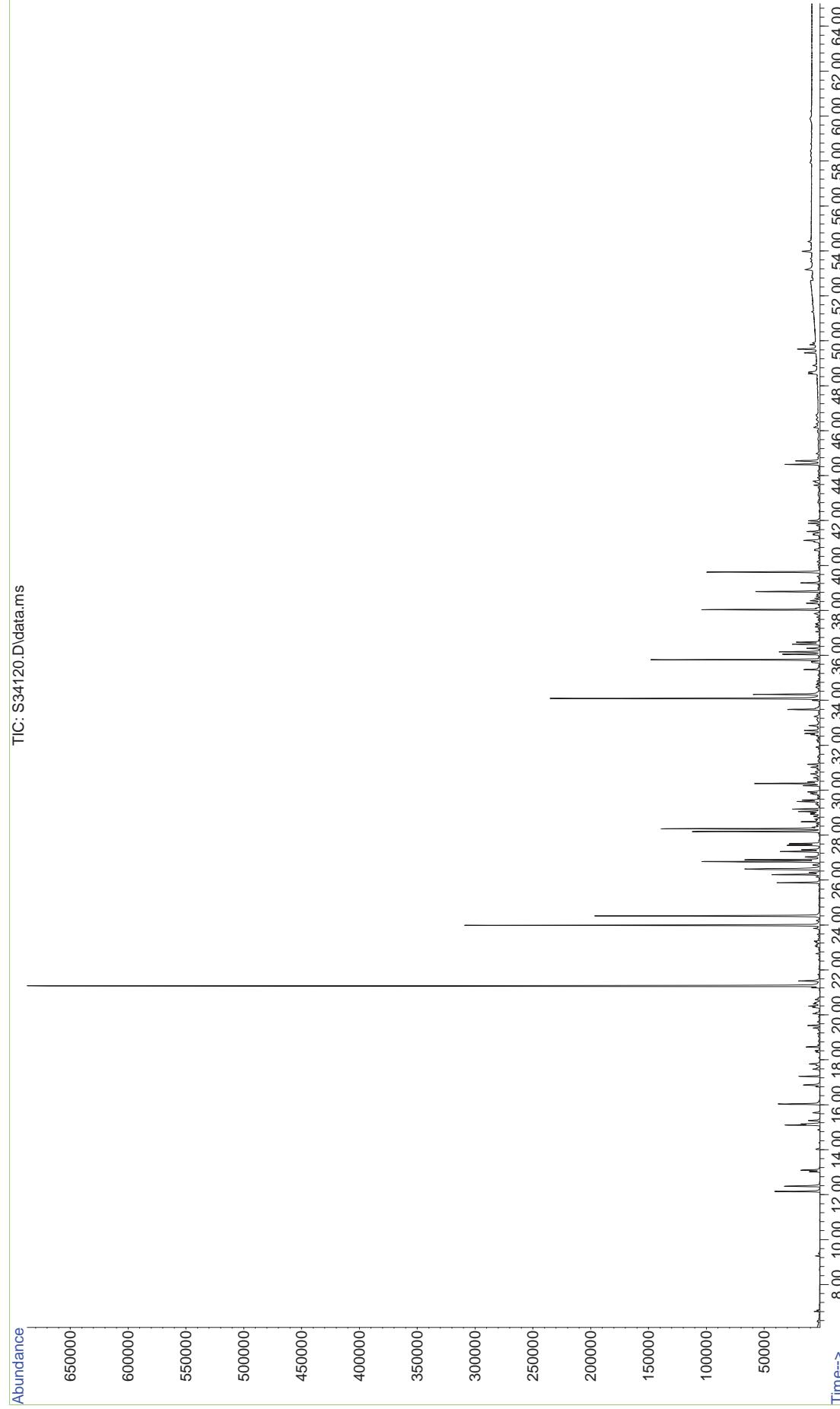
Date Acquired: 20 Nov 2013 5:53 pm

Method File: SAPAHSIM-MTBE.M

Sample Name: OP35834-dup1

Misc Info: OP35834,mss1523,0.0015,,,10,10

Abundance



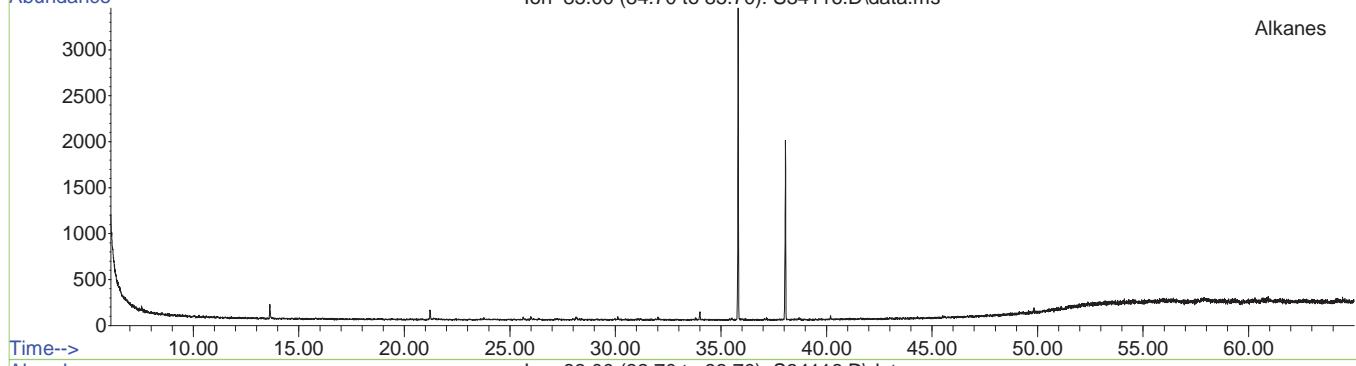
File: Z:\1\data\S131120\S34116.D

Date Acquired: 20 Nov 2013 1:04 pm

Sample Name: op35834-mb

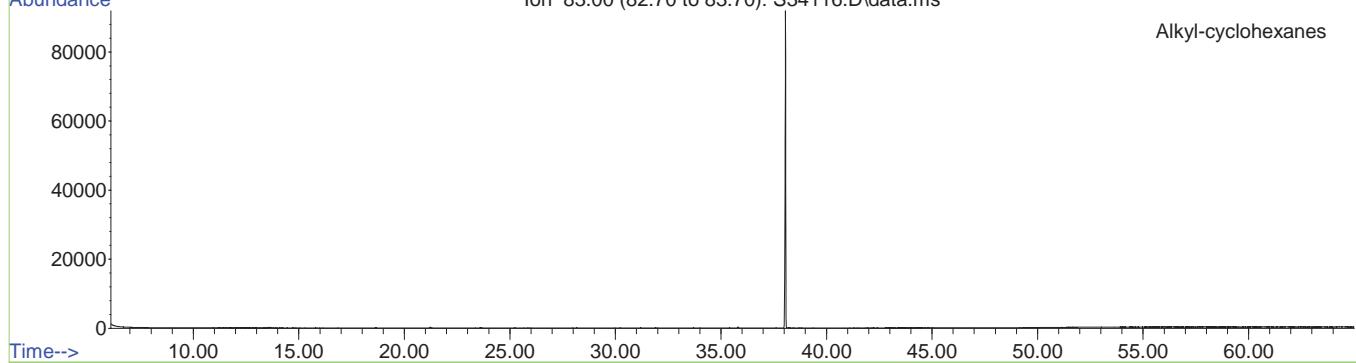
Misc Info: op35834,mss1523,0.01,,,2,1

Abundance Ion 85.00 (84.70 to 85.70): S34116.D\data.ms



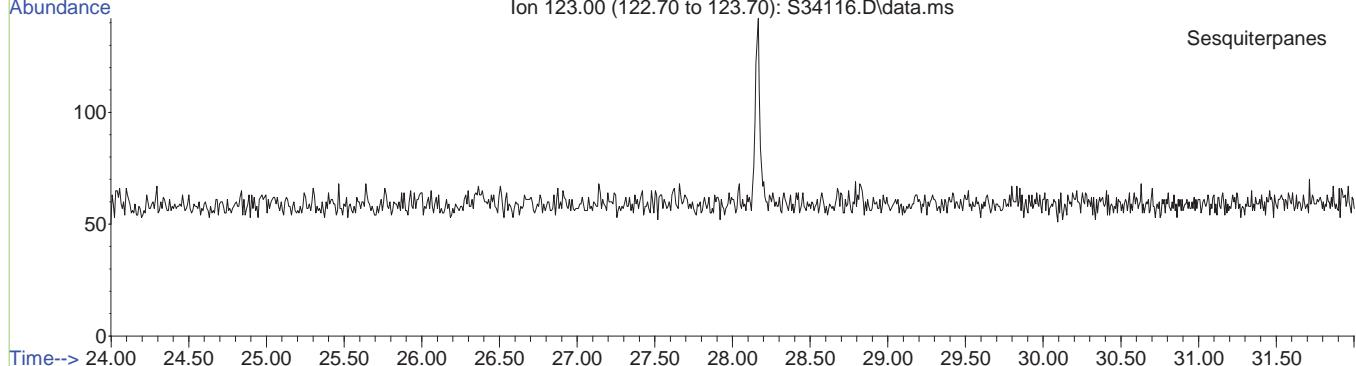
Alkanes

Abundance Ion 83.00 (82.70 to 83.70): S34116.D\data.ms



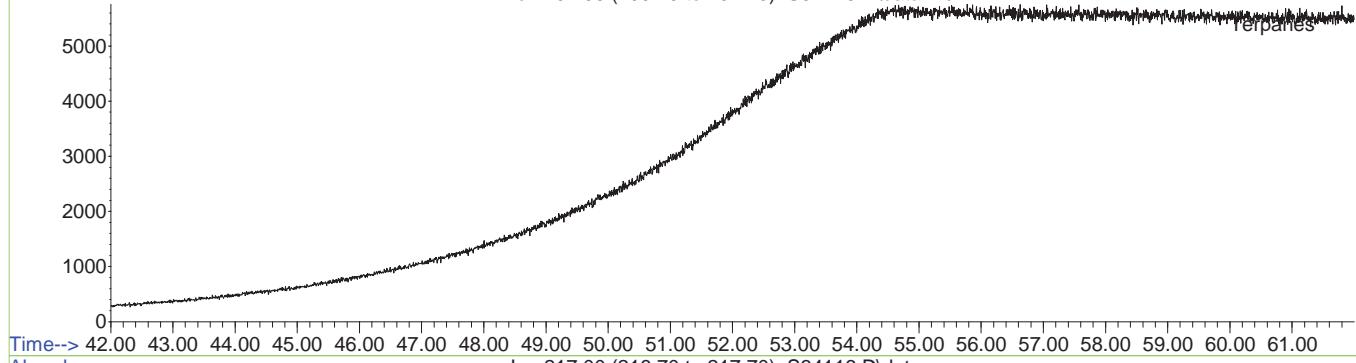
Alkyl-cyclohexanes

Abundance Ion 123.00 (122.70 to 123.70): S34116.D\data.ms



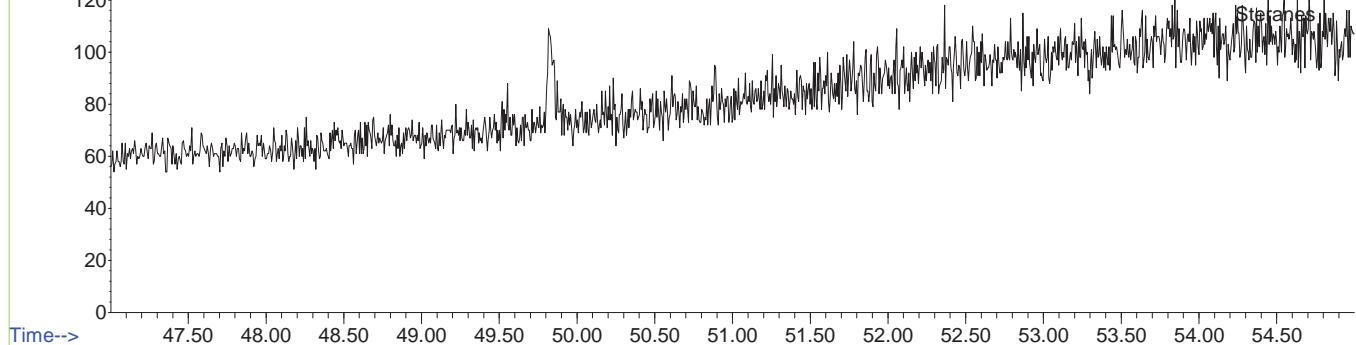
Sesquiterpanes

Abundance Ion 191.00 (190.70 to 191.70): S34116.D\data.ms



Terpanes

Abundance Ion 217.00 (216.70 to 217.70): S34116.D\data.ms



Steranes

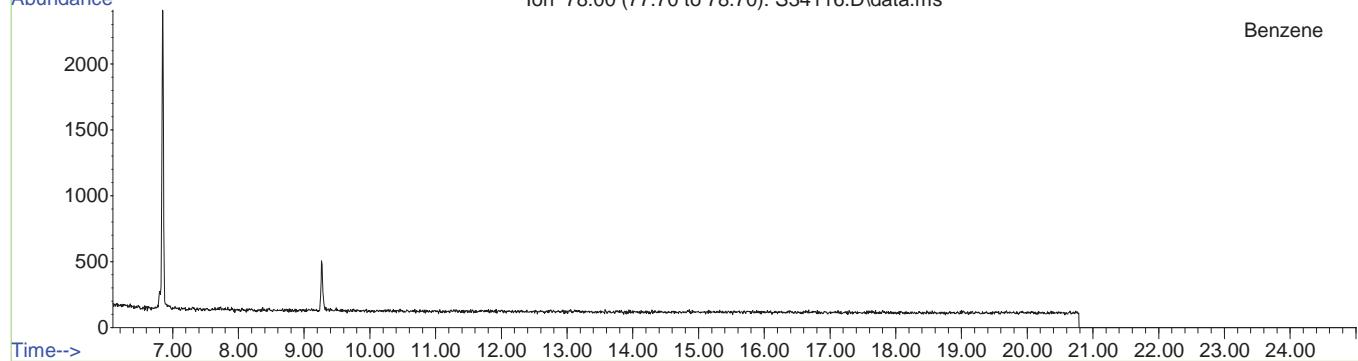
File: Z:\1\data\S131120\S34116.D

Date Acquired: 20 Nov 2013 1:04 pm

Sample Name: op35834-mb

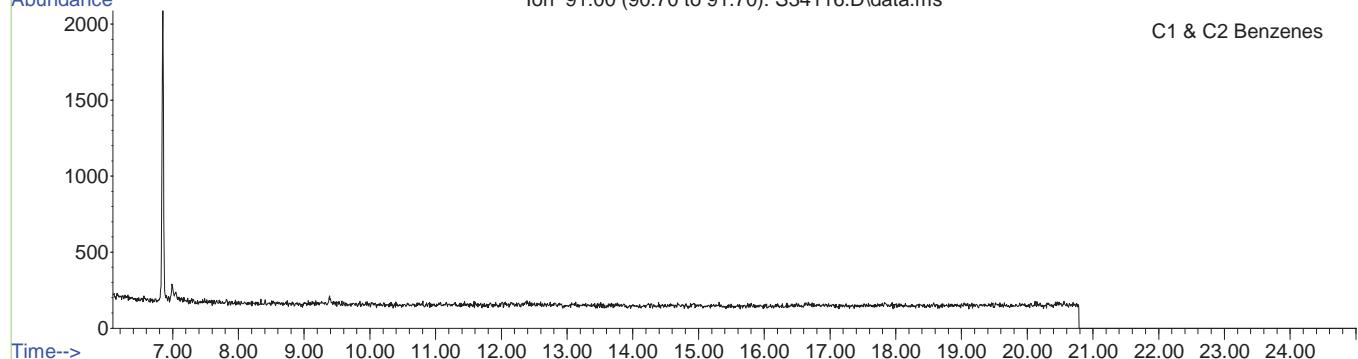
Misc Info: op35834,mss1523,0.01,,,2,,1

Abundance Ion 78.00 (77.70 to 78.70): S34116.D\data.ms



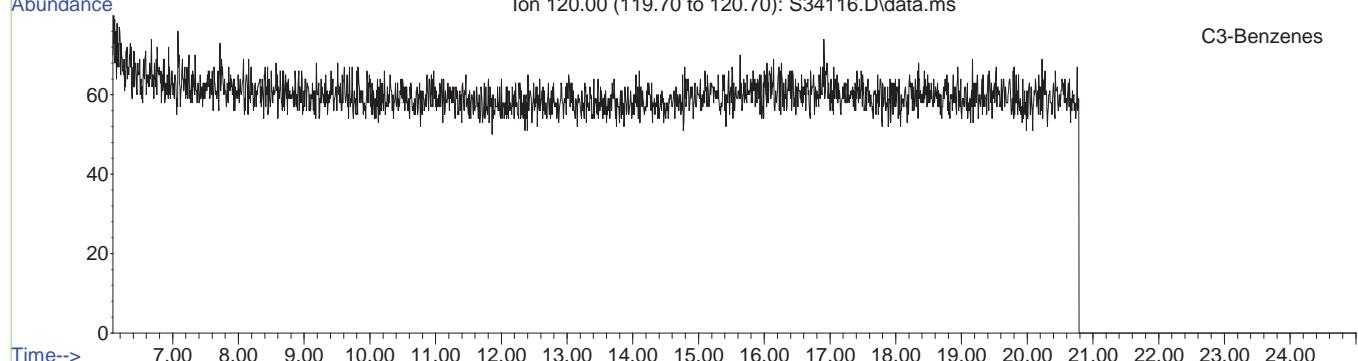
Benzene

Abundance Ion 91.00 (90.70 to 91.70): S34116.D\data.ms



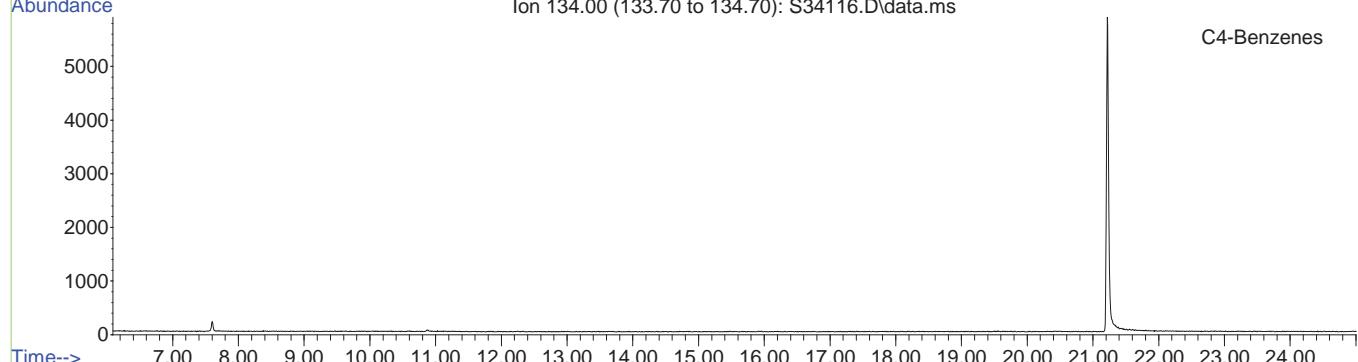
C1 & C2 Benzenes

Abundance Ion 120.00 (119.70 to 120.70): S34116.D\data.ms



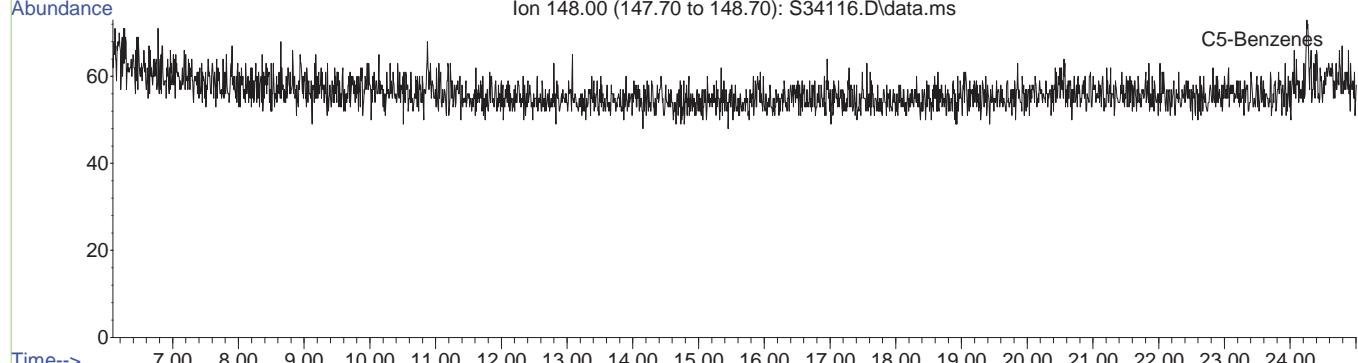
C3-Benzene

Abundance Ion 134.00 (133.70 to 134.70): S34116.D\data.ms



C4-Benzene

Abundance Ion 148.00 (147.70 to 148.70): S34116.D\data.ms



C5-Benzene

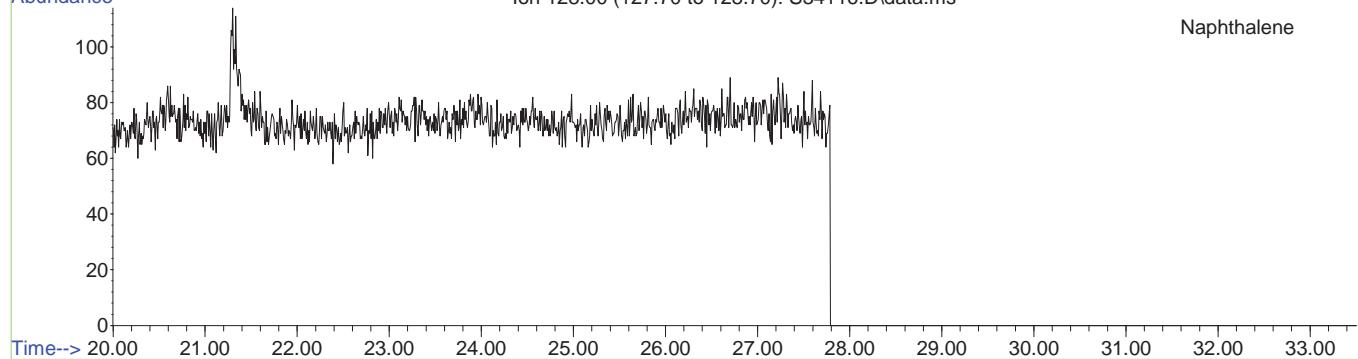
File: Z:\1\data\S131120\S34116.D

Date Acquired: 20 Nov 2013 1:04 pm

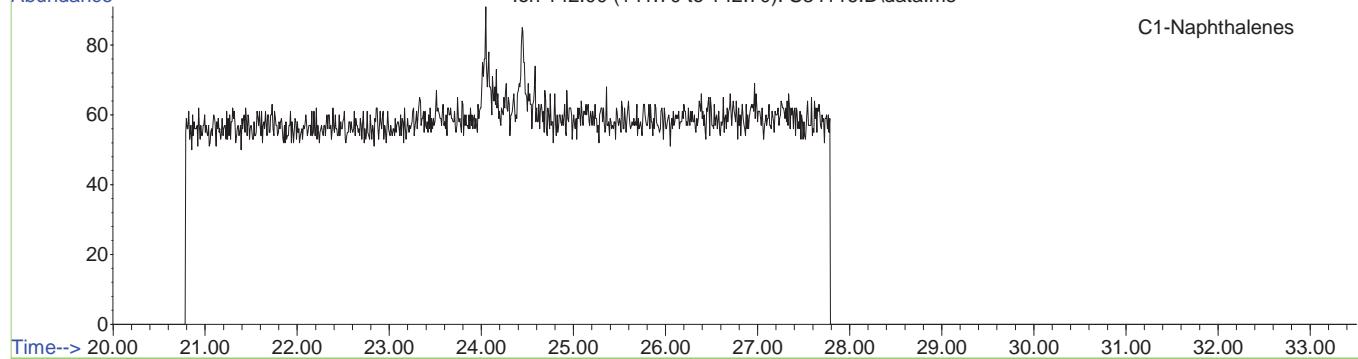
Sample Name: op35834-mb

Misc Info: op35834,mss1523,0.01,,,2,1

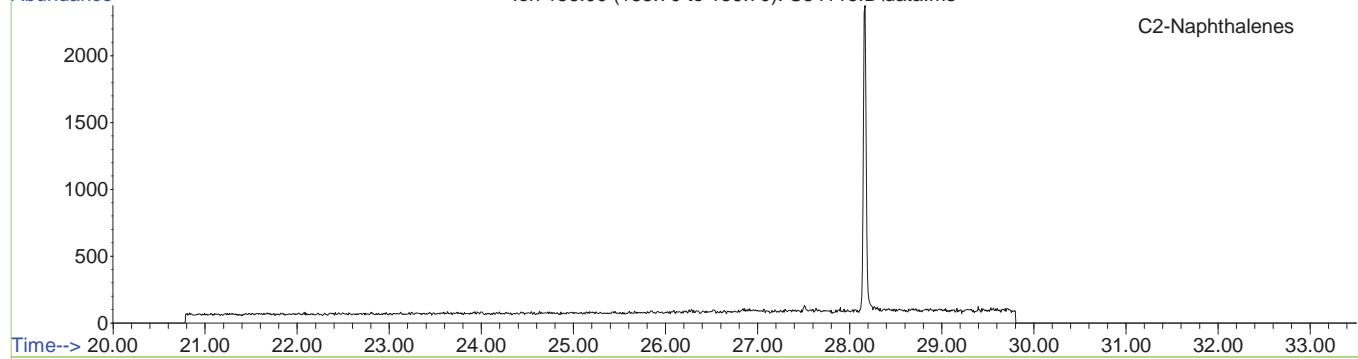
Abundance Ion 128.00 (127.70 to 128.70): S34116.D\data.ms



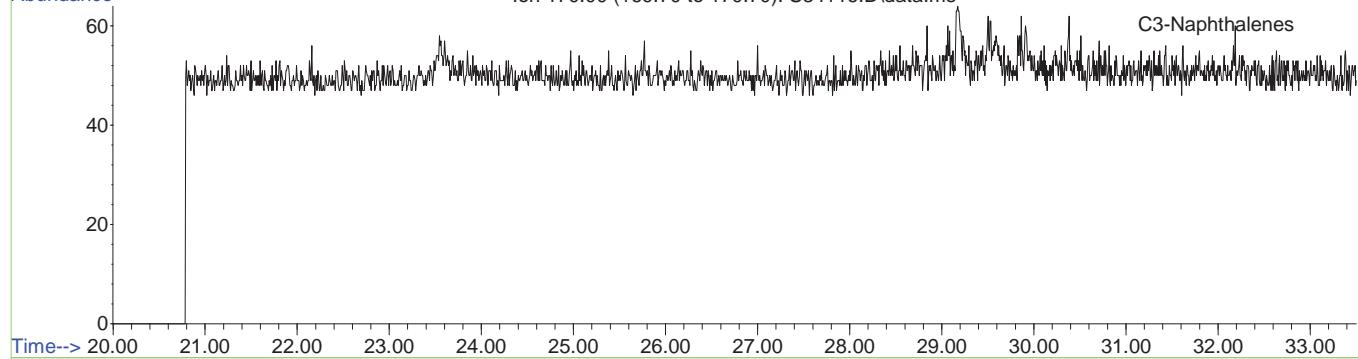
Abundance Ion 142.00 (141.70 to 142.70): S34116.D\data.ms



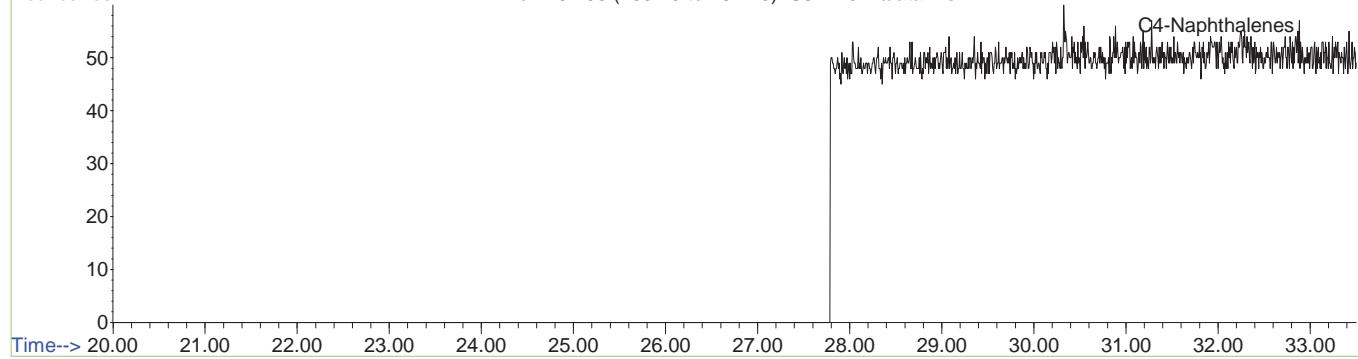
Abundance Ion 156.00 (155.70 to 156.70): S34116.D\data.ms



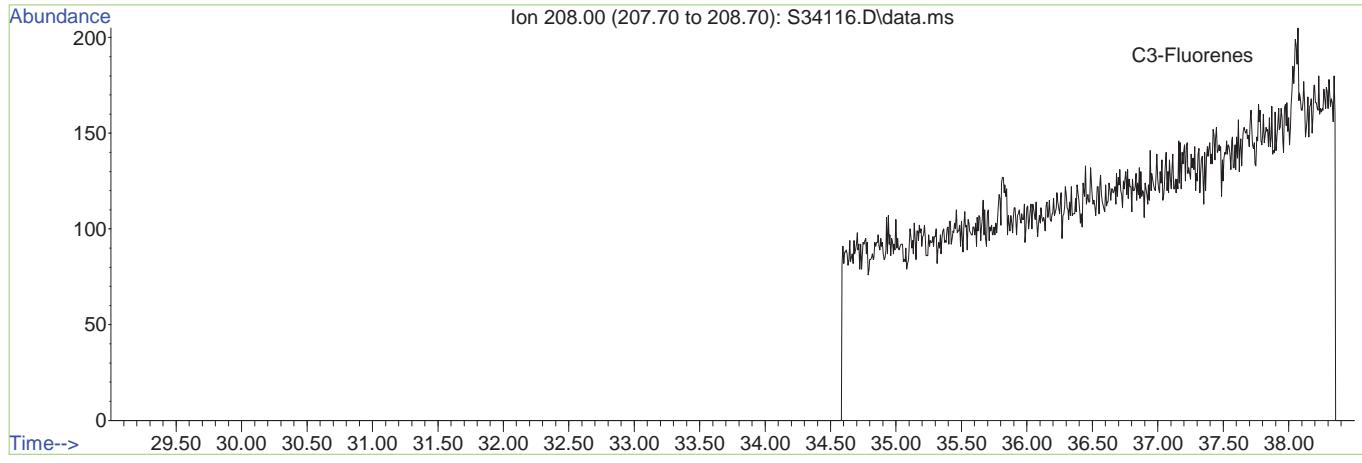
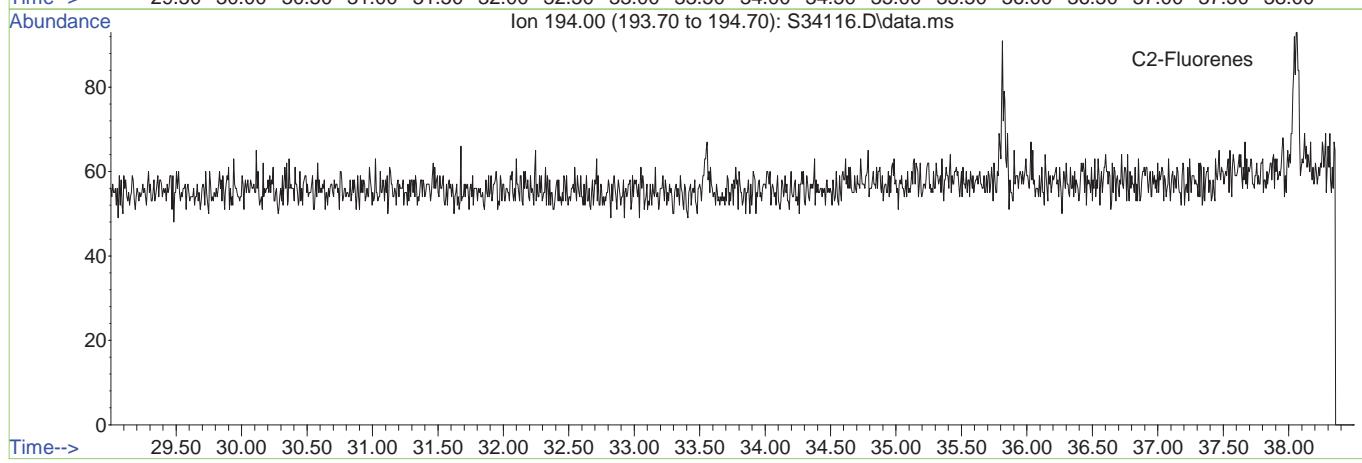
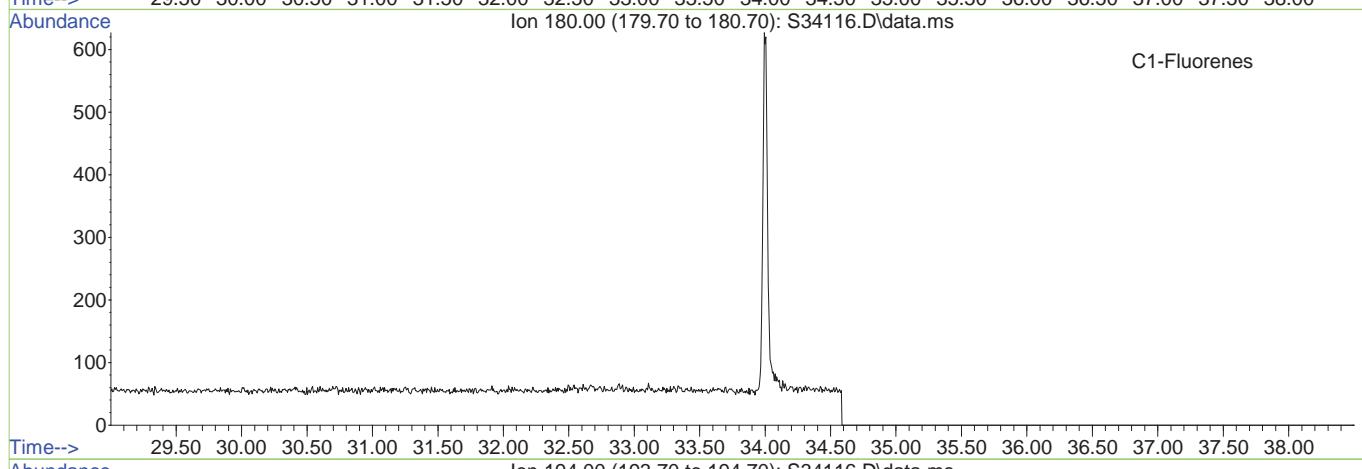
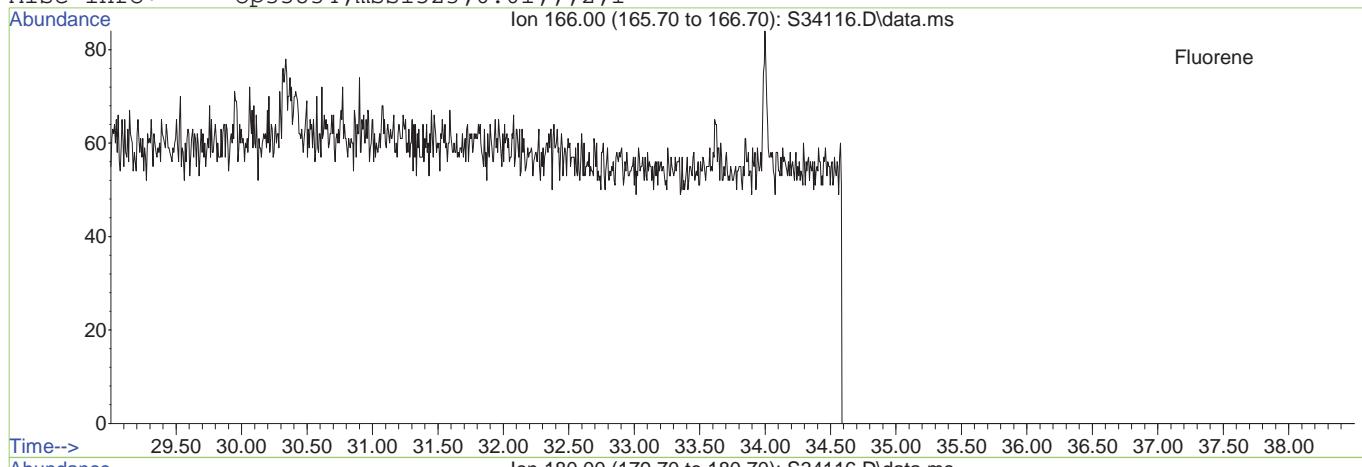
Abundance Ion 170.00 (169.70 to 170.70): S34116.D\data.ms



Abundance Ion 184.00 (183.70 to 184.70): S34116.D\data.ms



File: Z:\1\data\S131120\S34116.D  
Date Acquired: 20 Nov 2013 1:04 pm  
Method File: SAPAHSIM-MTBE.M  
Sample Name: op35834-mb  
Misc Info: op35834,mss1523,0.01,,,2,1



File: Z:\1\data\S131120\S34116.D

Date Acquired: 20 Nov 2013 1:04 pm

Sample Name: op35834-mb

Misc Info: op35834,mss1523,0.01,,,2,,1

Abundance

Ion 178.00 (177.70 to 178.70): S34116.D\data.ms

Phenanthrene & Anthracene

10000

5000

0

Time-->

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00

Abundance

Ion 192.00 (191.70 to 192.70): S34116.D\data.ms

C1-Phenanthrenes & Anthracenes

3000

2500

2000

1500

1000

500

0

Time-->

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00

Abundance

Ion 206.00 (205.70 to 206.70): S34116.D\data.ms

C2-Phenanthrenes & Anthracenes

3000

2500

2000

1500

1000

500

0

Time-->

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00

Abundance

Ion 220.00 (219.70 to 220.70): S34116.D\data.ms

C3-Phenanthrenes & Anthracenes

50

40

30

20

10

0

Time-->

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00

Abundance

Ion 234.00 (233.70 to 234.70): S34116.D\data.ms

C4-Phenanthrenes & Anthracenes

50

40

30

20

10

0

Time-->

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00

File: Z:\1\data\S131120\S34116.D

Date Acquired: 20 Nov 2013 1:04 pm

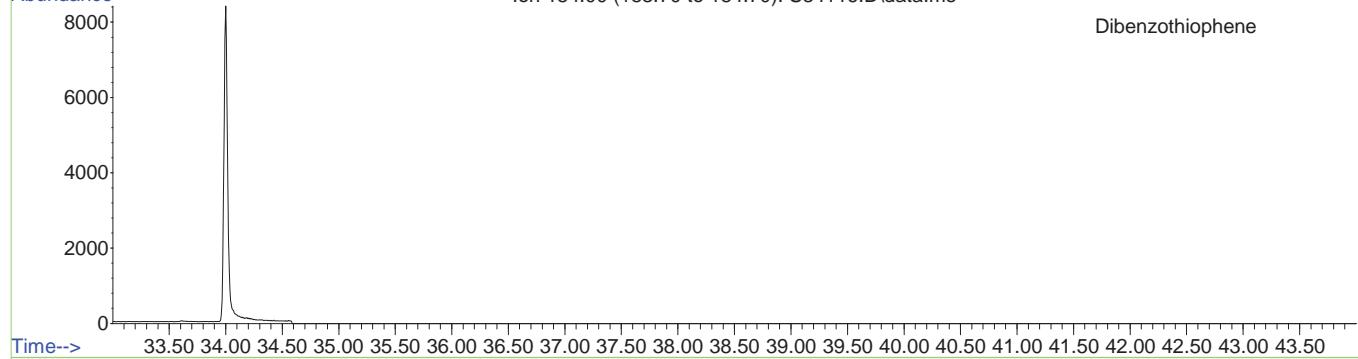
Sample Name: op35834-mb

Misc Info: op35834,mss1523,0.01,,,2,,1

Abundance

Ion 184.00 (183.70 to 184.70): S34116.D\data.ms

Dibenzothiophene



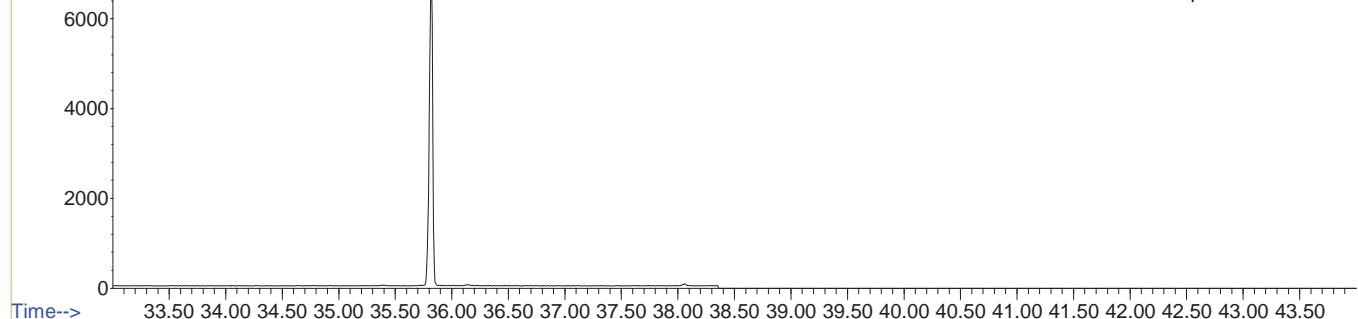
Time-->

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50

Abundance

Ion 198.00 (197.70 to 198.70): S34116.D\data.ms

C1-Dibenzothiophenes



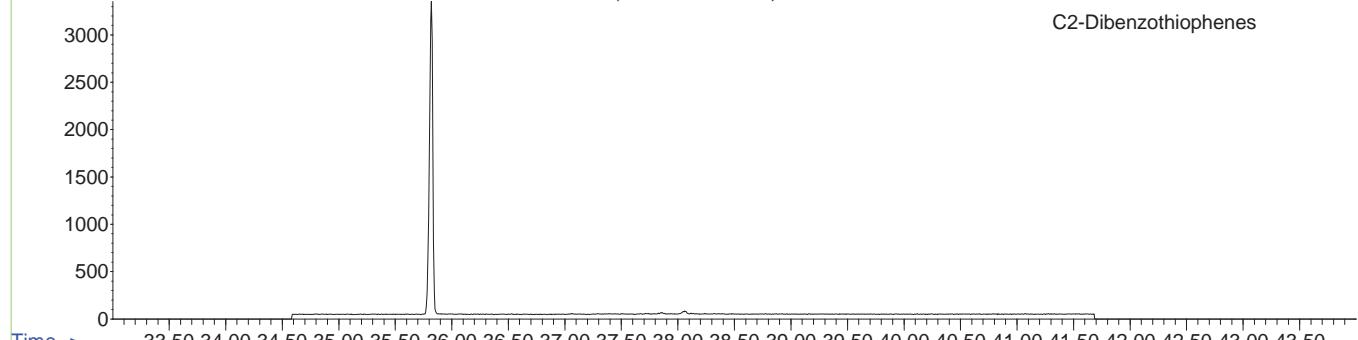
Time-->

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50

Abundance

Ion 212.00 (211.70 to 212.70): S34116.D\data.ms

C2-Dibenzothiophenes



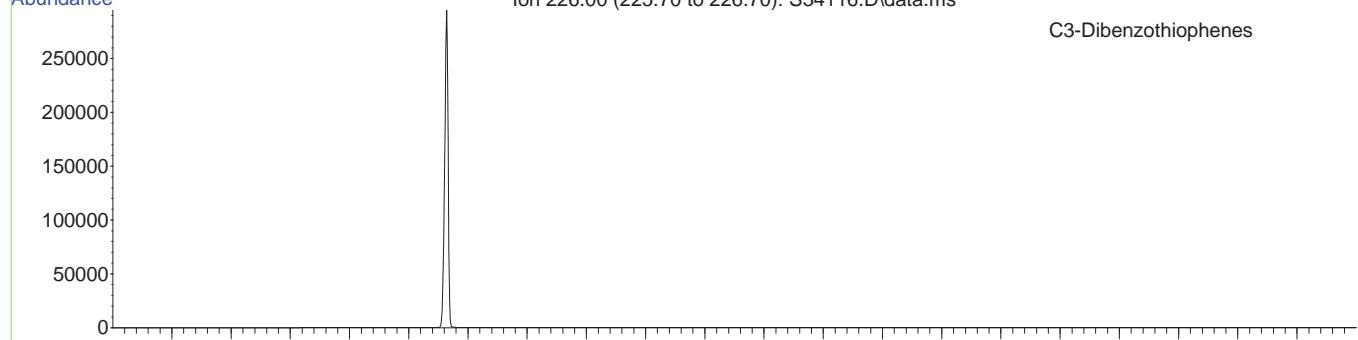
Time-->

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50

Abundance

Ion 226.00 (225.70 to 226.70): S34116.D\data.ms

C3-Dibenzothiophenes



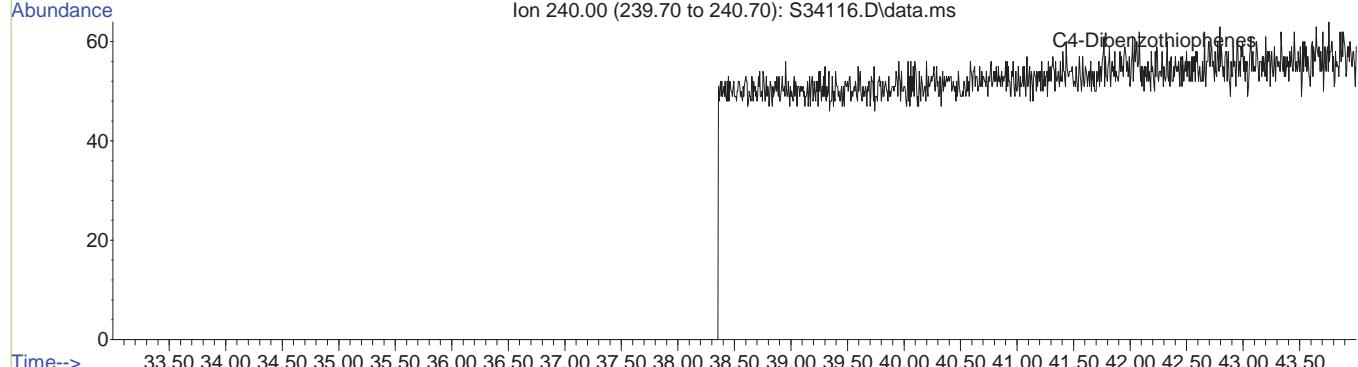
Time-->

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50

Abundance

Ion 240.00 (239.70 to 240.70): S34116.D\data.ms

C4-Dibenzothiophenes



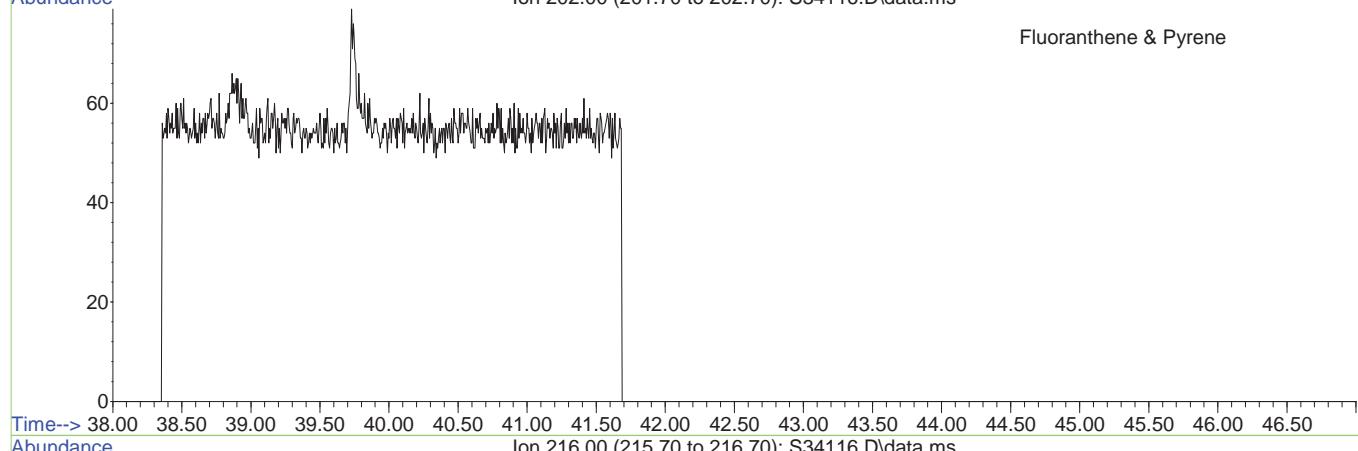
Time-->

33.50 34.00 34.50 35.00 35.50 36.00 36.50 37.00 37.50 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50

File: Z:\1\data\S131120\S34116.D  
Date Acquired: 20 Nov 2013 1:04 pm  
Method File: SAPAHSIM-MTBE.M  
Sample Name: op35834-mb  
Misc Info: op35834,mss1523,0.01,,,2,1

Abundance Ion 202.00 (201.70 to 202.70): S34116.D\data.ms

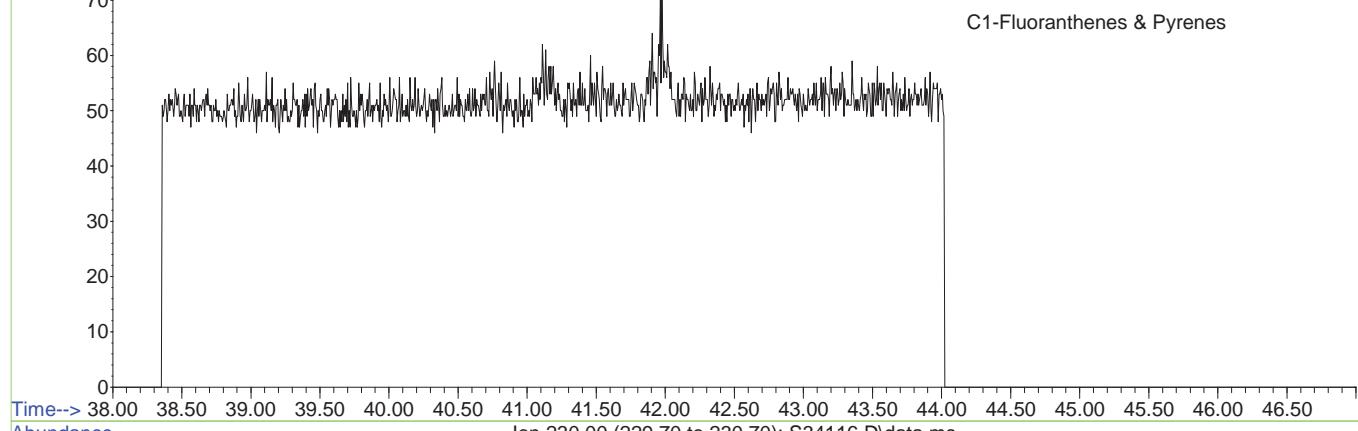
Fluoranthene & Pyrene



Time--> 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50 44.00 44.50 45.00 45.50 46.00 46.50

Abundance Ion 216.00 (215.70 to 216.70): S34116.D\data.ms

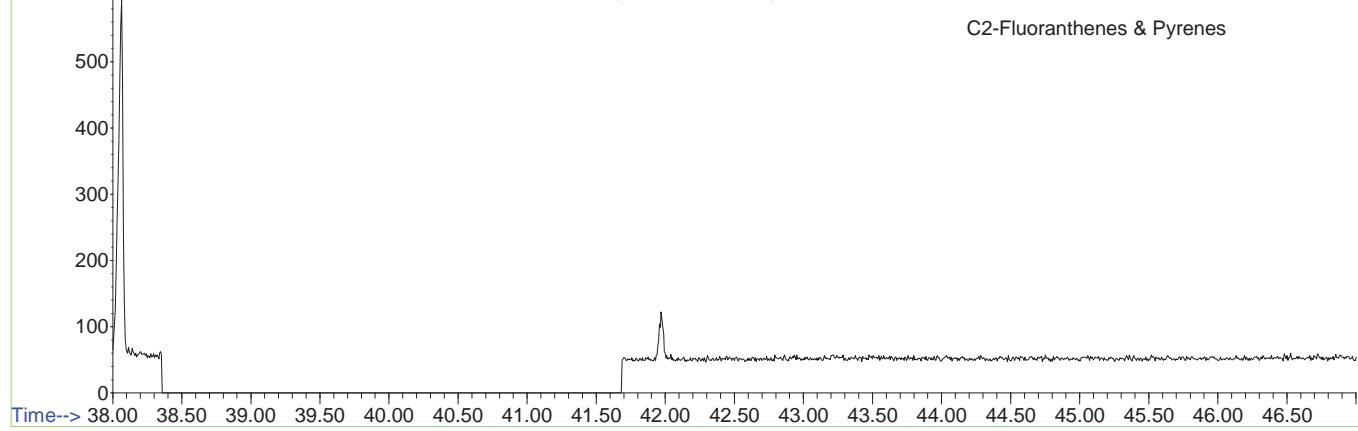
C1-Fluoranthenes & Pyrenes



Time--> 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50 44.00 44.50 45.00 45.50 46.00 46.50

Abundance Ion 230.00 (229.70 to 230.70): S34116.D\data.ms

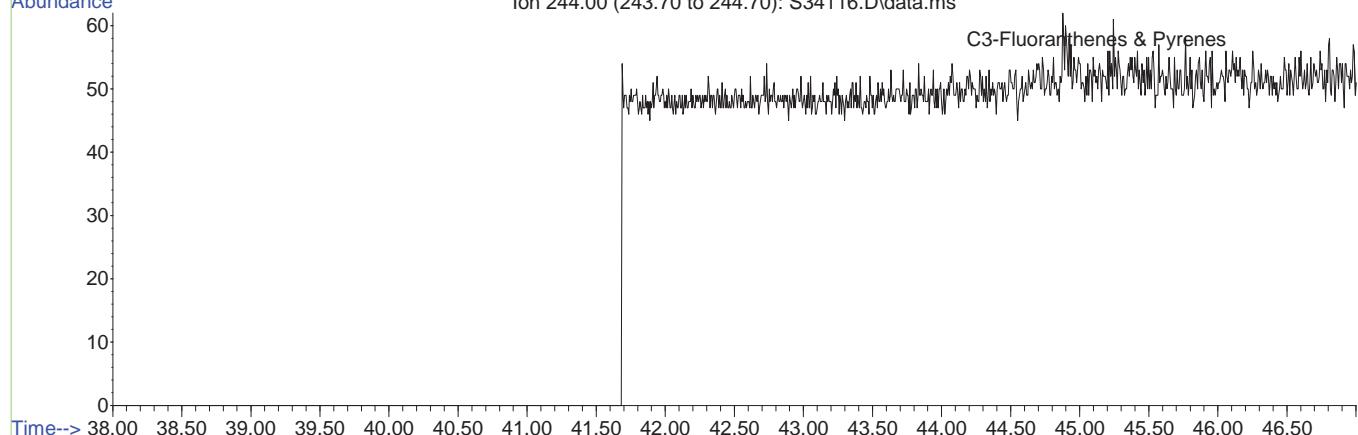
C2-Fluoranthenes & Pyrenes



Time--> 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50 44.00 44.50 45.00 45.50 46.00 46.50

Abundance Ion 244.00 (243.70 to 244.70): S34116.D\data.ms

C3-Fluoranthenes & Pyrenes



Time--> 38.00 38.50 39.00 39.50 40.00 40.50 41.00 41.50 42.00 42.50 43.00 43.50 44.00 44.50 45.00 45.50 46.00 46.50

File: Z:\1\data\S131120\S34116.D

Date Acquired: 20 Nov 2013 1:04 pm

Sample Name: op35834-mb

Misc Info: op35834,mss1523,0.01,,,2,,1

Abundance Ion 228.00 (227.70 to 228.70): S34116.D\data.ms

Benz(a)anthracene & Chrysene

Time--> 44.00 44.50 45.00 45.50 46.00 46.50 47.00 47.50 48.00 48.50 49.00 49.50 50.00 50.50 51.00 51.50 52.00 52.50

Abundance Ion 242.00 (241.70 to 242.70): S34116.D\data.ms

C1-Benz(a)anthracenes & Chrysenes

Time--> 44.00 44.50 45.00 45.50 46.00 46.50 47.00 47.50 48.00 48.50 49.00 49.50 50.00 50.50 51.00 51.50 52.00 52.50

Abundance Ion 256.00 (255.70 to 256.70): S34116.D\data.ms

C2-Benz(a)anthracenes & Chrysenes

Time--> 44.00 44.50 45.00 45.50 46.00 46.50 47.00 47.50 48.00 48.50 49.00 49.50 50.00 50.50 51.00 51.50 52.00 52.50

Abundance Ion 270.00 (269.70 to 270.70): S34116.D\data.ms

C3-Benz(a)anthracenes & Chrysenes

Time--> 44.00 44.50 45.00 45.50 46.00 46.50 47.00 47.50 48.00 48.50 49.00 49.50 50.00 50.50 51.00 51.50 52.00 52.50

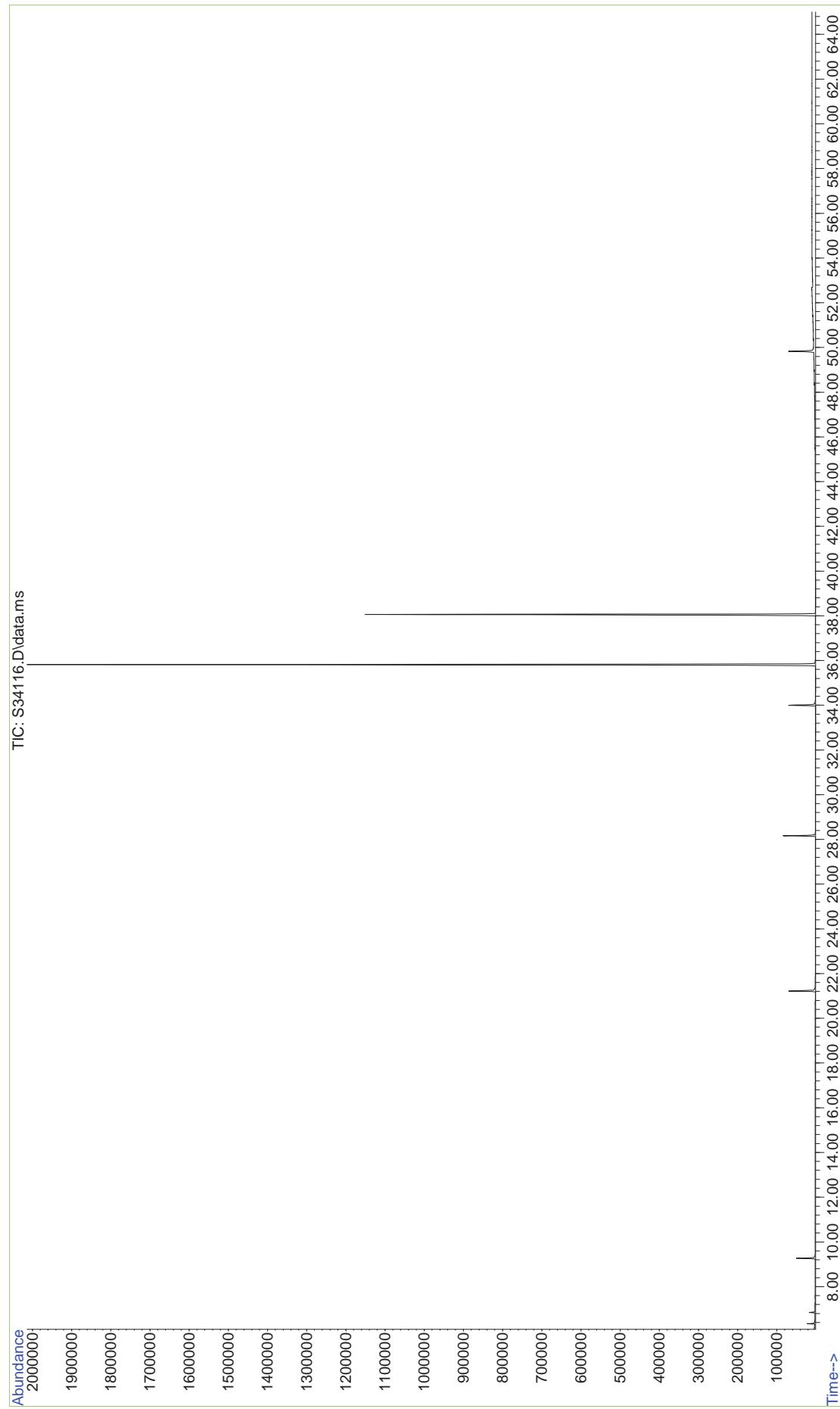
Abundance Ion 284.00 (283.70 to 284.70): S34116.D\data.ms

C4-Benz(a)anthracenes & Chrysenes

## ACCUTEST

## GC/MS TOTAL ION CHROMATOGRAM

File: Z:\1\data\S131120\S34116.D  
Date Acquired: 20 Nov 2013 1:04 pm  
Method File: SAPAHSIM-MTBE.M  
Sample Name: OP35834-mbo  
Misc Info: OP35834,mss1523,0.01,,,2,1



## **Appendix E**

# **Accutest Laboratory Report**



11/26/13



## Technical Report for

META Environmental, Inc.

Parsons, Halleck Street, NJ

Accutest Job Number: MC26103

Sampling Date: 10/31/13

### Report to:

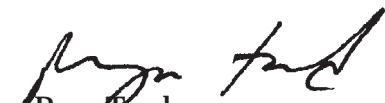
META Environmental, Inc.  
115 Dean Avenue Suite 300  
Franklin MA 02038, MA 02038  
dmauro@metaenv.com

ATTN: Dave Mauro

Total number of pages in report: 89



Test results contained within this data package meet the requirements of the National Environmental Laboratory Accreditation Program and/or state specific certification programs as applicable.



Reza Pand  
Lab Director

Client Service contact: Matthew Morrell 508-481-6200

Certifications: MA (M-MA136, SW846 NELAC) CT (PH-0109) NH (250210) RI (00071) ME (MA00136) FL (E87579)  
NY (11791) NJ (MA926) PA (6801121) ND (R-188) CO MN (11546AA) NC (653) IL (002337) WI (399080220)  
DoD ELAP (L-A-B L2235)

This report shall not be reproduced, except in its entirety, without the written approval of Accutest Laboratories.  
Test results relate only to samples analyzed.

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<table style="width: 100%; border-collapse: collapse;"> <tr><td><b>Section 1: Sample Summary .....</b></td><td style="text-align: right;"><b>3</b></td></tr> <tr><td><b>Section 2: Case Narrative/Conformance Summary .....</b></td><td style="text-align: right;"><b>4</b></td></tr> <tr><td><b>Section 3: Summary of Hits .....</b></td><td style="text-align: right;"><b>5</b></td></tr> <tr><td><b>Section 4: Sample Results .....</b></td><td style="text-align: right;"><b>7</b></td></tr> <tr><td>    <b>4.1: MC26103-1: SB-4(15-17)NAPL .....</b></td><td style="text-align: right;"><b>8</b></td></tr> <tr><td><b>Section 5: Misc. Forms .....</b></td><td style="text-align: right;"><b>12</b></td></tr> <tr><td>    <b>5.1: Chain of Custody .....</b></td><td style="text-align: right;"><b>13</b></td></tr> <tr><td><b>Section 6: GC/MS Semi-volatiles - QC Data Summaries .....</b></td><td style="text-align: right;"><b>15</b></td></tr> <tr><td>    <b>6.1: Method Blank Summary .....</b></td><td style="text-align: right;"><b>16</b></td></tr> <tr><td>    <b>6.2: Blank Spike Summary .....</b></td><td style="text-align: right;"><b>19</b></td></tr> <tr><td>    <b>6.3: Duplicate Summary .....</b></td><td style="text-align: right;"><b>22</b></td></tr> <tr><td>    <b>6.4: Surrogate Recovery Summaries .....</b></td><td style="text-align: right;"><b>25</b></td></tr> <tr><td><b>Section 7: GC/MS Semi-volatiles - Raw Data .....</b></td><td style="text-align: right;"><b>26</b></td></tr> <tr><td>    <b>7.1: Samples .....</b></td><td style="text-align: right;"><b>27</b></td></tr> <tr><td>    <b>7.2: Method Blanks .....</b></td><td style="text-align: right;"><b>74</b></td></tr> <tr><td><b>Section 8: GC Semi-volatiles - QC Data Summaries .....</b></td><td style="text-align: right;"><b>78</b></td></tr> <tr><td>    <b>8.1: Method Blank Summary .....</b></td><td style="text-align: right;"><b>79</b></td></tr> <tr><td>    <b>8.2: Blank Spike Summary .....</b></td><td style="text-align: right;"><b>80</b></td></tr> <tr><td>    <b>8.3: Duplicate Summary .....</b></td><td style="text-align: right;"><b>81</b></td></tr> <tr><td>    <b>8.4: Surrogate Recovery Summaries .....</b></td><td style="text-align: right;"><b>82</b></td></tr> <tr><td><b>Section 9: GC Semi-volatiles - Raw Data .....</b></td><td style="text-align: right;"><b>83</b></td></tr> <tr><td>    <b>9.1: Samples .....</b></td><td style="text-align: right;"><b>84</b></td></tr> <tr><td>    <b>9.2: Method Blanks .....</b></td><td style="text-align: right;"><b>87</b></td></tr> </table>	<b>Section 1: Sample Summary .....</b>	<b>3</b>	<b>Section 2: Case Narrative/Conformance Summary .....</b>	<b>4</b>	<b>Section 3: Summary of Hits .....</b>	<b>5</b>	<b>Section 4: Sample Results .....</b>	<b>7</b>	<b>4.1: MC26103-1: SB-4(15-17)NAPL .....</b>	<b>8</b>	<b>Section 5: Misc. Forms .....</b>	<b>12</b>	<b>5.1: Chain of Custody .....</b>	<b>13</b>	<b>Section 6: GC/MS Semi-volatiles - QC Data Summaries .....</b>	<b>15</b>	<b>6.1: Method Blank Summary .....</b>	<b>16</b>	<b>6.2: Blank Spike Summary .....</b>	<b>19</b>	<b>6.3: Duplicate Summary .....</b>	<b>22</b>	<b>6.4: Surrogate Recovery Summaries .....</b>	<b>25</b>	<b>Section 7: GC/MS Semi-volatiles - Raw Data .....</b>	<b>26</b>	<b>7.1: Samples .....</b>	<b>27</b>	<b>7.2: Method Blanks .....</b>	<b>74</b>	<b>Section 8: GC Semi-volatiles - QC Data Summaries .....</b>	<b>78</b>	<b>8.1: Method Blank Summary .....</b>	<b>79</b>	<b>8.2: Blank Spike Summary .....</b>	<b>80</b>	<b>8.3: Duplicate Summary .....</b>	<b>81</b>	<b>8.4: Surrogate Recovery Summaries .....</b>	<b>82</b>	<b>Section 9: GC Semi-volatiles - Raw Data .....</b>	<b>83</b>	<b>9.1: Samples .....</b>	<b>84</b>	<b>9.2: Method Blanks .....</b>	<b>87</b>	<table style="width: 100%; border-collapse: collapse;"> <tr><td style="background-color: #002060; color: white; text-align: center; padding: 2px;">1</td></tr> <tr><td style="background-color: #006400; color: white; text-align: center; padding: 2px;">2</td></tr> <tr><td style="background-color: #C8A234; color: black; text-align: center; padding: 2px;">3</td></tr> <tr><td style="background-color: #C00000; color: white; text-align: center; padding: 2px;">4</td></tr> <tr><td style="background-color: #6B8E23; color: black; text-align: center; padding: 2px;">5</td></tr> <tr><td style="background-color: #A9A9A9; color: black; text-align: center; padding: 2px;">6</td></tr> <tr><td style="background-color: #8B4513; color: white; text-align: center; padding: 2px;">7</td></tr> <tr><td style="background-color: #008000; color: black; text-align: center; padding: 2px;">8</td></tr> <tr><td style="background-color: #FF0000; color: white; text-align: center; padding: 2px;">9</td></tr> </table>	1	2	3	4	5	6	7	8	9
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Accutest Laboratories

## Sample Summary

META Environmental, Inc.

Parsons, Halleck Street, NJ

Job No: MC26103

Sample Number	Collected Date	Time By	Matrix Received	Code Type	Client Sample ID
MC26103-1	10/31/13	14:15 ZL	11/08/13	SO Oil	SB-4(15-17)NAPL

---

Soil samples reported on a dry weight basis unless otherwise indicated on result page.



## SAMPLE DELIVERY GROUP CASE NARRATIVE

**Client:** META Environmental, Inc.

**Job No** MC26103

**Site:** Parsons, Halleck Street, NJ

**Report Date** 11/26/2013 11:48:42 AM

1 Sample(s), 0 Trip Blank(s) and 0 Field Blank(s) were collected on 10/31/2013 and were received at Accutest on 11/08/2013 properly preserved, at 0.5 Deg. C and intact. These Samples received an Accutest job number of MC26103. A listing of the Laboratory Sample ID, Client Sample ID and dates of collection are presented in the Results Summary Section of this report.

Except as noted below, all method specified calibrations and quality control performance criteria were met for this job. For more information, please refer to QC summary pages.

### Extractables by GCMS By Method D5739-06/8270C SIM

<b>Matrix:</b> SO	<b>Batch ID:</b> OP35834
-------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) MC26103-1DUP were used as the QC samples indicated.
- RPD(s) for Duplicate for the majority of the compounds are outside control limits for sample OP35834-DUP1. High RPD due to possible sample heterogeneity.
- MC26103-1: Sample extracted beyond the recommended holding time for soils.

### Extractables by GC By Method ASTM D3328-06

<b>Matrix:</b> SO	<b>Batch ID:</b> OP35833
-------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) MC26103-1DUP were used as the QC samples indicated.
- MC26103-1: Sample extracted beyond the recommended holding time for soils.
- OP35833-DUP1, MC26103-1 for o-Terphenyl: Outside control limits due to dilution.

The Accutest Laboratories of New England certifies that all analysis were performed within method specification. It is further recommended that this report to be used in its entirety. The Accutest Laboratories of NE, Laboratory Director or assignee as verified by the signature on the cover page has authorized the release of this report(MC26103).

**Summary of Hits**

**Job Number:** MC26103  
**Account:** META Environmental, Inc.  
**Project:** Parsons, Halleck Street, NJ  
**Collected:** 10/31/13



Lab Sample ID	Client Sample ID	Result/ Analyte	Qual	RL	MDL	Units	Method
MC26103-1	SB-4(15-17)NAPL						
Benzene a	10900000	910000		ug/kg	D5739-06/8270C SIM		
C1-Benzene a	1580000	910000		ug/kg	D5739-06/8270C SIM		
C2-Benzenes a	72300000	910000		ug/kg	D5739-06/8270C SIM		
C3-Benzenes a	82600000	910000		ug/kg	D5739-06/8270C SIM		
C4-Benzenes a	43300000	910000		ug/kg	D5739-06/8270C SIM		
C5-Benzenes a	12800000	910000		ug/kg	D5739-06/8270C SIM		
Toluene a	2190000	910000		ug/kg	D5739-06/8270C SIM		
Ethylbenzene a	55300000	910000		ug/kg	D5739-06/8270C SIM		
m,p-Xylene a	52800000	910000		ug/kg	D5739-06/8270C SIM		
Styrene a	11300000	910000		ug/kg	D5739-06/8270C SIM		
o-Xylene a	23800000	910000		ug/kg	D5739-06/8270C SIM		
Isopropylbenzene a	6660000	910000		ug/kg	D5739-06/8270C SIM		
n-Propylbenzene a	2440000	910000		ug/kg	D5739-06/8270C SIM		
1,3,5-Trimethylbenzene a	13200000	910000		ug/kg	D5739-06/8270C SIM		
1,2,3-Trimethylbenzene a	13500000	910000		ug/kg	D5739-06/8270C SIM		
1,2,4-Trimethylbenzene a	40900000	910000		ug/kg	D5739-06/8270C SIM		
t-Butylbenzene a	5370000	910000		ug/kg	D5739-06/8270C SIM		
p-Isopropyltoluene a	6740000	910000		ug/kg	D5739-06/8270C SIM		
n-Butylbenzene a	1480000	910000		ug/kg	D5739-06/8270C SIM		
Benzo(b)thiophene a	26600000	910000		ug/kg	D5739-06/8270C SIM		
Naphthalene a	744000000	910000		ug/kg	D5739-06/8270C SIM		
2-Methylnaphthalene a	310000000	910000		ug/kg	D5739-06/8270C SIM		
1-Methylnaphthalene a	195000000	910000		ug/kg	D5739-06/8270C SIM		
C1-Naphthalenes a	331000000	910000		ug/kg	D5739-06/8270C SIM		
C2-Naphthalenes a	155000000	910000		ug/kg	D5739-06/8270C SIM		
C3-Naphthalenes a	51800000	910000		ug/kg	D5739-06/8270C SIM		
C4-Naphthalenes a	14800000	910000		ug/kg	D5739-06/8270C SIM		
Biphenyl a	27300000	910000		ug/kg	D5739-06/8270C SIM		
Acenaphthylene a	25200000	910000		ug/kg	D5739-06/8270C SIM		
Acenaphthene a	102000000	910000		ug/kg	D5739-06/8270C SIM		
Dibenzofuran a	9390000	910000		ug/kg	D5739-06/8270C SIM		
Fluorene a	48300000	910000		ug/kg	D5739-06/8270C SIM		
C1-Fluorennes a	23600000	910000		ug/kg	D5739-06/8270C SIM		
C2-Fluorennes a	13900000	910000		ug/kg	D5739-06/8270C SIM		
C3-Fluorennes a	6710000	910000		ug/kg	D5739-06/8270C SIM		
Dibenzothiophene a	28500000	910000		ug/kg	D5739-06/8270C SIM		
C1-Dibenzothiophenes a	26300000	910000		ug/kg	D5739-06/8270C SIM		
C2-Dibenzothiophenes a	16100000	910000		ug/kg	D5739-06/8270C SIM		
C3-Dibenzothiophenes a	6760000	910000		ug/kg	D5739-06/8270C SIM		
C4-Dibenzothiophenes a	1670000	910000		ug/kg	D5739-06/8270C SIM		
Phenanthrene a	179000000	910000		ug/kg	D5739-06/8270C SIM		
Anthracene a	53200000	910000		ug/kg	D5739-06/8270C SIM		
C1-Phenanthrenes/Anthracenes a	90900000	910000		ug/kg	D5739-06/8270C SIM		

**Summary of Hits**

**Job Number:** MC26103  
**Account:** META Environmental, Inc.  
**Project:** Parsons, Halleck Street, NJ  
**Collected:** 10/31/13



Lab Sample ID Analyte	Client Sample ID	Result/ Qual	RL	MDL	Units	Method
C2-Phenanthrenes/Anthracenes a	33900000	910000			ug/kg	D5739-06/8270C SIM
C3-Phenanthrenes/Anthracenes a	9630000	910000			ug/kg	D5739-06/8270C SIM
C4-Phenanthrenes/Anthracenes a	2610000	910000			ug/kg	D5739-06/8270C SIM
Retene a	1020000	910000			ug/kg	D5739-06/8270C SIM
Benzo(b)naphtho(2,1-d)thiophene a	4680000	910000			ug/kg	D5739-06/8270C SIM
Fluoranthene a	53200000	910000			ug/kg	D5739-06/8270C SIM
Pyrene a	87800000	910000			ug/kg	D5739-06/8270C SIM
C1-Fluoranthenes/Pyrenes a	45700000	910000			ug/kg	D5739-06/8270C SIM
C2-Fluoranthenes/Pyrenes a	11500000	910000			ug/kg	D5739-06/8270C SIM
C3-Fluoranthenes/Pyrenes a	2630000	910000			ug/kg	D5739-06/8270C SIM
Benzo(b)fluorene a	3540000	910000			ug/kg	D5739-06/8270C SIM
Benzo(c)fluorene a	2140000	910000			ug/kg	D5739-06/8270C SIM
2-Methylpyrene a	6480000	910000			ug/kg	D5739-06/8270C SIM
4-Methylpyrene a	5760000	910000			ug/kg	D5739-06/8270C SIM
1-Methylpyrene a	5610000	910000			ug/kg	D5739-06/8270C SIM
Benzo(a)anthracene a	22200000	910000			ug/kg	D5739-06/8270C SIM
Chrysene a	22000000	910000			ug/kg	D5739-06/8270C SIM
C1-Benzo(a)anthracenes/Chrysenes a	10300000	910000			ug/kg	D5739-06/8270C SIM
C2-Benzo(a)anthracenes/Chrysenes a	4160000	910000			ug/kg	D5739-06/8270C SIM
Benzo(b)fluoranthene a	8440000	910000			ug/kg	D5739-06/8270C SIM
Benzo(k)fluoranthene a	12200000	910000			ug/kg	D5739-06/8270C SIM
Benzo(e)pyrene a	11700000	910000			ug/kg	D5739-06/8270C SIM
Benzo(a)pyrene a	21800000	910000			ug/kg	D5739-06/8270C SIM
Perylene a	2850000	910000			ug/kg	D5739-06/8270C SIM
Indeno(1,2,3-cd)pyrene a	7000000	910000			ug/kg	D5739-06/8270C SIM
Dibenzo(a,h)anthracene a	2040000	910000			ug/kg	D5739-06/8270C SIM
Benzo(g,h,i)perylene a	9150000	910000			ug/kg	D5739-06/8270C SIM
Coronene a	2100000	910000			ug/kg	D5739-06/8270C SIM
TPH (C8-C40) a	2510000	450000			mg/kg	ASTM D3328-06

(a) Sample extracted beyond the recommended holding time for soils.

**4**

## Sample Results

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### Report of Analysis

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Accutest Laboratories

**Report of Analysis**

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<b>Client Sample ID:</b>	SB-4(15-17)NAPL	<b>Date Sampled:</b>	10/31/13
<b>Lab Sample ID:</b>	MC26103-1	<b>Date Received:</b>	11/08/13
<b>Matrix:</b>	SO - Oil	<b>Percent Solids:</b>	n/a
<b>Method:</b>	D5739-06/8270C SIM SW846 3580A		
<b>Project:</b>	Parsons, Halleck Street, NJ		

Run #1 <sup>a</sup>	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	S34121.D	10	11/20/13	RP	11/19/13	OP35834	MSS1523
Run #2							

	Initial Weight	Final Volume
Run #1	0.00110 g	10.0 ml
Run #2		

**Alkylated PAHs**

CAS No.	Compound	Result	RL	Units	Q
71-43-2	Benzene	10900000	910000	ug/kg	
	C1-Benzene	1580000	910000	ug/kg	
	C2-Benzenes	72300000	910000	ug/kg	
	C3-Benzenes	82600000	910000	ug/kg	
	C4-Benzenes	43300000	910000	ug/kg	
	C5-Benzenes	12800000	910000	ug/kg	
108-88-3	Toluene	2190000	910000	ug/kg	
100-41-4	Ethylbenzene	55300000	910000	ug/kg	
	m,p-Xylene	52800000	910000	ug/kg	
100-42-5	Styrene	11300000	910000	ug/kg	
95-47-6	o-Xylene	23800000	910000	ug/kg	
98-82-8	Isopropylbenzene	6660000	910000	ug/kg	
103-65-1	n-Propylbenzene	2440000	910000	ug/kg	
108-67-8	1,3,5-Trimethylbenzene	13200000	910000	ug/kg	
526-73-8	1,2,3-Trimethylbenzene	13500000	910000	ug/kg	
95-63-6	1,2,4-Trimethylbenzene	40900000	910000	ug/kg	
98-06-6	t-Butylbenzene	5370000	910000	ug/kg	
135-98-8	sec-Butylbenzene	ND	910000	ug/kg	
99-87-6	p-Isopropyltoluene	6740000	910000	ug/kg	
104-51-8	n-Butylbenzene	1480000	910000	ug/kg	
493-02-7	trans-Decalin	ND	910000	ug/kg	
493-01-6	cis-Decalin	ND	910000	ug/kg	
11095-43-5	Benzo(b)thiophene	26600000	910000	ug/kg	
91-20-3	Naphthalene	744000000	910000	ug/kg	
91-57-6	2-Methylnaphthalene	310000000	910000	ug/kg	
90-12-0	1-Methylnaphthalene	195000000	910000	ug/kg	
	C1-Naphthalenes	331000000	910000	ug/kg	
	C2-Naphthalenes	155000000	910000	ug/kg	
	C3-Naphthalenes	51800000	910000	ug/kg	
	C4-Naphthalenes	14800000	910000	ug/kg	
92-52-4	Biphenyl	27300000	910000	ug/kg	
208-96-8	Acenaphthylene	25200000	910000	ug/kg	

ND = Not detected

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

**Report of Analysis**

Page 2 of 3

<b>Client Sample ID:</b>	SB-4(15-17)NAPL	<b>Date Sampled:</b>	10/31/13
<b>Lab Sample ID:</b>	MC26103-1	<b>Date Received:</b>	11/08/13
<b>Matrix:</b>	SO - Oil	<b>Percent Solids:</b>	n/a
<b>Method:</b>	D5739-06/8270C SIM SW846 3580A		
<b>Project:</b>	Parsons, Halleck Street, NJ		

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**Alkylated PAHs**

CAS No.	Compound	Result	RL	Units	Q
83-32-9	Acenaphthene	102000000	910000	ug/kg	
132-64-9	Dibenzofuran	9390000	910000	ug/kg	
86-73-7	Fluorene	48300000	910000	ug/kg	
	C1-Fluorennes	23600000	910000	ug/kg	
	C2-Fluorennes	13900000	910000	ug/kg	
	C3-Fluorennes	6710000	910000	ug/kg	
132-65-0	Dibenzothiophene	28500000	910000	ug/kg	
	C1-Dibenzothiophenes	26300000	910000	ug/kg	
	C2-Dibenzothiophenes	16100000	910000	ug/kg	
	C3-Dibenzothiophenes	6760000	910000	ug/kg	
	C4-Dibenzothiophenes	1670000	910000	ug/kg	
85-01-8	Phenanthrene	179000000	910000	ug/kg	
120-12-7	Anthracene	53200000	910000	ug/kg	
	C1-Phenanthrenes/Anthracene	90900000	910000	ug/kg	
	C2-Phenanthrenes/Anthracene	33900000	910000	ug/kg	
	C3-Phenanthrenes/Anthracene	9630000	910000	ug/kg	
	C4-Phenanthrenes/Anthracene	2610000	910000	ug/kg	
483-65-8	Retene	1020000	910000	ug/kg	
239-35-0	Benzo(b)naphtho(2,1-d)thioph	4680000	910000	ug/kg	
206-44-0	Fluoranthene	53200000	910000	ug/kg	
129-00-0	Pyrene	87800000	910000	ug/kg	
	C1-Fluoranthenes/Pyrenes	45700000	910000	ug/kg	
	C2-Fluoranthenes/Pyrenes	11500000	910000	ug/kg	
	C3-Fluoranthenes/Pyrenes	2630000	910000	ug/kg	
243-17-4	Benzo(b)fluorene	3540000	910000	ug/kg	
205-12-9	Benzo(c)fluorene	2140000	910000	ug/kg	
3442-78-2	2-Methylpyrene	6480000	910000	ug/kg	
3353-12-6	4-Methylpyrene	5760000	910000	ug/kg	
2381-21-7	1-Methylpyrene	5610000	910000	ug/kg	
56-55-3	Benzo(a)anthracene	22200000	910000	ug/kg	
218-01-9	Chrysene	22000000	910000	ug/kg	
	C1-Benzo(a)anthracenes/Chrys	10300000	910000	ug/kg	
	C2-Benzo(a)anthracenes/Chrys	4160000	910000	ug/kg	
	C3-Benzo(a)anthracenes/Chrys	ND	910000	ug/kg	
	C4-Benzo(a)anthracenes/Chrys	ND	910000	ug/kg	
205-99-2	Benzo(b)fluoranthene	8440000	910000	ug/kg	
207-08-9	Benzo(k)fluoranthene	12200000	910000	ug/kg	
192-97-2	Benzo(e)pyrene	11700000	910000	ug/kg	
50-32-8	Benzo(a)pyrene	21800000	910000	ug/kg	
198-55-0	Perylene	2850000	910000	ug/kg	

ND = Not detected

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

**Report of Analysis**

Page 3 of 3

<b>Client Sample ID:</b>	SB-4(15-17)NAPL	<b>Date Sampled:</b>	10/31/13
<b>Lab Sample ID:</b>	MC26103-1	<b>Date Received:</b>	11/08/13
<b>Matrix:</b>	SO - Oil	<b>Percent Solids:</b>	n/a
<b>Method:</b>	D5739-06/8270C SIM SW846 3580A		
<b>Project:</b>	Parsons, Halleck Street, NJ		

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**Alkylated PAHs**

CAS No.	Compound	Result	RL	Units	Q
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193-39-5	Indeno(1,2,3-cd)pyrene	7000000	910000	ug/kg	
53-70-3	Dibenzo(a,h)anthracene	2040000	910000	ug/kg	
191-24-2	Benzo(g,h,i)perylene	9150000	910000	ug/kg	
191-07-1	Coronene	2100000	910000	ug/kg	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
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2037-26-5	Toluene-D8	85%		40-140%
1146-65-2	Naphthalene-d8	76%		40-140%
1517-22-2	Phenanthrene-d10	70%		40-140%
	Perylene-d12	58%		40-140%

(a) Sample extracted beyond the recommended holding time for soils.

ND = Not detected

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

**Report of Analysis**

Page 1 of 1

<b>Client Sample ID:</b> SB-4(15-17)NAPL	<b>Date Sampled:</b> 10/31/13
<b>Lab Sample ID:</b> MC26103-1	<b>Date Received:</b> 11/08/13
<b>Matrix:</b> SO - Oil	<b>Percent Solids:</b> n/a
<b>Method:</b> ASTM D3328-06 SW846 3580A	
<b>Project:</b> Parsons, Halleck Street, NJ	

	<b>File ID</b>	<b>DF</b>	<b>Analyzed</b>	<b>By</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analytical Batch</b>
Run #1 <sup>a</sup>	BG42080.D	1	11/20/13	RP	11/19/13	OP35833	GBG1601
Run #2							

	<b>Initial Weight</b>	<b>Final Volume</b>
Run #1	0.00220 g	10.0 ml
Run #2		

<b>CAS No.</b>	<b>Compound</b>	<b>Result</b>	<b>RL</b>	<b>Units</b>	<b>Q</b>
	TPH (C8-C40)	2510000	450000	mg/kg	

<b>CAS No.</b>	<b>Surrogate Recoveries</b>	<b>Run# 1</b>	<b>Run# 2</b>	<b>Limits</b>
84-15-1	o-Terphenyl	430% <sup>b</sup>		40-140%

(a) Sample extracted beyond the recommended holding time for soils.

(b) Outside control limits due to dilution.

ND = Not detected

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

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4



## Misc. Forms

5

### Custody Documents and Other Forms

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Includes the following where applicable:

- Chain of Custody

**PROJECT** Halleck St  
**CONTACT** Eric Gaulin  
**COMPANY** Parsons  
**ADDRESS** 200 Octagonal Ln.  
**EMAIL** somerset, NJ 08873  
**PHONE** 973 537 3570 FAX

**CHAIN OF CUSTODY RECORD**

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<b>Turn Around Time</b>	
<b>Standard</b>	<input checked="" type="checkbox"/>
If Authorized *	
1 Week	<input type="checkbox"/>
Other	<input type="checkbox"/>

META



## Environmental, Inc.

115 Dean Avenue Suite 300 Franklin MA 02038  
Tel (508) 541-9146 - Fax (508) 541-1035 - [www.metaenv.com](http://www.metaenv.com)

*100-76103*

mc26103

### Parameters

PHONE 732 537 3570			FAX									
SAMPLED BY  Print Name <u>Zohar Levy</u> Sign <u>ZL</u>												
Samp #	Date	Time	Field Sample ID		Container			Matrix	Preserv.	Comments		
			Size	G/P	Grab	Composite	# of Containers					
-1	10/31/13	1415	SG-4(15-17) NAPL		4oz	<input checked="" type="checkbox"/>	<input type="checkbox"/>	1	Soil	-	X	Hydrocarbon Figure
												20E,
Relinquished by <u>ZL</u>			Date & Time 11/6/13 @ 1500	Relinquished by <u>VPS</u>			Date & Time		Relinquished by <u>WES</u>		Date & Time 11/8/13	
Received by <u>Jewell R.</u>			Date & Time 11/8/13 10:34	Received by <u>Jewell R.</u>			Received by		Received by		Date & Time	
Shipping Info.			Remarks <u>Rec 0.50</u>			Temp °C _____						
* Surcharges may apply												

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MC26103: Chain of Custody  
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## Accutest Laboratories Sample Receipt Summary

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Accutest Job Number: MC26103

Client: META

Immediate Client Services Action Required: No

Date / Time Received: 11/8/2013

Delivery Method:

Client Service Action Required at Login: No

Project: HALLOCK ST

No. Coolers: 1

Airbill #'s:

**Cooler Security****Y or N**

1. Custody Seals Present:      3. COC Present:    
 2. Custody Seals Intact:      4. Smpl Dates/Time OK

**Cooler Temperature****Y or N**

1. Temp criteria achieved:    
 2. Cooler temp verification: Infrared gun  
 3. Cooler media: Ice (bag)

**Quality Control Preservation****Y or N****N/A**

1. Trip Blank present / cooler:     
 2. Trip Blank listed on COC:     
 3. Samples preserved properly:    
 4. VOCs headspace free:

**Sample Integrity - Documentation****Y or N**

1. Sample labels present on bottles:    
 2. Container labeling complete:    
 3. Sample container label / COC agree:

**Sample Integrity - Condition****Y or N**

1. Sample recvd within HT:    
 2. All containers accounted for:    
 3. Condition of sample: Intact

**Sample Integrity - Instructions****Y or N****N/A**

1. Analysis requested is clear:    
 2. Bottles received for unspecified tests:    
 3. Sufficient volume recvd for analysis:    
 4. Compositing instructions clear:     
 5. Filtering instructions clear:

Comments

Accutest Laboratories  
V:508.481.6200495 Technology Center West, Bldg One  
F: 508.481.7753Marlborough, MA  
www.accutest.com**MC26103: Chain of Custody****Page 2 of 2**

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## GC/MS Semi-volatiles

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### QC Data Summaries

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Includes the following where applicable:

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries
- Surrogate Recovery Summaries

**Method Blank Summary**

Page 1 of 3

**Job Number:** MC26103  
**Account:** METAMAW META Environmental, Inc.  
**Project:** Parsons, Halleck Street, NJ

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP35834-MB	S34116.D	1	11/20/13	RP	11/19/13	OP35834	MSS1523

The QC reported here applies to the following samples:

Method: D5739-06/8270C SIM

MC26103-1

CAS No.	Compound	Result	RL	Units	Q
71-43-2	Benzene	9220	2000	ug/kg	
	C1-Benzene	ND	2000	ug/kg	
	C2-Benzenes	ND	2000	ug/kg	
	C3-Benzenes	ND	2000	ug/kg	
	C4-Benzenes	ND	2000	ug/kg	
	C5-Benzenes	ND	2000	ug/kg	
108-88-3	Toluene	ND	2000	ug/kg	
100-41-4	Ethylbenzene	ND	2000	ug/kg	
	m,p-Xylene	ND	2000	ug/kg	
100-42-5	Styrene	ND	2000	ug/kg	
95-47-6	o-Xylene	ND	2000	ug/kg	
98-82-8	Isopropylbenzene	ND	2000	ug/kg	
103-65-1	n-Propylbenzene	ND	2000	ug/kg	
108-67-8	1,3,5-Trimethylbenzene	ND	2000	ug/kg	
526-73-8	1,2,3-Trimethylbenzene	ND	2000	ug/kg	
95-63-6	1,2,4-Trimethylbenzene	ND	2000	ug/kg	
98-06-6	t-Butylbenzene	ND	2000	ug/kg	
135-98-8	sec-Butylbenzene	ND	2000	ug/kg	
99-87-6	p-Isopropyltoluene	ND	2000	ug/kg	
104-51-8	n-Butylbenzene	ND	2000	ug/kg	
493-02-7	trans-Decalin	ND	2000	ug/kg	
493-01-6	cis-Decalin	ND	2000	ug/kg	
11095-43-5	Benzo(b)thiophene	ND	2000	ug/kg	
91-20-3	Naphthalene	ND	2000	ug/kg	
91-57-6	2-Methylnaphthalene	ND	2000	ug/kg	
90-12-0	1-Methylnaphthalene	ND	2000	ug/kg	
	C1-Naphthalenes	ND	2000	ug/kg	
	C2-Naphthalenes	ND	2000	ug/kg	
	C3-Naphthalenes	ND	2000	ug/kg	
	C4-Naphthalenes	ND	2000	ug/kg	
92-52-4	Biphenyl	ND	2000	ug/kg	
208-96-8	Acenaphthylene	ND	2000	ug/kg	
83-32-9	Acenaphthene	ND	2000	ug/kg	
132-64-9	Dibenzofuran	ND	2000	ug/kg	
86-73-7	Fluorene	ND	2000	ug/kg	
	C1-Fluorenes	ND	2000	ug/kg	

6.1.1  
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**Method Blank Summary**

Page 2 of 3

**Job Number:** MC26103  
**Account:** METAMAW META Environmental, Inc.  
**Project:** Parsons, Halleck Street, NJ

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP35834-MB	S34116.D	1	11/20/13	RP	11/19/13	OP35834	MSS1523

The QC reported here applies to the following samples:

Method: D5739-06/8270C SIM

MC26103-1

CAS No.	Compound	Result	RL	Units	Q
132-65-0	C2-Fluorenes	ND	2000	ug/kg	
	C3-Fluorenes	ND	2000	ug/kg	
	Dibenzothiophene	ND	2000	ug/kg	
	C1-Dibenzothiophenes	ND	2000	ug/kg	
	C2-Dibenzothiophenes	ND	2000	ug/kg	
	C3-Dibenzothiophenes	ND	2000	ug/kg	
85-01-8	C4-Dibenzothiophenes	ND	2000	ug/kg	
	Phenanthrene	ND	2000	ug/kg	
120-12-7	Anthracene	ND	2000	ug/kg	
	C1-Phenanthrenes/Anthracene	ND	2000	ug/kg	
	C2-Phenanthrenes/Anthracene	ND	2000	ug/kg	
	C3-Phenanthrenes/Anthracene	ND	2000	ug/kg	
	C4-Phenanthrenes/Anthracene	ND	2000	ug/kg	
483-65-8	Retene	ND	2000	ug/kg	
239-35-0	Benzo(b)naphtho(2,1-d)thioph	ND	2000	ug/kg	
206-44-0	Fluoranthene	ND	2000	ug/kg	
129-00-0	Pyrene	ND	2000	ug/kg	
	C1-Fluoranthenes/Pyrenes	ND	2000	ug/kg	
	C2-Fluoranthenes/Pyrenes	ND	2000	ug/kg	
	C3-Fluoranthenes/Pyrenes	ND	2000	ug/kg	
243-17-4	Benzo(b)fluorene	ND	2000	ug/kg	
205-12-9	Benzo(c)fluorene	ND	2000	ug/kg	
3442-78-2	2-Methylpyrene	ND	2000	ug/kg	
3353-12-6	4-Methylpyrene	ND	2000	ug/kg	
2381-21-7	1-Methylpyrene	ND	2000	ug/kg	
56-55-3	Benzo(a)anthracene	ND	2000	ug/kg	
218-01-9	Chrysene	ND	2000	ug/kg	
	C1-Benzo(a)anthracenes/Chrys	ND	2000	ug/kg	
	C2-Benzo(a)anthracenes/Chrys	ND	2000	ug/kg	
	C3-Benzo(a)anthracenes/Chrys	ND	2000	ug/kg	
	C4-Benzo(a)anthracenes/Chrys	ND	2000	ug/kg	
205-99-2	Benzo(b)fluoranthene	ND	2000	ug/kg	
207-08-9	Benzo(k)fluoranthene	ND	2000	ug/kg	
192-97-2	Benzo(e)pyrene	ND	2000	ug/kg	
50-32-8	Benzo(a)pyrene	ND	2000	ug/kg	
198-55-0	Perylene	ND	2000	ug/kg	

6.1.1  
6

**Method Blank Summary**

Page 3 of 3

**Job Number:** MC26103  
**Account:** METAMAW META Environmental, Inc.  
**Project:** Parsons, Halleck Street, NJ

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP35834-MB	S34116.D	1	11/20/13	RP	11/19/13	OP35834	MSS1523

The QC reported here applies to the following samples:

Method: D5739-06/8270C SIM

MC26103-1

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6

CAS No.	Compound	Result	RL	Units	Q
193-39-5	Indeno(1,2,3-cd)pyrene	ND	2000	ug/kg	
53-70-3	Dibenz(a,h)anthracene	ND	2000	ug/kg	
191-24-2	Benzo(g,h,i)perylene	ND	2000	ug/kg	
191-07-1	Coronene	ND	2000	ug/kg	

CAS No.	Surrogate Recoveries	Limits
2037-26-5	Toluene-D8	108% 40-140%
1146-65-2	Naphthalene-d8	90% 40-140%
1517-22-2	Phenanthrene-d10	88% 40-140%
	Perylene-d12	88% 40-140%

**Blank Spike Summary**

Page 1 of 3

**Job Number:** MC26103**Account:** METAMAW META Environmental, Inc.**Project:** Parsons, Halleck Street, NJ

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP35834-BS2	S34117.D	1	11/20/13	RP	11/19/13	OP35834	MSS1523

**The QC reported here applies to the following samples:****Method:** D5739-06/8270C SIM

MC26103-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
71-43-2	Benzene	1000000	966000	97	50-150 <sup>a</sup>
	C1-Benzene		ND		50-150 <sup>a</sup>
	C2-Benzene		ND		50-150 <sup>a</sup>
	C3-Benzene		ND		50-150 <sup>a</sup>
	C4-Benzene		ND		50-150 <sup>a</sup>
	C5-Benzene		ND		50-150 <sup>a</sup>
108-88-3	Toluene	1000000	918000	92	50-150 <sup>a</sup>
100-41-4	Ethylbenzene	1000000	888000	89	50-150 <sup>a</sup>
	m,p-Xylene	1000000	896000	90	50-150 <sup>a</sup>
100-42-5	Styrene	1000000	878000	88	50-150 <sup>a</sup>
95-47-6	o-Xylene	1000000	907000	91	50-150 <sup>a</sup>
98-82-8	Isopropylbenzene	1000000	930000	93	50-150 <sup>a</sup>
103-65-1	n-Propylbenzene	1000000	916000	92	50-150 <sup>a</sup>
108-67-8	1,3,5-Trimethylbenzene	1000000	920000	92	50-150 <sup>a</sup>
526-73-8	1,2,3-Trimethylbenzene	1000000	922000	92	50-150 <sup>a</sup>
95-63-6	1,2,4-Trimethylbenzene	1000000	922000	92	50-150 <sup>a</sup>
98-06-6	t-Butylbenzene	1000000	937000	94	50-150 <sup>a</sup>
135-98-8	sec-Butylbenzene	1000000	947000	95	50-150 <sup>a</sup>
99-87-6	p-Isopropyltoluene	1000000	951000	95	50-150 <sup>a</sup>
104-51-8	n-Butylbenzene	1000000	922000	92	50-150 <sup>a</sup>
493-02-7	trans-Decalin	1000000	909000	91	50-150 <sup>a</sup>
493-01-6	cis-Decalin	1000000	933000	93	50-150 <sup>a</sup>
11095-43-5	Benzo(b)thiophene	1000000	996000	100	50-150 <sup>a</sup>
91-20-3	Naphthalene	1000000	951000	95	50-150
91-57-6	2-Methylnaphthalene	1000000	923000	92	50-150
90-12-0	1-Methylnaphthalene	1000000	943000	94	50-150
	C1-Naphthalenes		ND		50-150 <sup>a</sup>
	C2-Naphthalenes		ND		50-150 <sup>a</sup>
	C3-Naphthalenes		ND		50-150 <sup>a</sup>
	C4-Naphthalenes		ND		50-150 <sup>a</sup>
92-52-4	Biphenyl	1000000	897000	90	50-150 <sup>a</sup>
208-96-8	Acenaphthylene	1000000	1050000	105	50-150
83-32-9	Acenaphthene	1000000	991000	99	50-150
132-64-9	Dibenzofuran	1000000	998000	100	50-150
86-73-7	Fluorene	1000000	957000	96	50-150
	C1-Fluorenes		ND		50-150 <sup>a</sup>

\* = Outside of Control Limits.

6.2.1  
**6**

**Blank Spike Summary**

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**Job Number:** MC26103**Account:** METAMAW META Environmental, Inc.**Project:** Parsons, Halleck Street, NJ

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP35834-BS2	S34117.D	1	11/20/13	RP	11/19/13	OP35834	MSS1523

**The QC reported here applies to the following samples:****Method:** D5739-06/8270C SIM

MC26103-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits	
132-65-0	C2-Fluorenes		ND		50-150 <sup>a</sup>	
	C3-Fluorenes		ND		50-150 <sup>a</sup>	
	Dibenzothiophene	1000000	968000	97	50-150 <sup>a</sup>	
	C1-Dibenzothiophenes		ND		50-150 <sup>a</sup>	
	C2-Dibenzothiophenes		ND		50-150 <sup>a</sup>	
	C3-Dibenzothiophenes		ND		50-150 <sup>a</sup>	
85-01-8	C4-Dibenzothiophenes		ND		50-150 <sup>a</sup>	
	Phenanthrene	1000000	954000	95	50-150	
	Anthracene	1000000	998000	100	50-150	
	C1-Phenanthrenes/Anthracene		ND		50-150 <sup>a</sup>	
	C2-Phenanthrenes/Anthracene		ND		50-150 <sup>a</sup>	
	C3-Phenanthrenes/Anthracene		ND		50-150 <sup>a</sup>	
483-65-8	C4-Phenanthrenes/Anthracene		ND		50-150 <sup>a</sup>	
	Retene	1000000	932000	93	50-150 <sup>a</sup>	
	239-35-0	Benzo(b)naphtho(2,1-d)thioph	1000000	850000	85	50-150 <sup>a</sup>
	206-44-0	Fluoranthene	1000000	946000	95	50-150
	129-00-0	Pyrene	1000000	928000	93	50-150
	C1-Fluoranthenes/Pyrenes		ND		50-150 <sup>a</sup>	
243-17-4	C2-Fluoranthenes/Pyrenes		ND		50-150 <sup>a</sup>	
	C3-Fluoranthenes/Pyrenes		ND		50-150 <sup>a</sup>	
	Benzo(b)fluorene		ND		50-150 <sup>a</sup>	
	205-12-9	Benzo(c)fluorene		ND	50-150 <sup>a</sup>	
	3442-78-2	2-Methylpyrene		ND	50-150 <sup>a</sup>	
	3353-12-6	4-Methylpyrene		ND	50-150 <sup>a</sup>	
2381-21-7	2381-21-7	1-Methylpyrene		ND	50-150 <sup>a</sup>	
	56-55-3	Benzo(a)anthracene	1000000	840000	84	50-150
	218-01-9	Chrysene	1000000	845000	85	50-150
	C1-Benzo(a)anthracenes/Chrys		ND		50-150 <sup>a</sup>	
	C2-Benzo(a)anthracenes/Chrys		ND		50-150 <sup>a</sup>	
	C3-Benzo(a)anthracenes/Chrys		ND		50-150 <sup>a</sup>	
205-99-2	C4-Benzo(a)anthracenes/Chrys		ND		50-150 <sup>a</sup>	
	Benzo(b)fluoranthene	1000000	861000	86	50-150	
	207-08-9	Benzo(k)fluoranthene	1000000	846000	85	50-150
	192-97-2	Benzo(e)pyrene	1000000	849000	85	50-150 <sup>a</sup>
	50-32-8	Benzo(a)pyrene	1000000	824000	82	50-150
	198-55-0	Perylene	1000000	846000	85	50-150 <sup>a</sup>

\* = Outside of Control Limits.

6.2.1  
**6**

**Blank Spike Summary**

Page 3 of 3

**Job Number:** MC26103**Account:** METAMAW META Environmental, Inc.**Project:** Parsons, Halleck Street, NJ

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP35834-BS2	S34117.D	1	11/20/13	RP	11/19/13	OP35834	MSS1523

**The QC reported here applies to the following samples:****Method:** D5739-06/8270C SIM

MC26103-1

6.2.1  
6

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
193-39-5	Indeno(1,2,3-cd)pyrene	1000000	843000	84	50-150
53-70-3	Dibenz(a,h)anthracene	1000000	727000	73	50-150
191-24-2	Benzo(g,h,i)perylene	1000000	838000	84	50-150
191-07-1	Coronene	1000000	820000	82	50-150 <sup>a</sup>

CAS No.	Surrogate Recoveries	BSP	Limits
2037-26-5	Toluene-D8	102%	40-140%
1146-65-2	Naphthalene-d8	95%	40-140%
1517-22-2	Phenanthrene-d10	93%	40-140%
	Perylene-d12	79%	40-140%

(a) Advisory control limits.

\* = Outside of Control Limits.

**Duplicate Summary**

Page 1 of 3

**Job Number:** MC26103**Account:** METAMAW META Environmental, Inc.**Project:** Parsons, Halleck Street, NJ

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP35834-DUP1	S34120.D	10	11/20/13	RP	11/19/13	OP35834	MSS1523
MC26103-1 <sup>a</sup>	S34121.D	10	11/20/13	RP	11/19/13	OP35834	MSS1523

**The QC reported here applies to the following samples:****Method:** D5739-06/8270C SIM

MC26103-1

6.3.1  
6

CAS No.	Compound	MC26103-1 DUP					
		ug/kg	Q	ug/kg	Q	RPD	Limits
71-43-2	Benzene	10900000	3930000	94* b	30 c		
	C1-Benzene	1580000	745000	72* b	30 c		
	C2-Benzenes	72300000	28700000	86* b	30 c		
	C3-Benzenes	82600000	35800000	79* b	30 c		
	C4-Benzenes	43300000	21800000	66* b	30 c		
	C5-Benzenes	12800000	7390000	54* b	30 c		
108-88-3	Toluene	2190000	999000	75* b	30 c		
100-41-4	Ethylbenzene	55300000	21300000	89* b	30 c		
	m,p-Xylene	52800000	20600000	88* b	30 c		
100-42-5	Styrene	11300000	7770000	37* b	30 c		
95-47-6	o-Xylene	23800000	9760000	84* b	30 c		
98-82-8	Isopropylbenzene	6660000	2710000	84* b	30 c		
103-65-1	n-Propylbenzene	2440000	1080000	77* b	30 c		
108-67-8	1,3,5-Trimethylbenzene	13200000	5830000	77* b	30 c		
526-73-8	1,2,3-Trimethylbenzene	13500000	6160000	75* b	30 c		
95-63-6	1,2,4-Trimethylbenzene	40900000	18100000	77* b	30 c		
98-06-6	t-Butylbenzene	5370000	2530000	72* b	30 c		
135-98-8	sec-Butylbenzene	ND	ND	nc	30 c		
99-87-6	p-Isopropyltoluene	6740000	3070000	75* b	30 c		
104-51-8	n-Butylbenzene	1480000	802000	59* b	30 c		
493-02-7	trans-Decalin	ND	ND	nc	30 c		
493-01-6	cis-Decalin	ND	ND	nc	30 c		
11095-43-5	Benzo(b)thiophene	26600000	13600000	65* b	30 c		
91-20-3	Naphthalene	744000000	374000000	66* b	30		
91-57-6	2-Methylnaphthalene	310000000	158000000	65* b	30		
90-12-0	1-Methylnaphthalene	195000000	107000000	58* b	30		
	C1-Naphthalenes	331000000	174000000	62* b	30 c		
	C2-Naphthalenes	155000000	89500000	54* b	30 c		
	C3-Naphthalenes	51800000	30800000	51* b	30 c		
	C4-Naphthalenes	14800000	8960000	49* b	30 c		
92-52-4	Biphenyl	27300000	14700000	60* b	30 c		
208-96-8	Acenaphthylene	25200000	17600000	36* b	30		
83-32-9	Acenaphthene	102000000	56700000	57* b	30		
132-64-9	Dibenzofuran	9390000	5330000	55* b	30		
86-73-7	Fluorene	48300000	27400000	55* b	30		
	C1-Fluorennes	23600000	14100000	50* b	30 c		

\* = Outside of Control Limits.

**Duplicate Summary**

Page 2 of 3

**Job Number:** MC26103**Account:** METAMAW META Environmental, Inc.**Project:** Parsons, Halleck Street, NJ

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP35834-DUP1	S34120.D	10	11/20/13	RP	11/19/13	OP35834	MSS1523
MC26103-1 <sup>a</sup>	S34121.D	10	11/20/13	RP	11/19/13	OP35834	MSS1523

**The QC reported here applies to the following samples:****Method:** D5739-06/8270C SIM

MC26103-1

6.3.1  
6

CAS No.	Compound	MC26103-1 DUP					
		ug/kg	Q	ug/kg	Q	RPD	Limits
132-65-0	C2-Fluorenes	13900000	8690000	46* b	30 c		
	C3-Fluorenes	6710000	4790000	33* b	30 c		
	Dibenzothiophene	28500000	17100000	50* b	30 c		
	C1-Dibenzothiophenes	26300000	16600000	45* b	30 c		
	C2-Dibenzothiophenes	16100000	10500000	42* b	30 c		
	C3-Dibenzothiophenes	6760000	4510000	40* b	30 c		
85-01-8	C4-Dibenzothiophenes	1670000	1300000	25	30 c		
	Phenanthrene	179000000	109000000	49* b	30		
120-12-7	Anthracene	53200000	31900000	50* b	30		
	C1-Phenanthrenes/Anthracene	90900000	56500000	47* b	30 c		
	C2-Phenanthrenes/Anthracene	33900000	22100000	42* b	30 c		
	C3-Phenanthrenes/Anthracene	9630000	6430000	40* b	30 c		
	C4-Phenanthrenes/Anthracene	2610000	2000000	26	30 c		
483-65-8	Retene	1020000	1150000	12	30 c		
239-35-0	Benzo(b)naphtho(2,1-d)thioph	4680000	3150000	39* b	30 c		
206-44-0	Fluoranthene	53200000	34700000	42* b	30		
129-00-0	Pyrene	87800000	58400000	40* b	30		
	C1-Fluoranthenes/Pyrenes	45700000	30000000	41* b	30 c		
	C2-Fluoranthenes/Pyrenes	11500000	7710000	39* b	30 c		
	C3-Fluoranthenes/Pyrenes	2630000	2230000	16	30 c		
243-17-4	Benzo(b)fluorene	3540000	1910000	60* b	30 c		
205-12-9	Benzo(c)fluorene	2140000	1420000	40* b	30 c		
3442-78-2	2-Methylpyrene	6480000	4350000	39* b	30 c		
3353-12-6	4-Methylpyrene	5760000	3740000	43* b	30 c		
2381-21-7	1-Methylpyrene	5610000	3760000	39* b	30 c		
56-55-3	Benzo(a)anthracene	22200000	14200000	44* b	30		
218-01-9	Chrysene	22000000	14800000	39* b	30		
	C1-Benzo(a)anthracenes/Chrys	10300000	6890000	40* b	30 c		
	C2-Benzo(a)anthracenes/Chrys	4160000	3070000	30	30 c		
	C3-Benzo(a)anthracenes/Chrys	ND	ND	nc	30 c		
	C4-Benzo(a)anthracenes/Chrys	ND	ND	nc	30 c		
205-99-2	Benzo(b)fluoranthene	8440000	5580000	41* b	30		
207-08-9	Benzo(k)fluoranthene	12200000	8280000	38* b	30		
192-97-2	Benzo(e)pyrene	11700000	8010000	37* b	30 c		
50-32-8	Benzo(a)pyrene	21800000	14200000	42* b	30		
198-55-0	Perylene	2850000	2070000	32* b	30 c		

\* = Outside of Control Limits.

**Duplicate Summary**

Page 3 of 3

**Job Number:** MC26103**Account:** METAMAW META Environmental, Inc.**Project:** Parsons, Halleck Street, NJ

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP35834-DUP1	S34120.D	10	11/20/13	RP	11/19/13	OP35834	MSS1523
MC26103-1 <sup>a</sup>	S34121.D	10	11/20/13	RP	11/19/13	OP35834	MSS1523

**The QC reported here applies to the following samples:****Method:** D5739-06/8270C SIM

MC26103-1

6.3.1  
6

CAS No.	Compound	MC26103-1 DUP					
		ug/kg	Q	ug/kg	Q	RPD	Limits
193-39-5	Indeno(1,2,3-cd)pyrene	7000000		4830000		37* <sup>b</sup>	30
53-70-3	Dibenzo(a,h)anthracene	2040000		1590000		25	30
191-24-2	Benzo(g,h,i)perylene	9150000		6310000		37* <sup>b</sup>	30
191-07-1	Coronene	2100000		1420000		39* <sup>b</sup>	30 <sup>c</sup>

CAS No.	Surrogate Recoveries	DUP	MC26103-1 Limits		
2037-26-5	Toluene-D8	82%	85%	40-140%	
1146-65-2	Naphthalene-d8	73%	76%	40-140%	
1517-22-2	Phenanthrene-d10	69%	70%	40-140%	
	Perylene-d12	66%	58%	40-140%	

(a) Sample extracted beyond the recommended holding time for soils.

(b) High RPD due to possible sample heterogeneity.

(c) Advisory control limits.

\* = Outside of Control Limits.

**Semivolatile Surrogate Recovery Summary**

Page 1 of 1

**Job Number:** MC26103**Account:** METAMAW META Environmental, Inc.**Project:** Parsons, Halleck Street, NJ**Method:** D5739-06/8270C SIM**Matrix:** SO**Samples and QC shown here apply to the above method**

<b>Lab Sample ID</b>	<b>Lab File ID</b>	<b>S1</b>	<b>S2</b>	<b>S3</b>	<b>S4</b>
MC26103-1	S34121.D	85.0	76.0	70.0	58.0
OP35834-BS2	S34117.D	102.0	95.0	93.0	79.0
OP35834-DUP1	S34120.D	82.0	73.0	69.0	66.0
OP35834-MB	S34116.D	108.0	90.0	88.0	88.0

<b>Surrogate Compounds</b>	<b>Recovery Limits</b>
--------------------------------	----------------------------

<b>S1</b> = Toluene-D8	40-140%
<b>S2</b> = Naphthalene-d8	40-140%
<b>S3</b> = Phenanthrene-d10	40-140%
<b>S4</b> = Perylene-d12	40-140%

6.4.1

6



## GC/MS Semi-volatiles

---

Raw Data

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2

Data Path : Z:\1\data\S131120\  
 Data File : S34121.D  
 Acq On : 20 Nov 2013 7:05 pm  
 Operator : RUBENP  
 Sample : mc26103-1  
 Misc : op35834,mss1523,0.0011,,,10,10  
 ALS Vial : 8 Sample Multiplier: 1

Manual Integrations  
 APPROVED  
 71 of 133  
 (compounds with "m" flag)

James Roush  
 11/25/13 13:58

Quant Time: Nov 21 09:29:26 2013  
 Quant Method : Z:\1\methods\S131114-MAHPAHEXT.M  
 Quant Title : PAHs & Alkylated PAHs by GC/MS/SIM  
 QLast Update : Fri Nov 15 10:17:28 2013  
 Response via : Initial Calibration

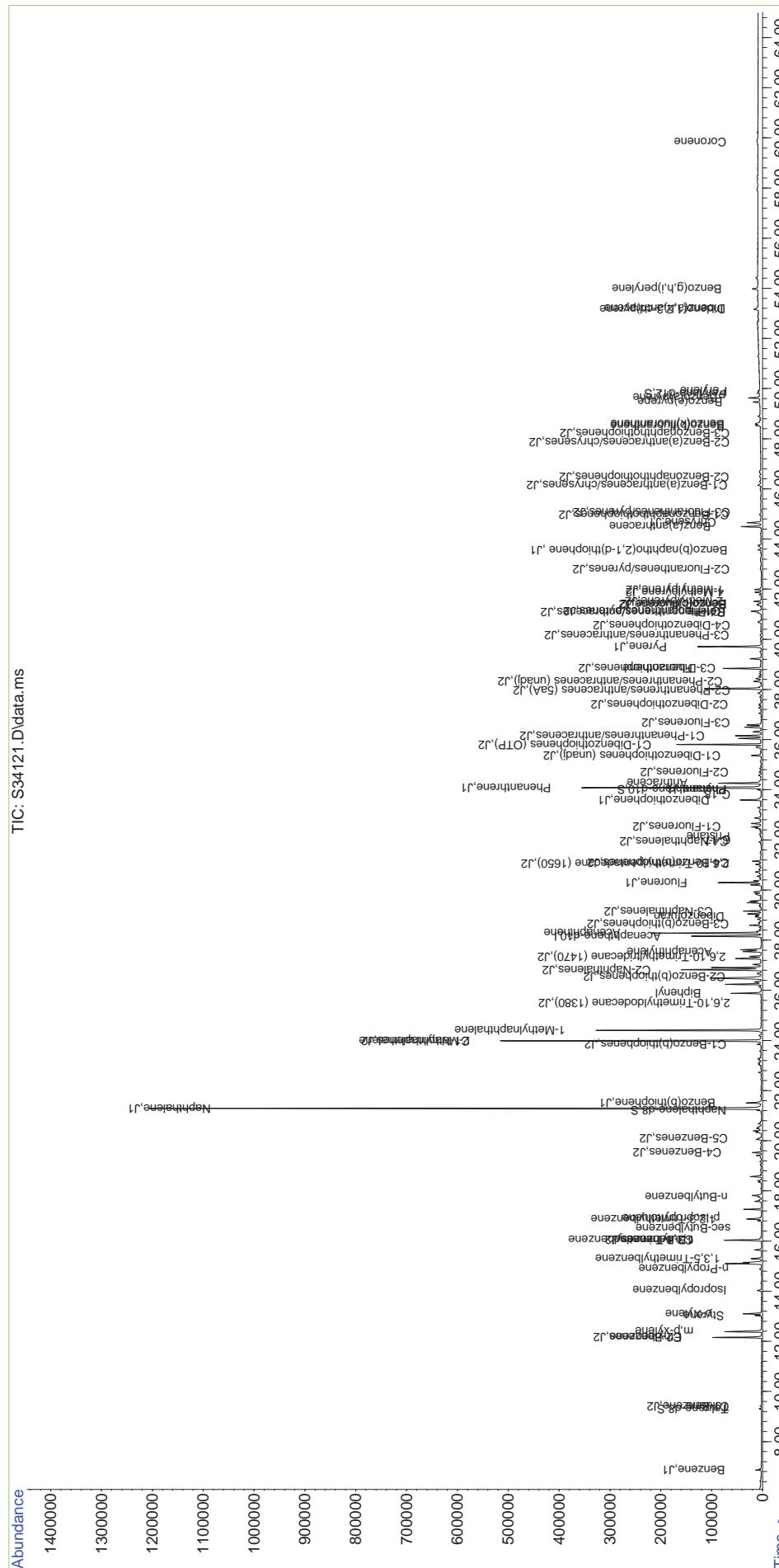
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<hr/>						
Internal Standards						
1) Acenaphthene-d10	28.158	164	120191	1100.00	ng/mL	0.00
<hr/>						
System Monitoring Compounds						
2) Toluene-d8	9.307	98	9456	84.93	µg/mL	0.00
Spiked Amount 1000.000			Recovery =	8.49%		
3) Naphthalene-d8	21.209	136	15849	76.25	ng/mL	0.00
Spiked Amount 1000.000	Range 1 - 0		Recovery =	7.63%#		
4) Phenanthrene-d10	33.992	188	13833	69.53	ng/mL	0.00
Spiked Amount 1000.000			Recovery =	6.95%		
5) Perylene-d12	49.822	264	10085	58.06	ng/mL	-0.01
Spiked Amount 1000.000			Recovery =	5.81%		
<hr/>						
Target Compounds						
7) Benzene	6.865	78	14211	119.487	µg/mL	96
8) C1-Benzene	9.415	92	2073	17.430	µg/mL	97
9) C2-Benzenes	12.152	106	94637m	795.712	µg/mL	
10) C3-Benzene	16.035	120	108018m	908.220	µg/mL	
11) C4-Benzene	19.528	134	56594m	475.845	µg/mL	
12) C5-Benzene	20.097	148	16710m	140.498	µg/mL	
14) Toluene	9.415	91	3708	24.070	µg/mL	98
15) Ethylbenzene	12.152	91	120813	608.249	µg/mL	97
16) m,p-xylene	12.383	91	93177	580.275	µg/mL	98
17) Styrene	13.031	104	14443	124.221	µg/mL	98
18) o-Xylene	13.096	91	41864	261.337	µg/mL	97
19) Isopropylbenzene	14.025	105	13102	73.218	µg/mL	100
20) n-Propylbenzene	14.882	91	6330	26.872	µg/mL	96
21) 1,3,5-Trimethylbenzene	15.300	105	25275	145.739	µg/mL	100
22) t-Butylbenzene	16.035	119	10280	59.123	µg/mL#	55
23) 1,2,4-Trimethylbenzene	16.035	105	79015	449.801	µg/mL	97
24) sec-Butylbenzene	16.525	105	594	2.786	µg/mL#	57
25) 1,2,3-Trimethylbenzene	16.877	105	27598	148.973	µg/mL	99
26) p-Isopropyltoluene	16.892	119	15318	74.127	µg/mL	93
27) n-Butylbenzene	17.771	91	2993	16.278	µg/mL#	1
34) Benzo(b)thiophene	21.506	134	59298	292.581	µg/mL	99
35) C1-Benzo(b)thiophenes	23.849	148	50004m	246.723	µg/mL	
36) C2-Benzo(b)thiophenes	26.538	162	43026m	212.293	µg/mL	
37) C3-Benzo(b)thiophenes	28.640	176	16346m	80.652	µg/mL	
38) C4-Benzo(b)thiophenes	31.092	190	7227m	35.659	µg/mL	
44) Naphthalene	21.284	128	2033676	8188.789	ng/mL	100
45) 2-Methylnaphthalene	23.989	142	558449	3407.199	ng/mL	100
46) 1-Methylnaphthalene	24.401	142	343419	2145.763	ng/mL	100
47) C1-Naphthalenes	23.989	142	905257m	3645.103	ng/mL	
48) C2-Naphthalenes	26.818	156	424526m	1709.394	ng/mL	
49) C3-Naphthalenes	29.158	170	141598m	570.158	ng/mL	
50) C4-Naphthalenes	31.918	184	40532m	163.206	ng/mL	
51) Biphenyl	25.886	154	66901	300.246	µg/mL	100
52) Acenaphthylene	27.544	152	64940	276.928	ng/mL	95
53) Acenaphthene	28.280	154	159791	1120.237	ng/mL	99
54) Dibenzofuran	28.935	168	22453	103.239	ng/mL	99
55) Fluorene	30.292	166	91616	530.858	ng/mL	100
56) C1-Fluorenes	32.520	180	44887m	260.092	ng/mL	
57) C2-Fluorenes	34.703	194	26382m	152.867	ng/mL	
58) C3-Fluorenes	36.685	208	12737m	73.803	ng/mL	
59) Dibenzothiophene	33.589	184	76682	313.116	ng/mL#	83
60) C1-Dibenzothiophenes (...	35.364	198	72602m	296.456	ng/mL	
61) C1-Dibenzothiophenes (...	35.800	198	1669m	6.815	µg/mL	
62) C2-Dibenzothiophenes	37.410	212	43395m	177.195	ng/mL	
63) C3-Dibenzothiophenes	38.859	226	18216m	74.382	ng/mL	
64) C4-Dibenzothiophenes	40.558	240	4491m	18.338	ng/mL	
65) Phenanthrene	34.082	178	493163	1971.726	ng/mL	100
66) Anthracene	34.261	178	132394	585.517	ng/mL	99
67) C1-Phenanthrenes/anthr...	36.153	192	250005m	999.551	ng/mL	

Data Path : Z:\1\data\S131120\  
 Data File : S34121.D  
 Acq On : 20 Nov 2013 7:05 pm  
 Operator : RUBENP  
 Sample : mc26103-1  
 Misc : op35834,mss1523,0.0011,,,10,10  
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: Nov 21 09:29:26 2013  
 Quant Method : Z:\1\methods\S131114-MAHPAHEXT.M  
 Quant Title : PAHs & Alkylated PAHs by GC/MS/SIM  
 QLast Update : Fri Nov 15 10:17:28 2013  
 Response via : Initial Calibration

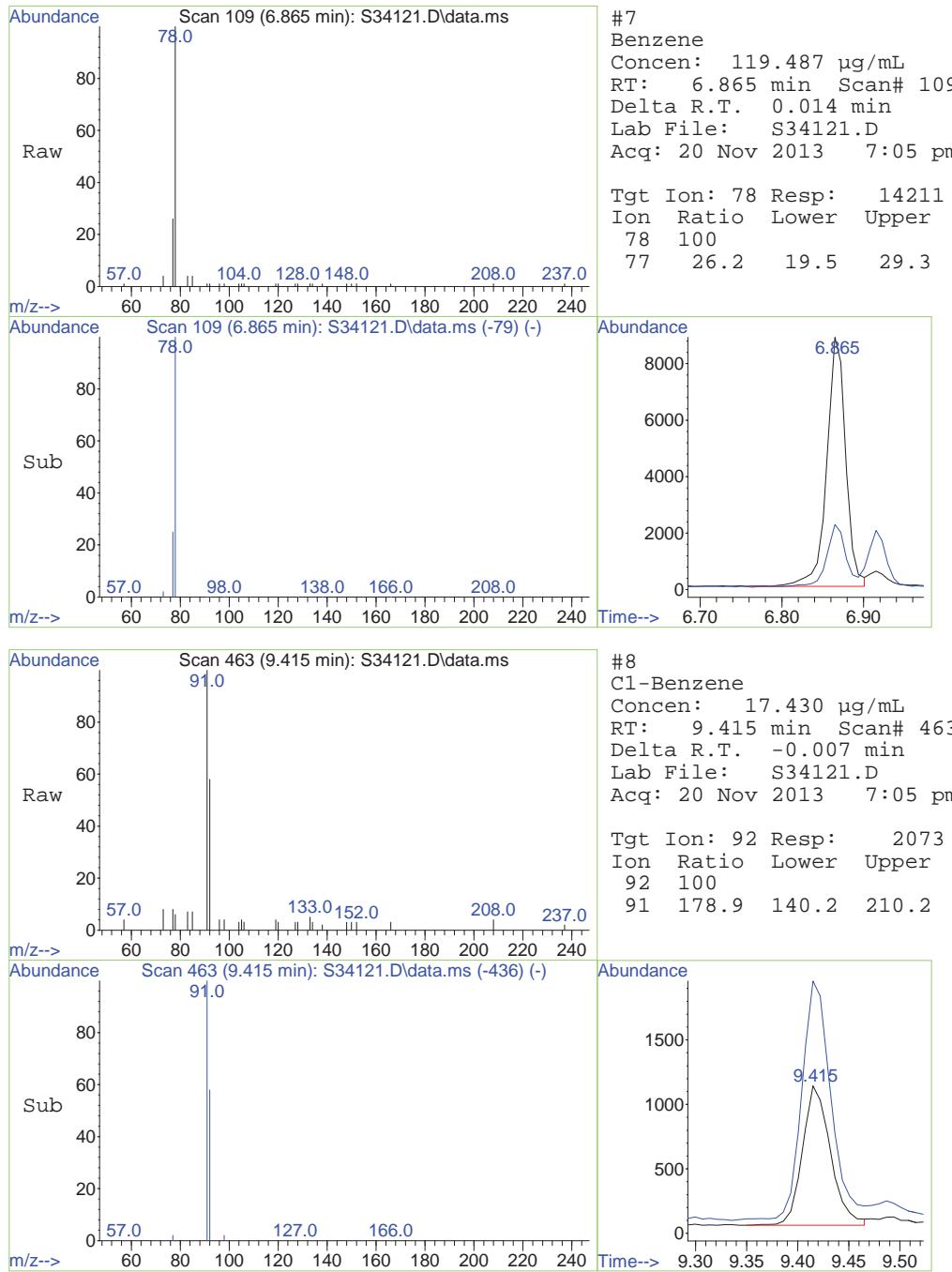
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
68) C2-Phenanthrenes/anthr...	38.320	206	93851m	375.228	ng/mL	
69) C2-Phenanthrenes/anthr...	37.942	206	470m	1.879	µg/mL	
70) C3-Phenanthrenes/anthr...	40.164	220	26487m	105.898	ng/mL	
71) C4-Phenanthrenes/anthr...	41.093	234	7176m	28.691	ng/mL	
72) Retene	41.093	234	396	11.230	µg/mL	99
73) Benzo(b)naphtho(2,1-d)...	43.563	234	11879	51.519	µg/mL	100
74) C1-Benzonaphthothiophenes	44.970	248	12772m	55.392	µg/mL	
75) C2-Benzonaphthothiophenes	46.480	262	5127m	22.236	µg/mL	
76) C3-Benzonaphthothiophenes	48.213	276	2883m	12.504	µg/mL	
78) Fluoranthene	38.841	202	151671	585.291	ng/mL	98
79) Pyrene	39.711	202	254706	966.243	ng/mL	98
80) C1-Fluoranthenes/pyrenes	41.117	216	132561m	502.878	ng/mL	
81) Benzo(b)fluorene	41.364	216	10264m	38.937	ng/mL	
82) Benzo(c)fluorene	41.405	216	6209	23.554	µg/mL#	1
83) 2-Methylpyrene	41.511	216	18795	71.300	ng/mL#	57
84) 4-Methylpyrene	41.875	216	16702m	63.360	ng/mL	
85) 1-Methylpyrene	41.987	216	16268m	61.714	ng/mL	
86) C2-Fluoranthenes/pyrenes	42.799	230	33370m	126.591	ng/mL	
87) C3-Fluoranthenes/pyrenes	45.057	244	7637m	28.971	ng/mL	
88) Benz(a)anthracene	44.496	228	54738m	244.165	ng/mL	
89) Chrysene	44.657	228	51902	242.378	ng/mL	100
90) C1-Benz(a)anthracenes/...	46.146	242	24173m	112.886	ng/mL	
91) C2-Benz(a)anthracenes/...	47.855	256	9795m	45.742	ng/mL	
94) Benzo(b)fluoranthene	48.534	252	22478	92.879	ng/mL	99
95) Benzo(k)fluoranthene	48.596	252	33226	133.898	ng/mL	96
96) Benzo(e)pyrene	49.462	252	29034	128.200	ng/mL	99
97) Benzo(a)pyrene	49.631	252	50241	239.760	ng/mL	99
98) Perylene	49.901	252	6459	31.385	ng/mL	99
99) Indeno(1,2,3-cd)pyrene	53.163	276	19394m	77.053	ng/mL	
100) Dibenz(a,h)anthracene	53.201	278	5487	22.421	ng/mL#	59
101) Benzo(g,h,i)perylene	53.978	276	26489	100.628	ng/mL	100
102) Coronene	59.847	300	5929	23.124	µg/mL#	86
103) C-17	31.989	85	131m	2.654	µg/mL	
104) Pristane	32.123	85	1101	29.196	µg/mL#	60
105) C-18	33.800	85	160m	3.196	µg/mL	
106) Phytane	33.992	85	773	14.341	µg/mL#	64
107) 2,6,10-Trimethyldodeca...	25.490	85	973	18.051	µg/mL#	61
108) 2,6,10-Trimethyltridec...	27.329	85	1253	23.245	µg/mL	99
109) 2,6,10-Trimethylpentad...	31.054	85	805m	14.934	µg/mL	

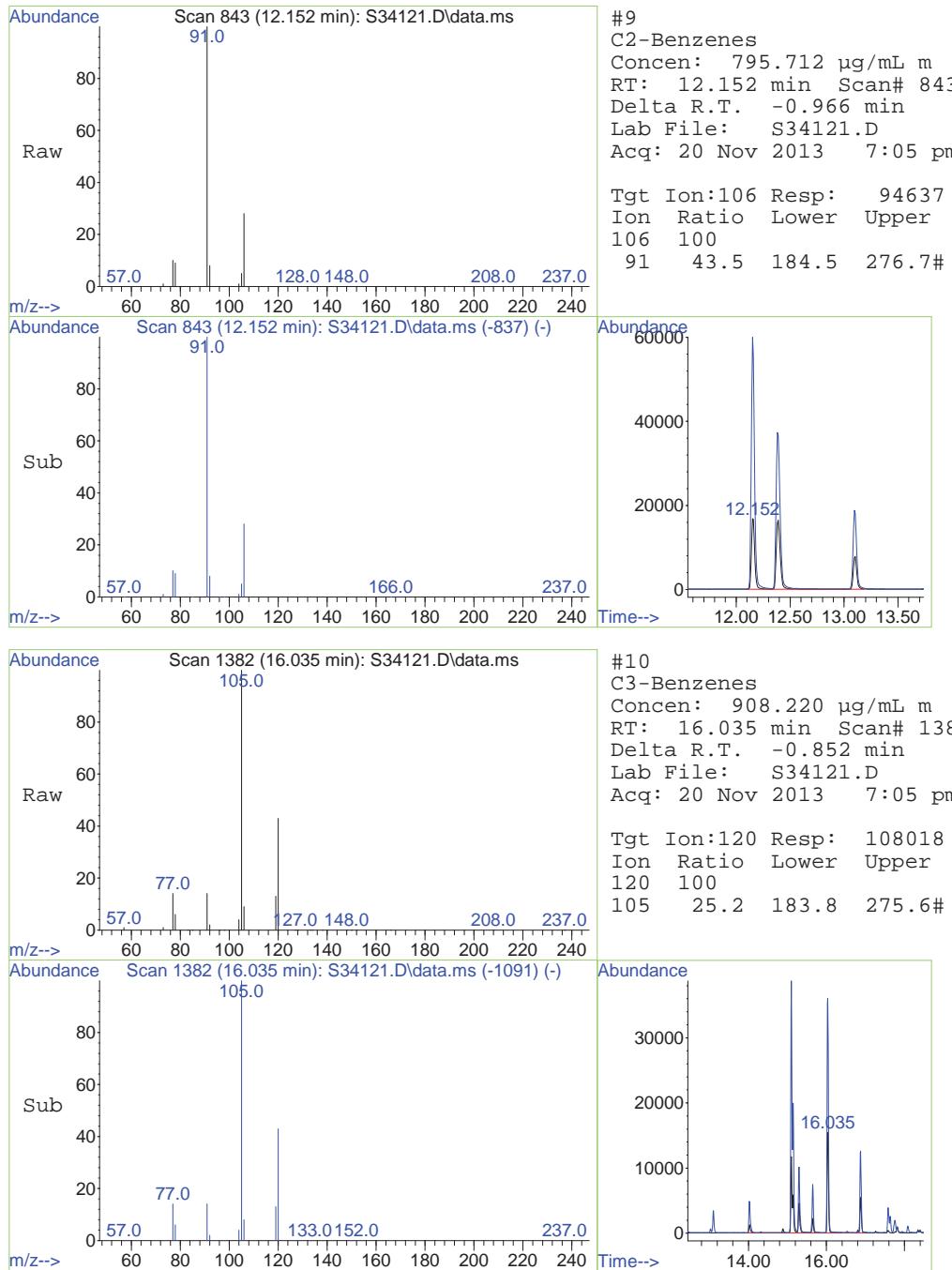
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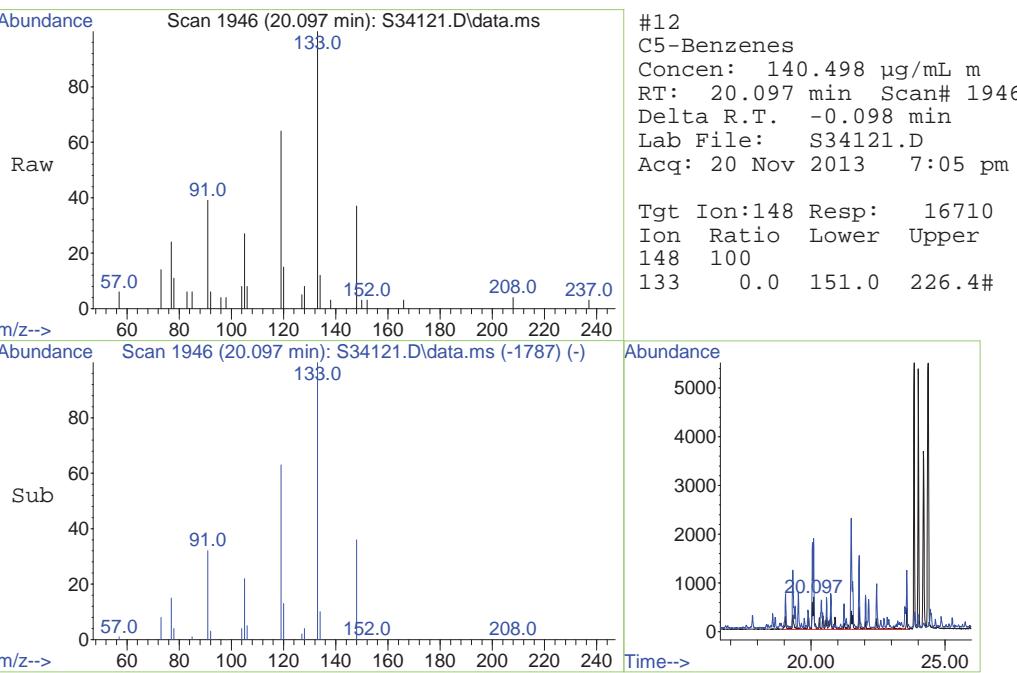
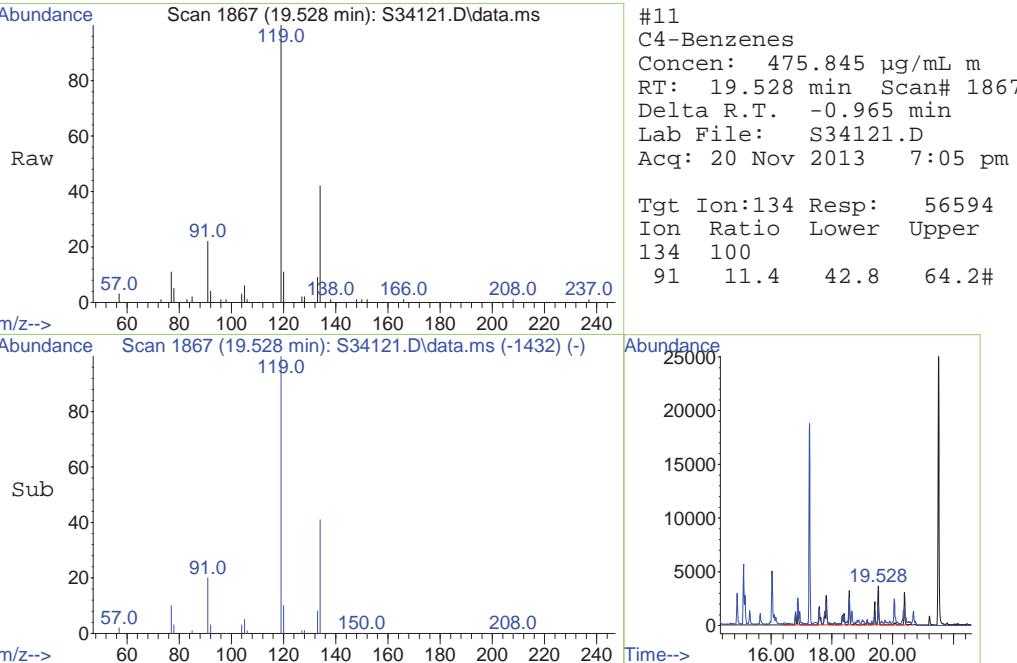


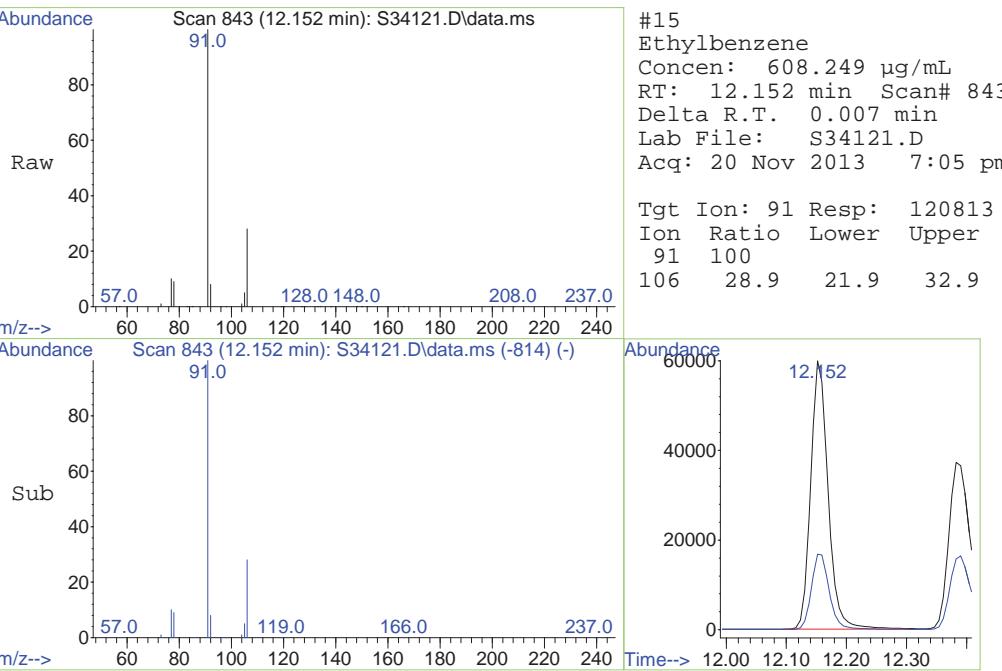
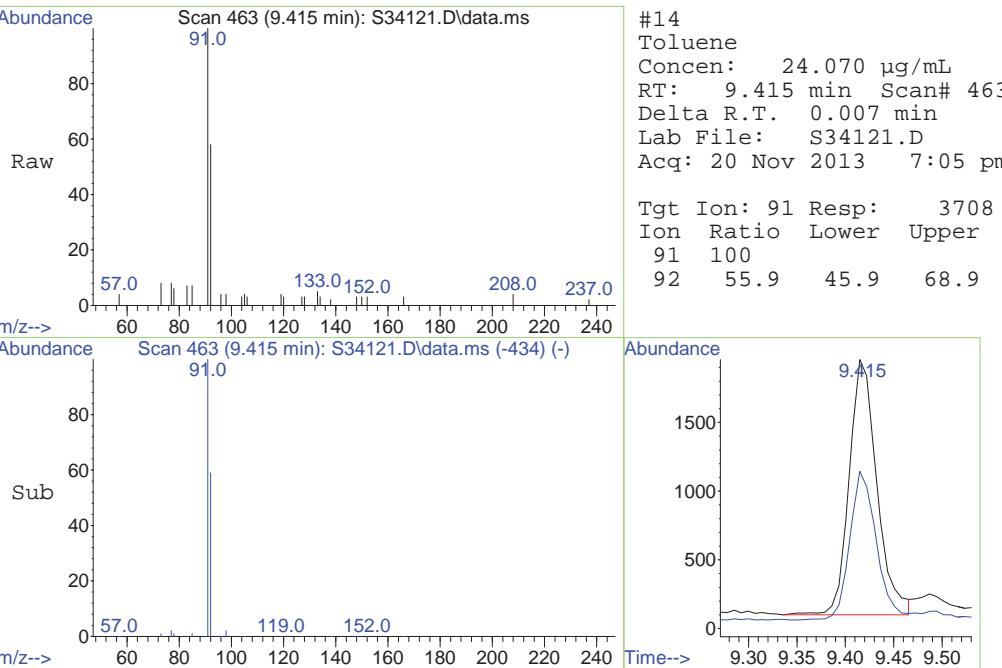
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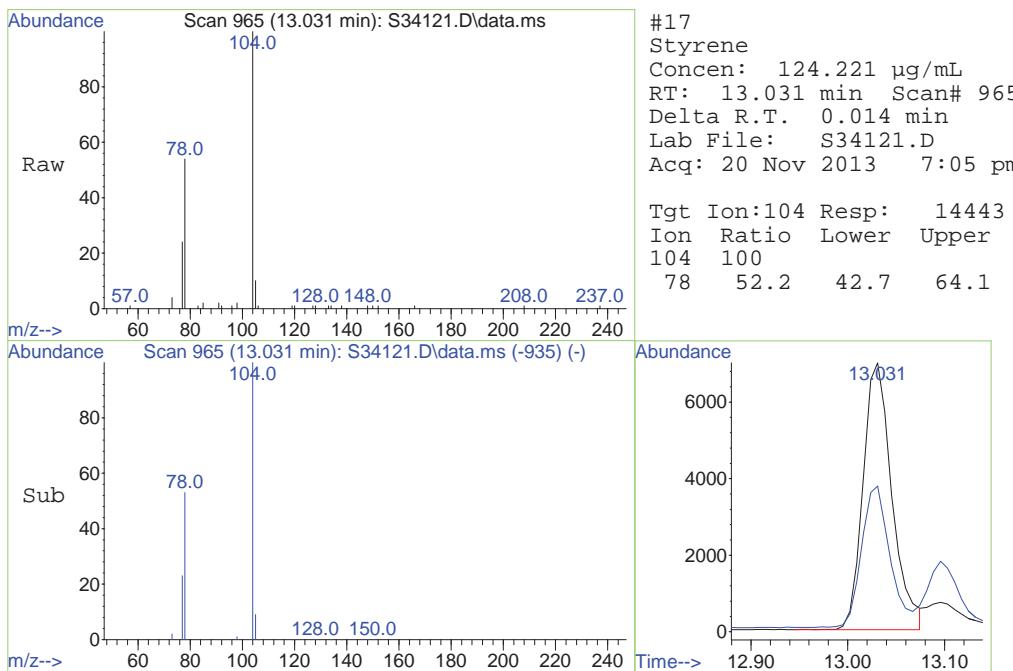
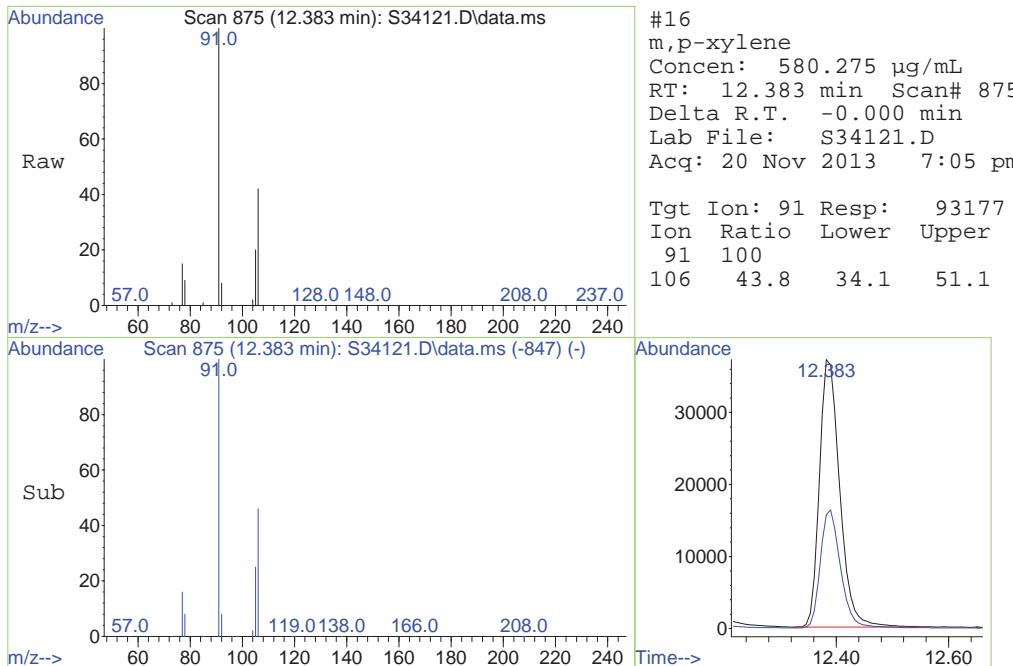
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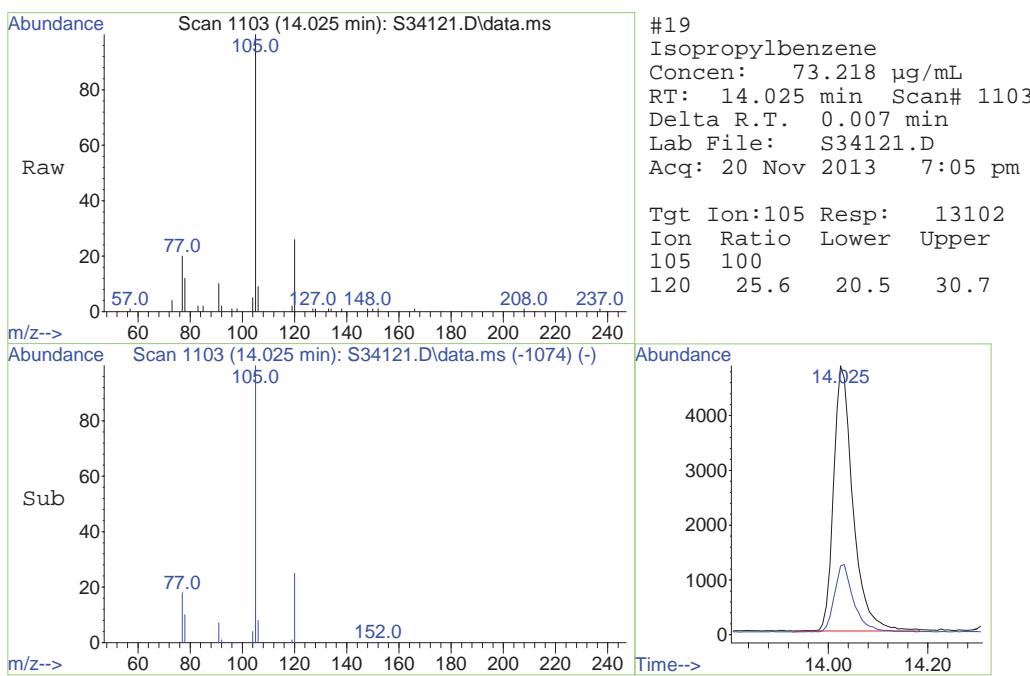
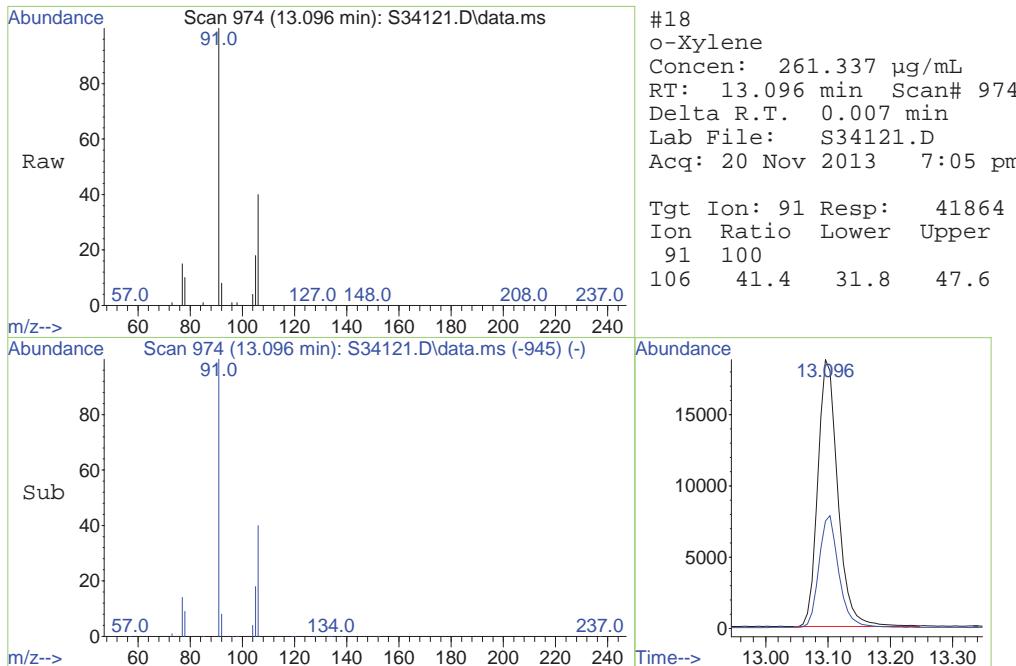


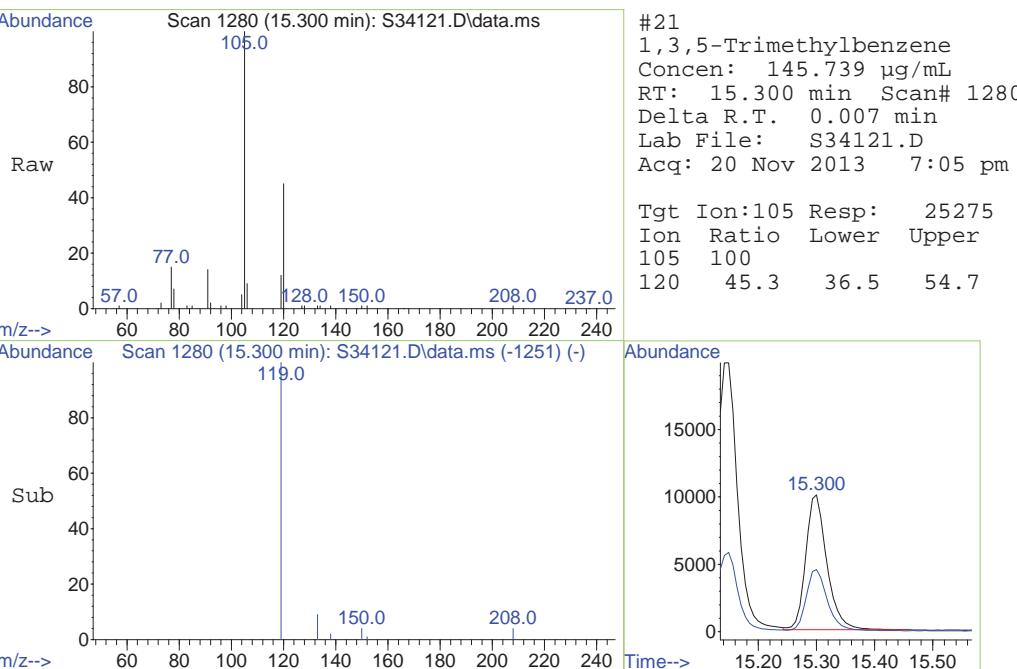
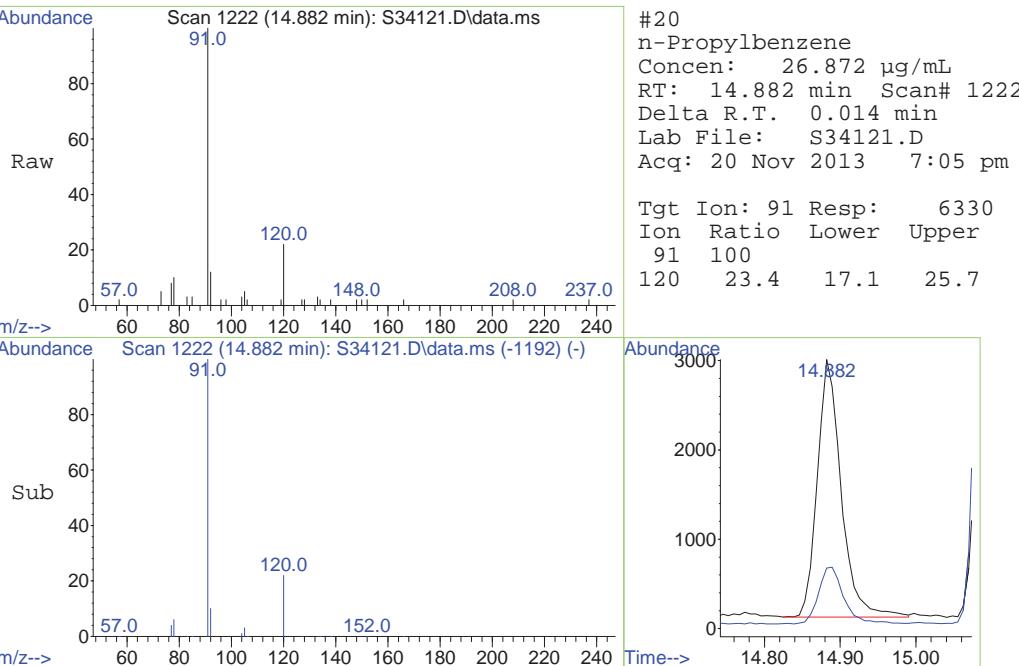


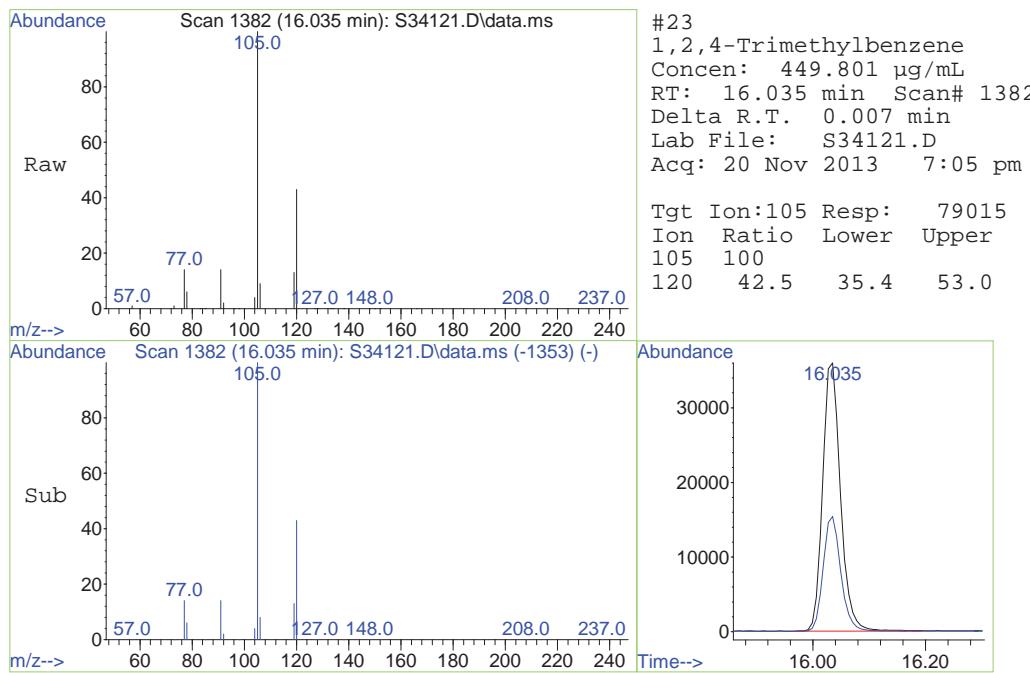
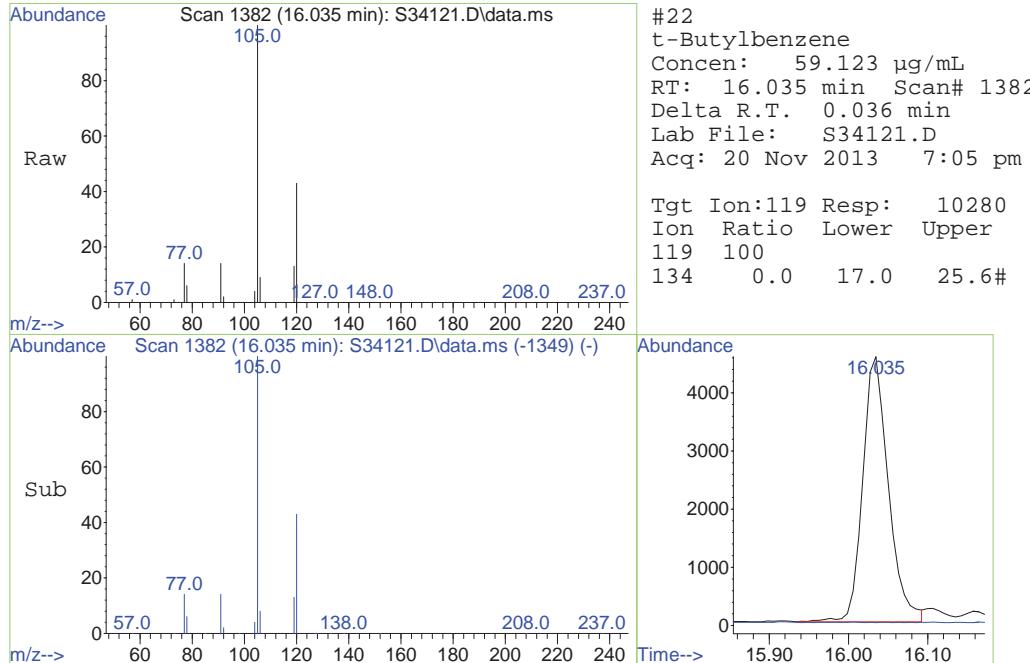


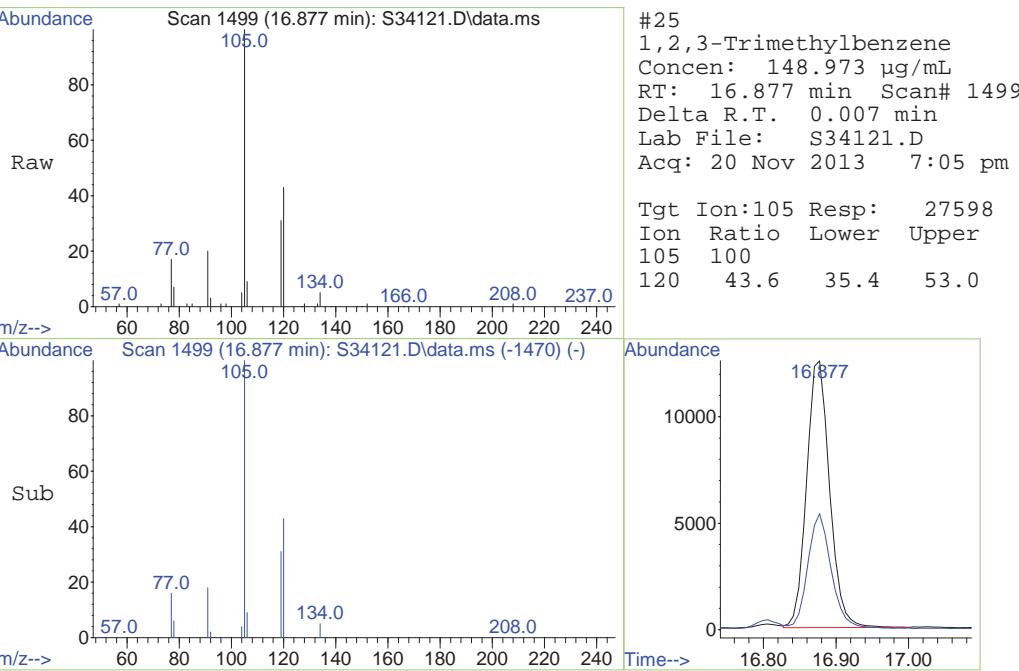
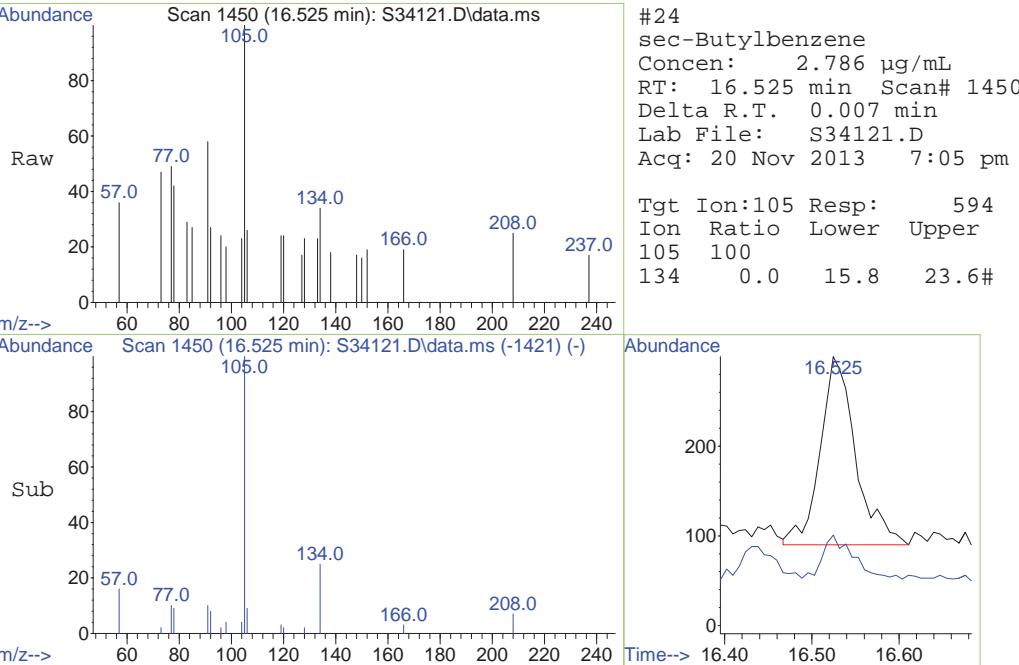


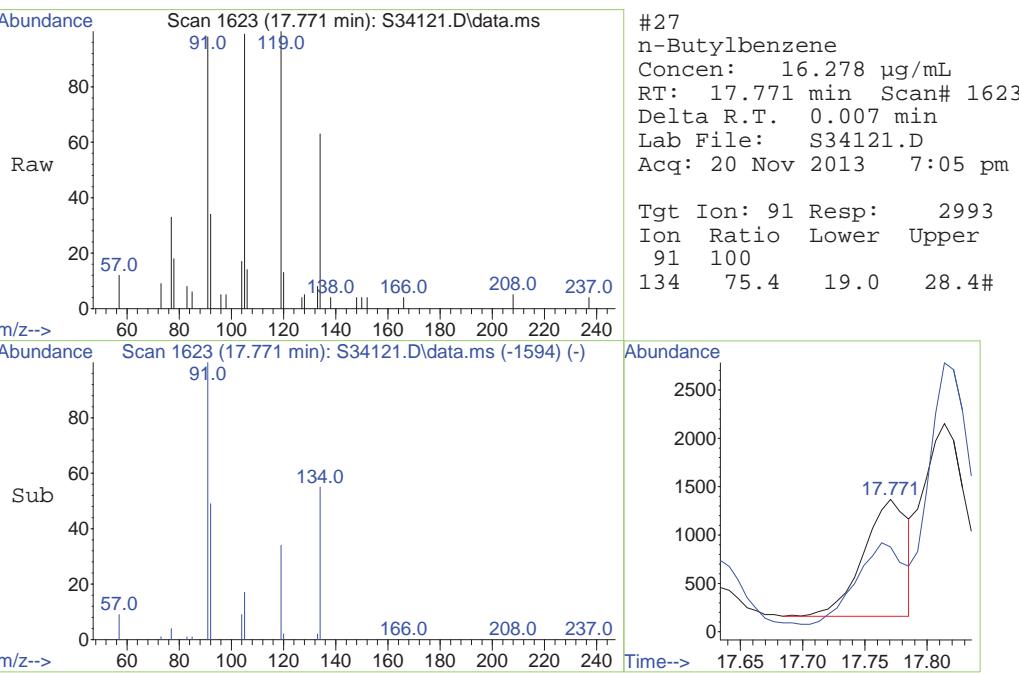
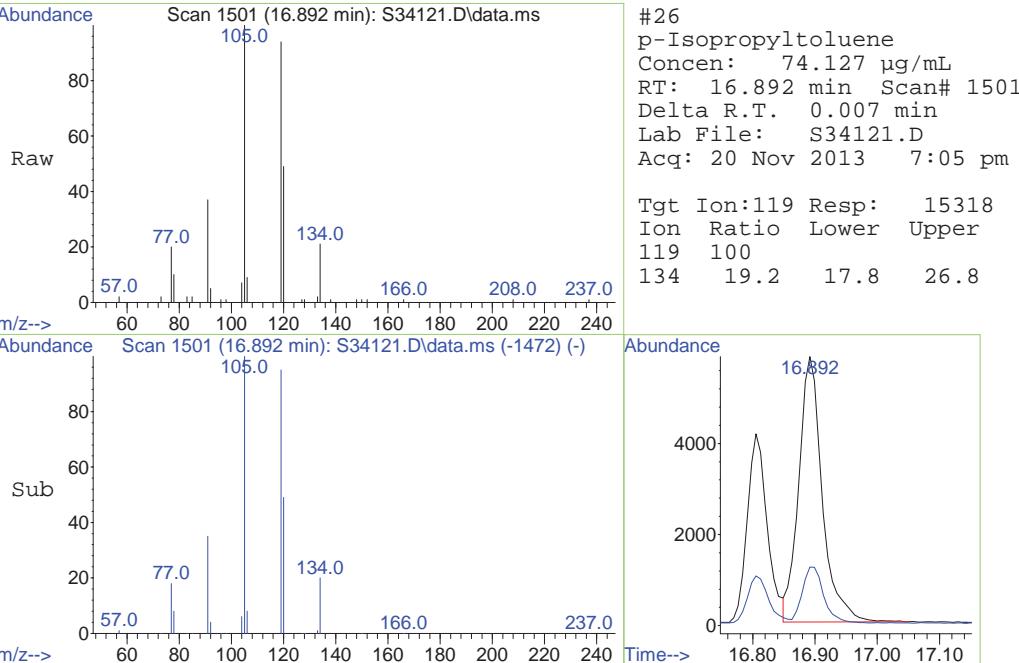


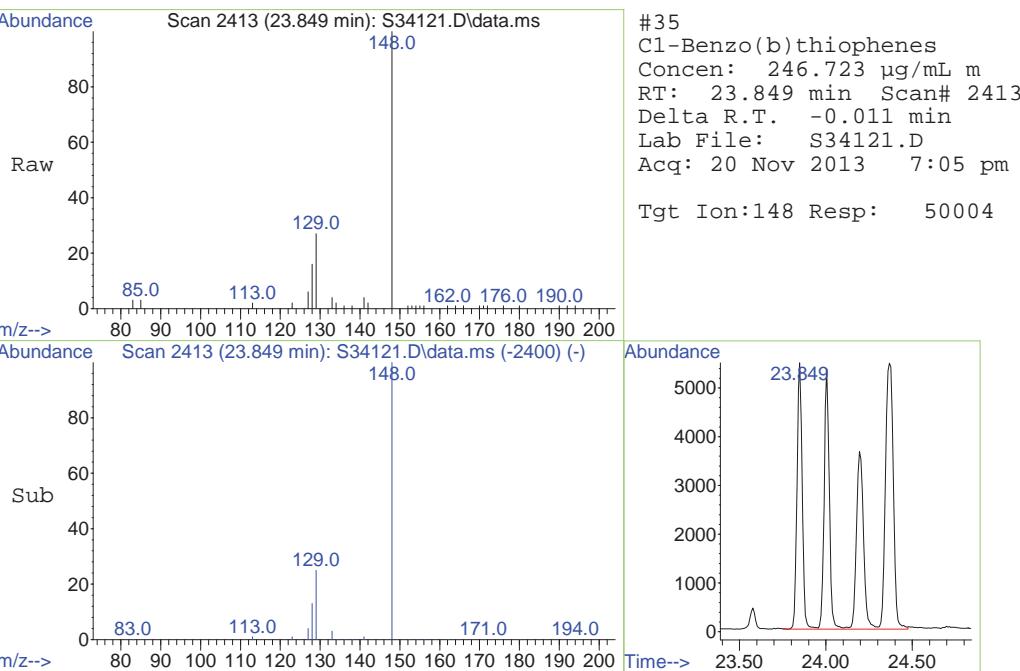
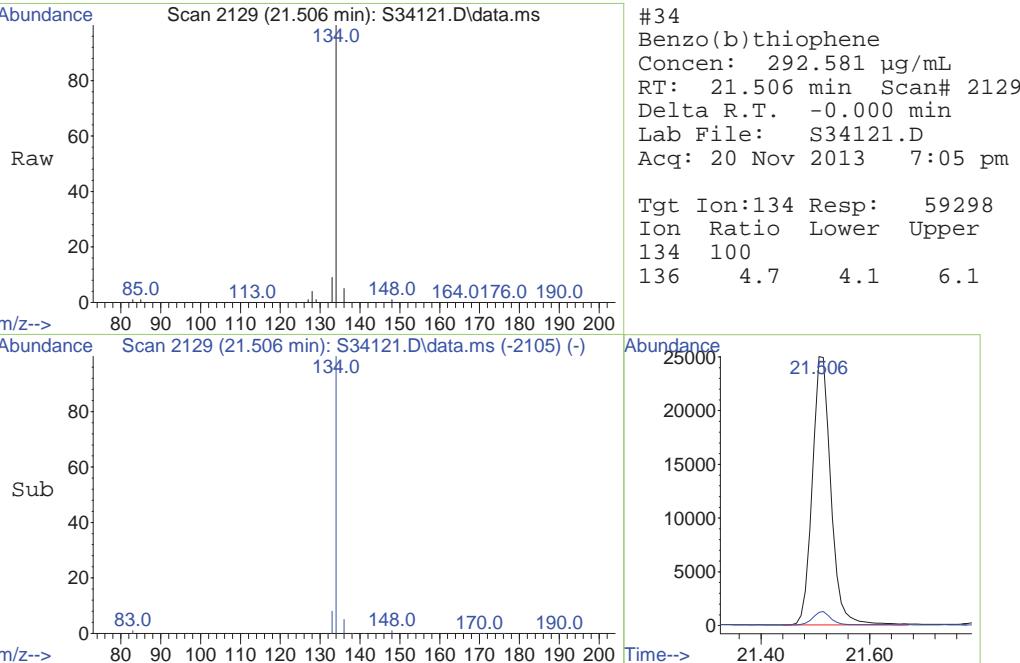


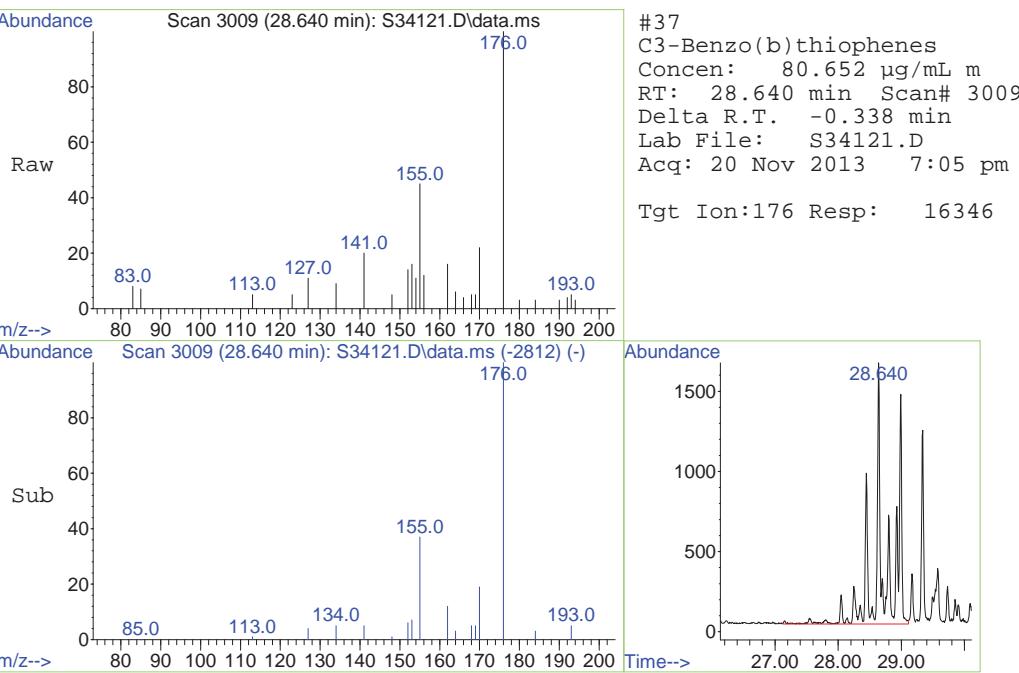
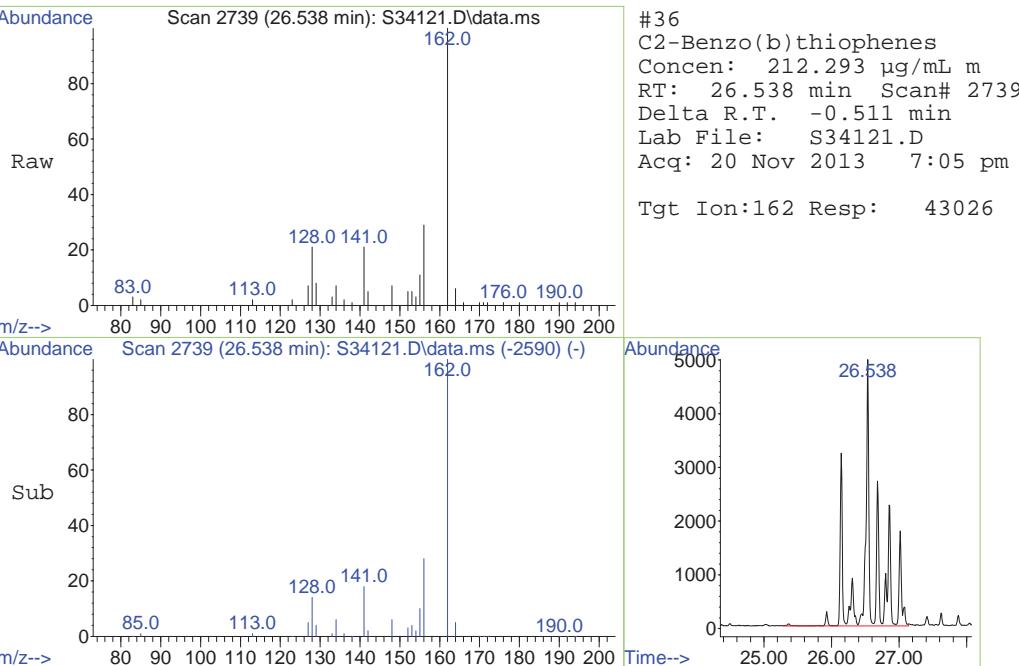


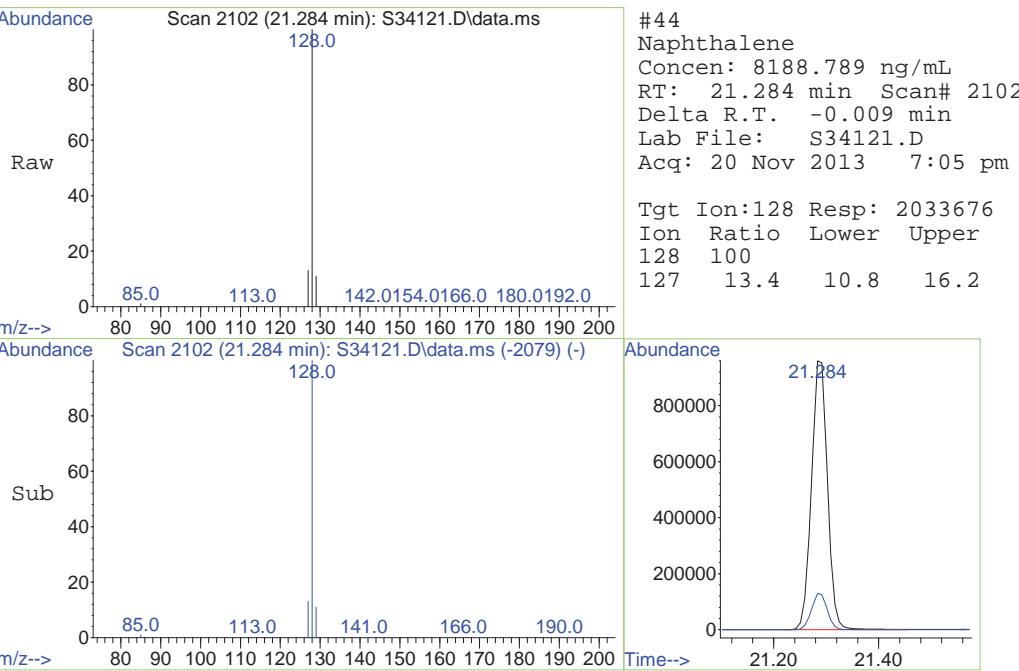
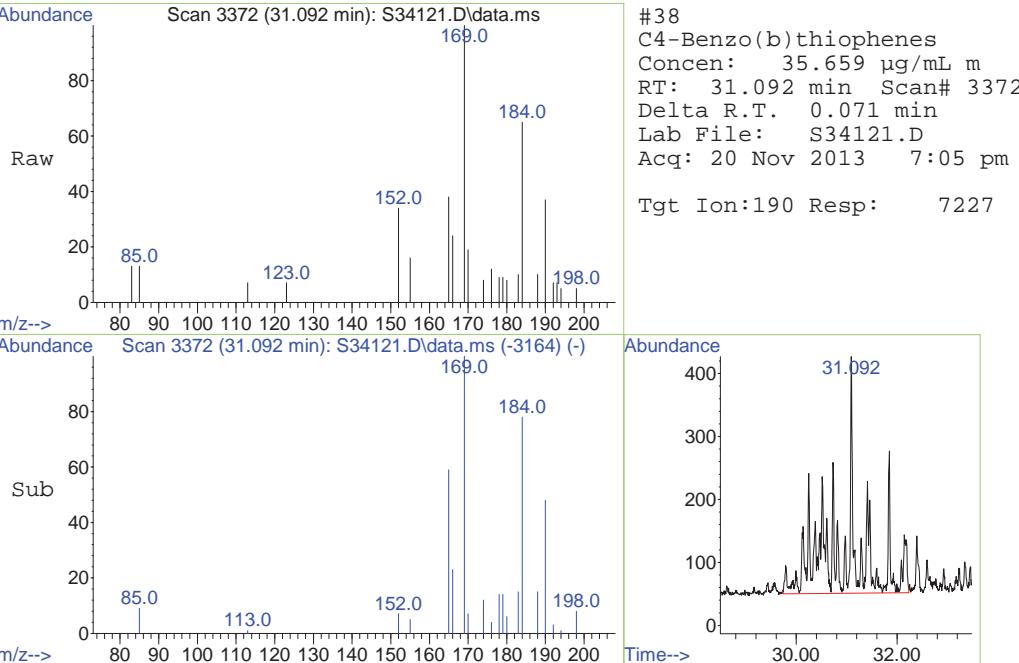


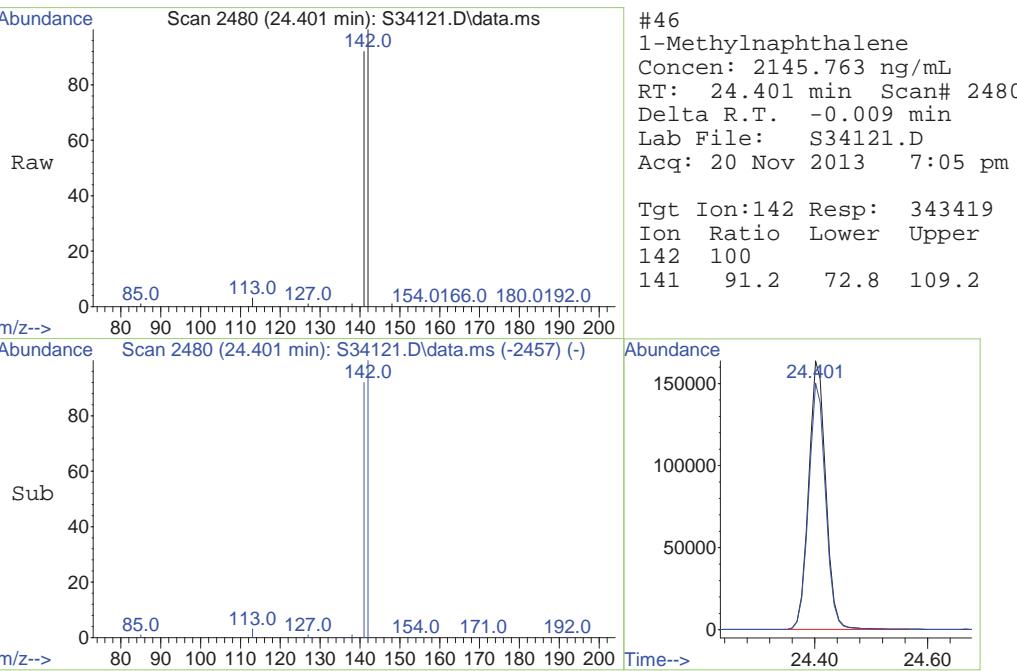
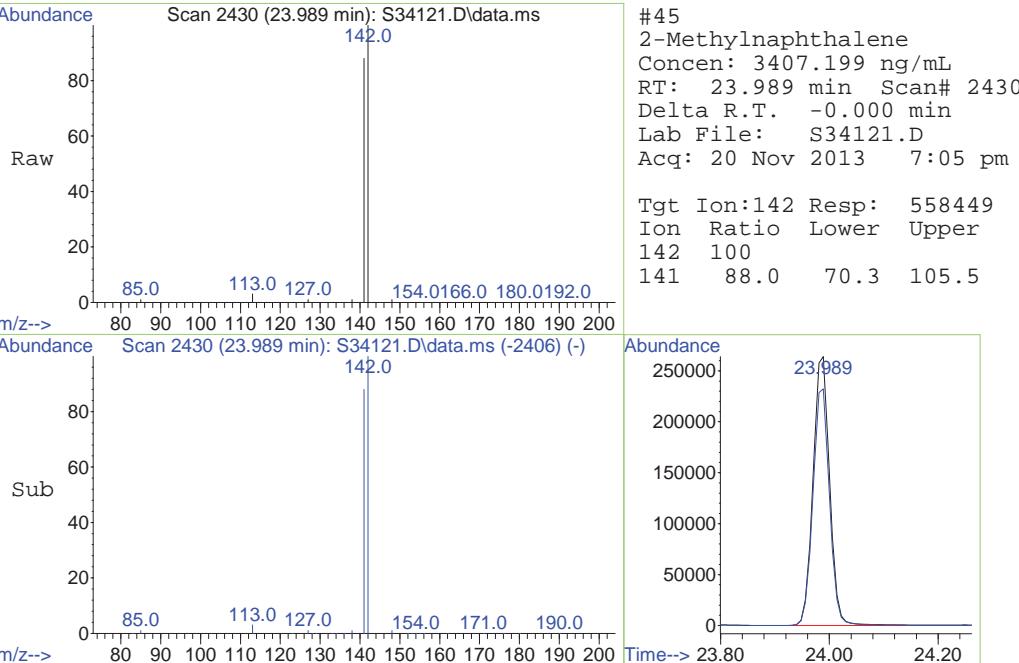


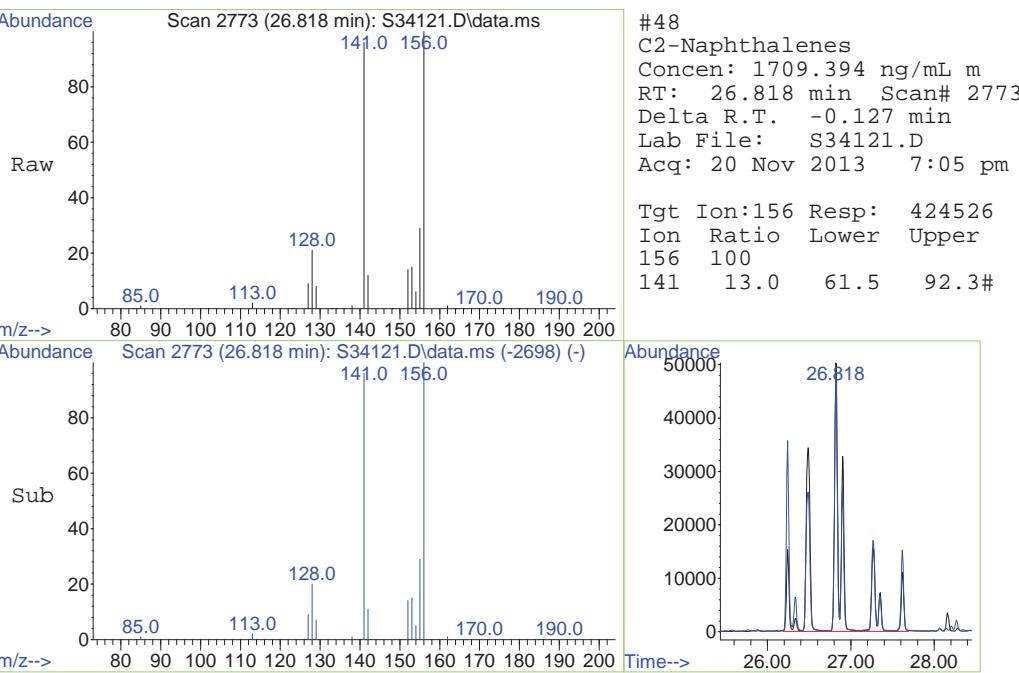
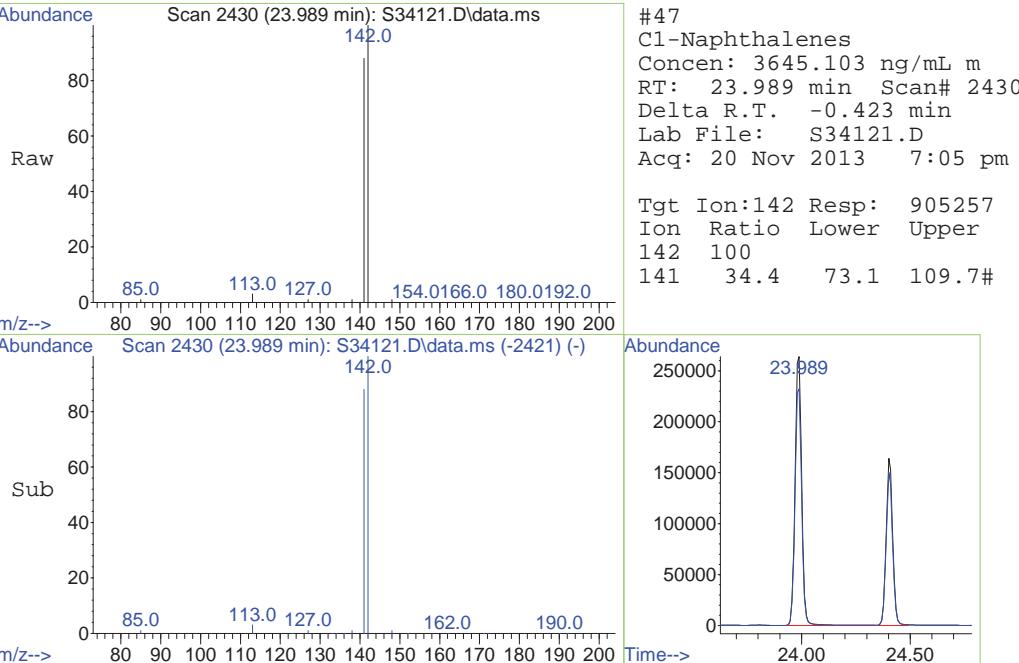


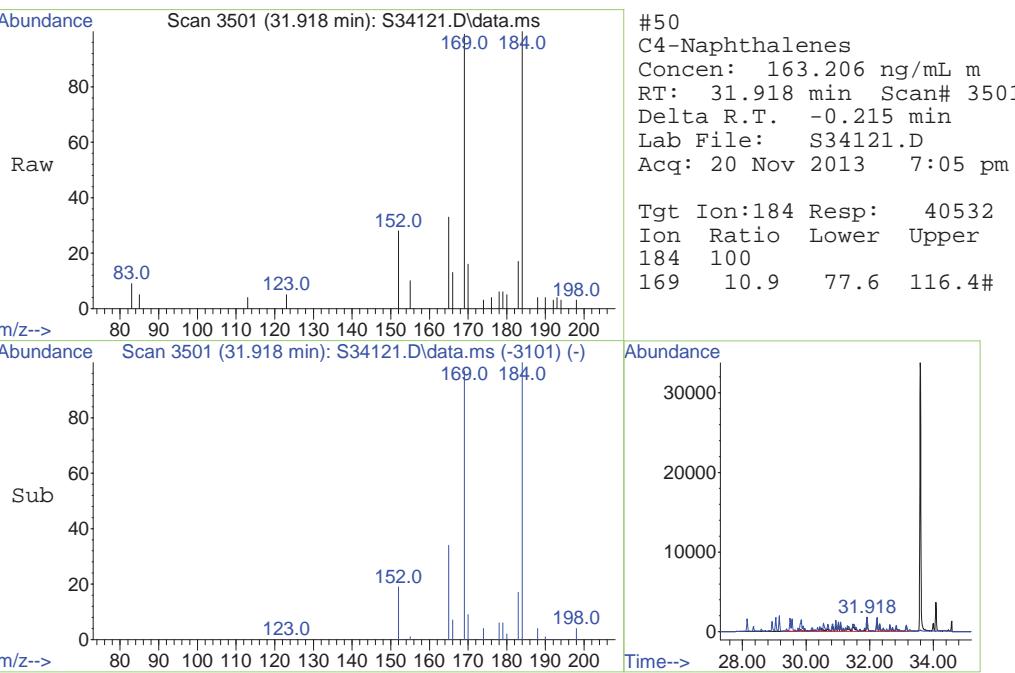
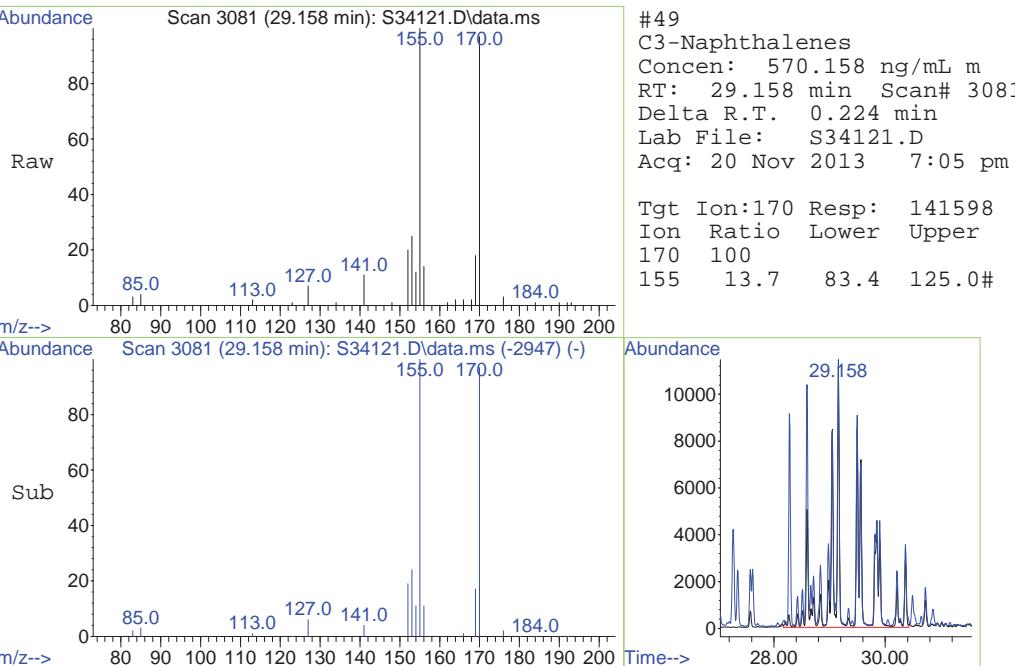


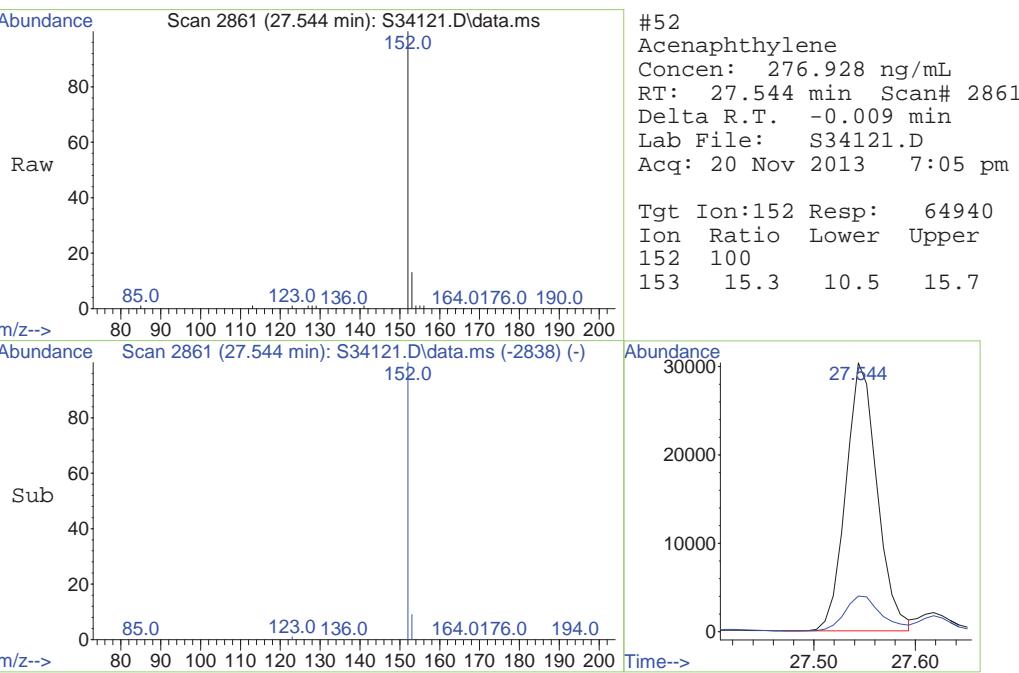
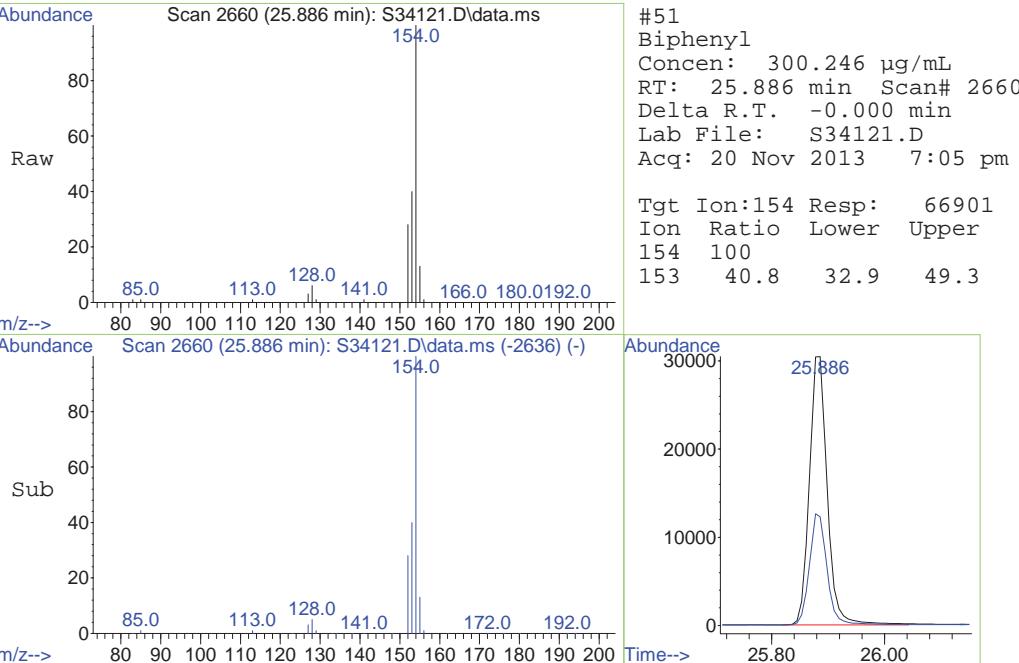


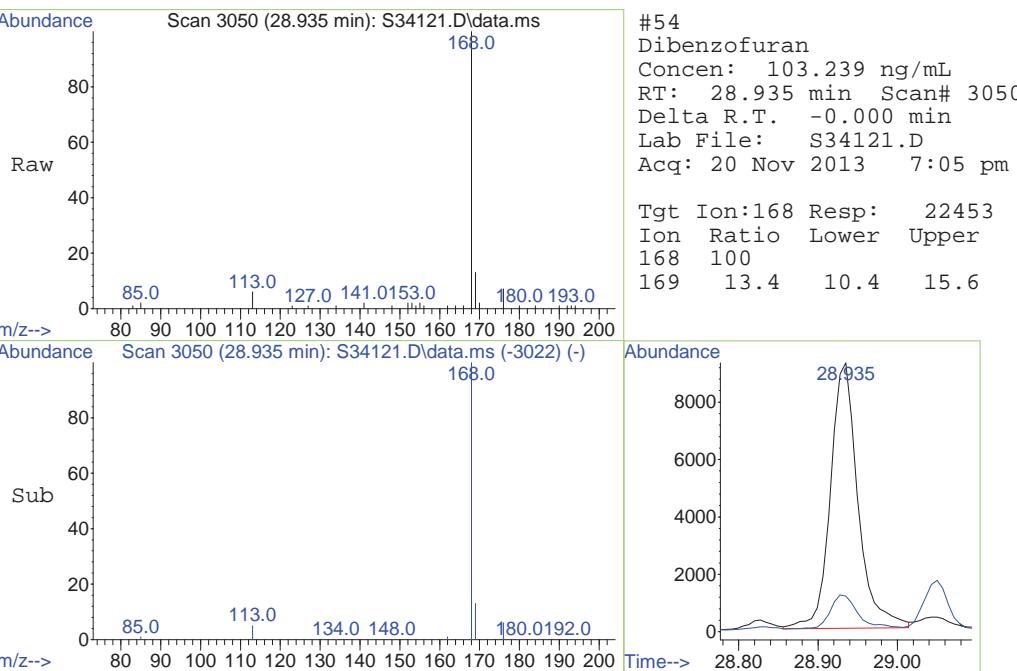
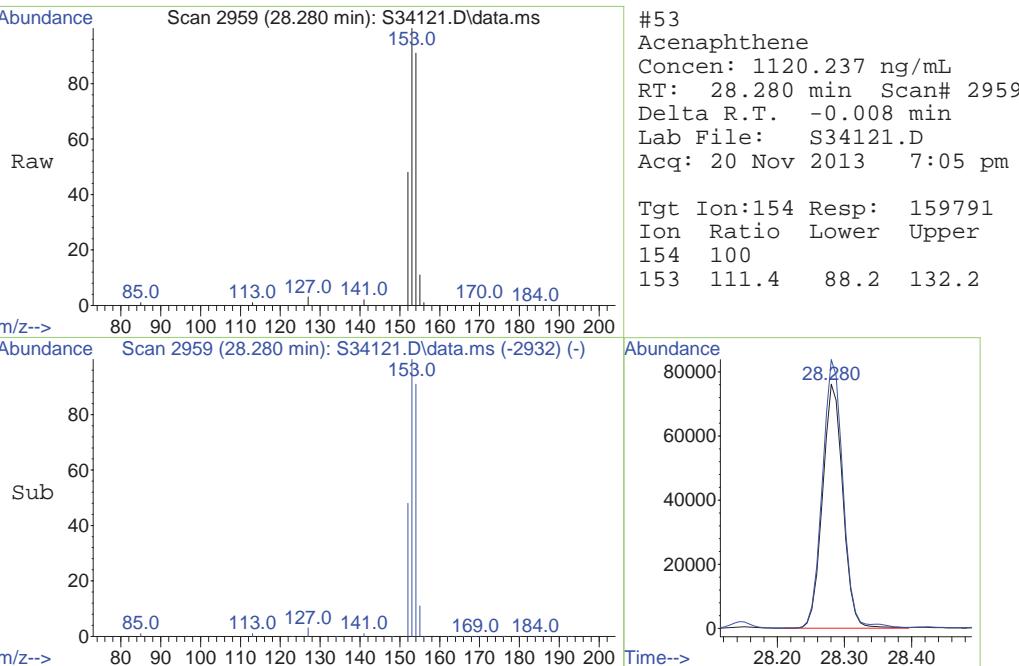


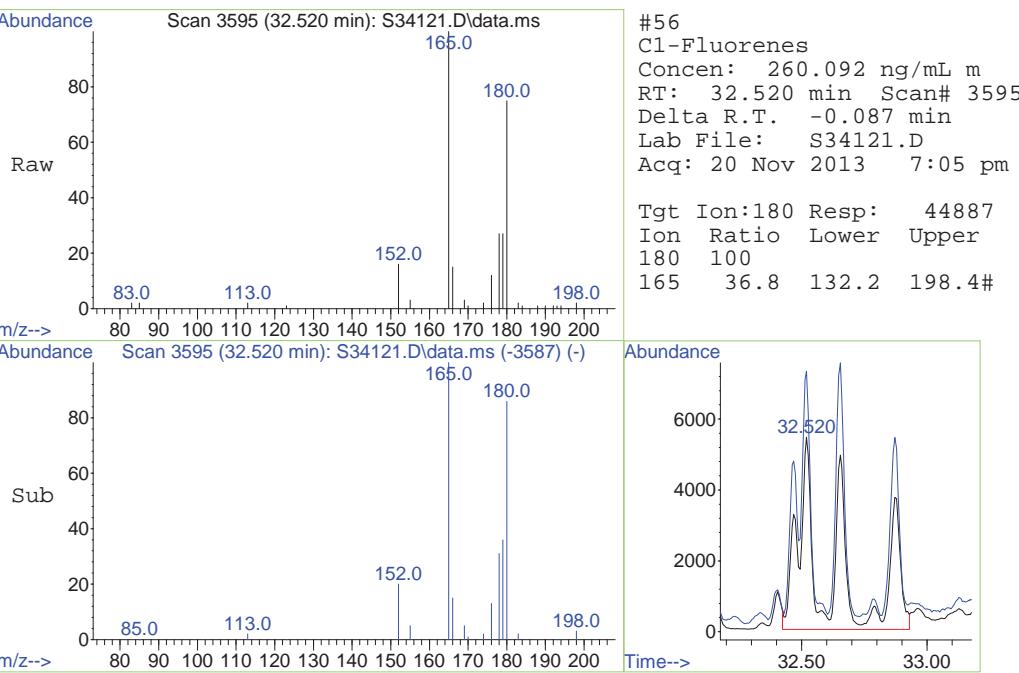
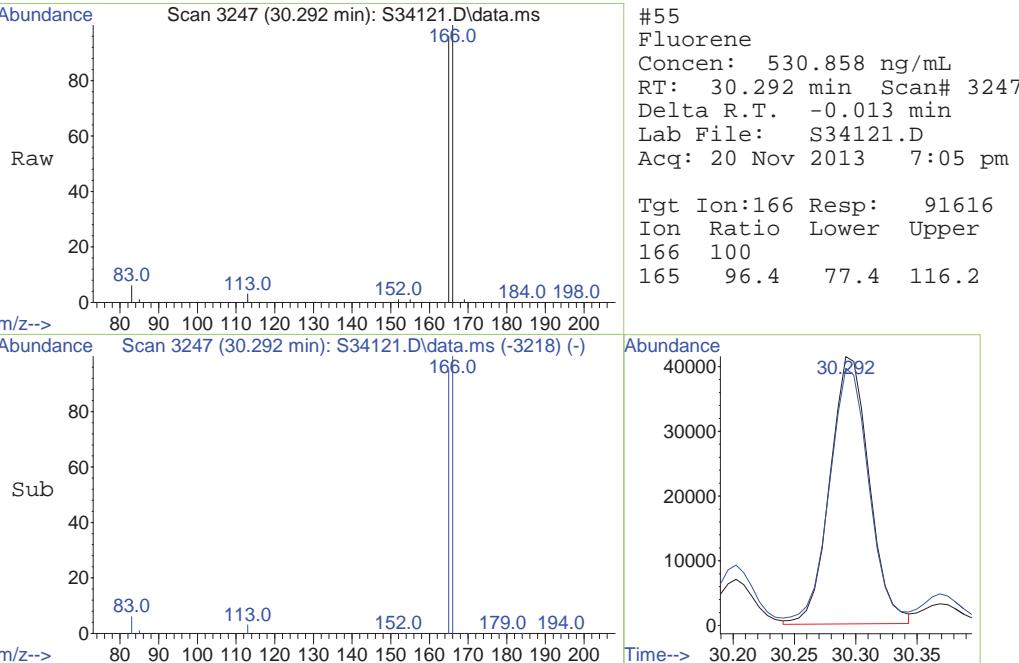


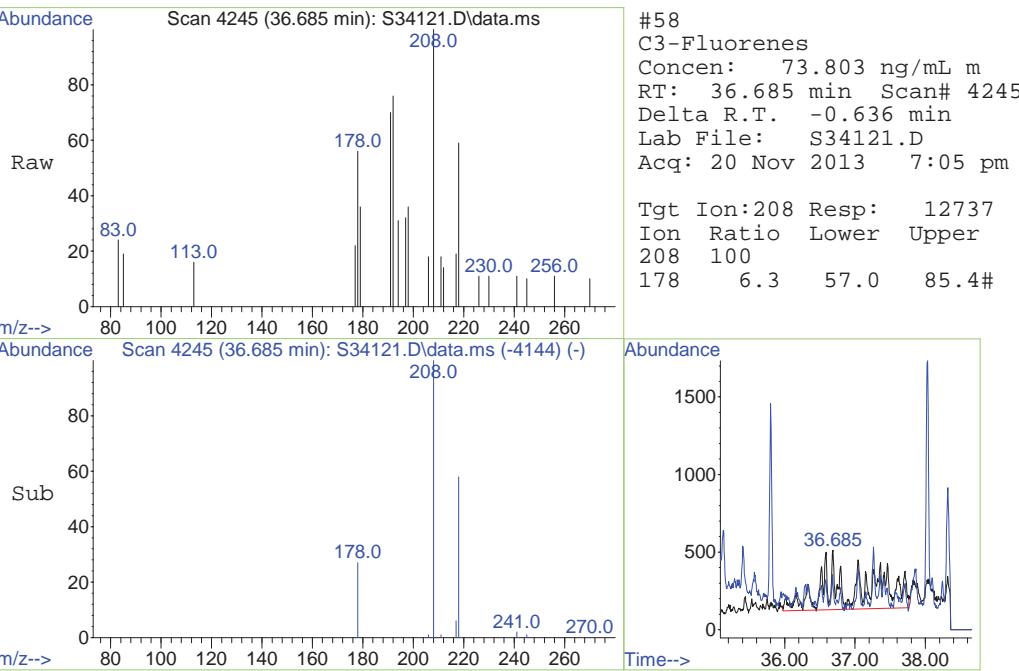
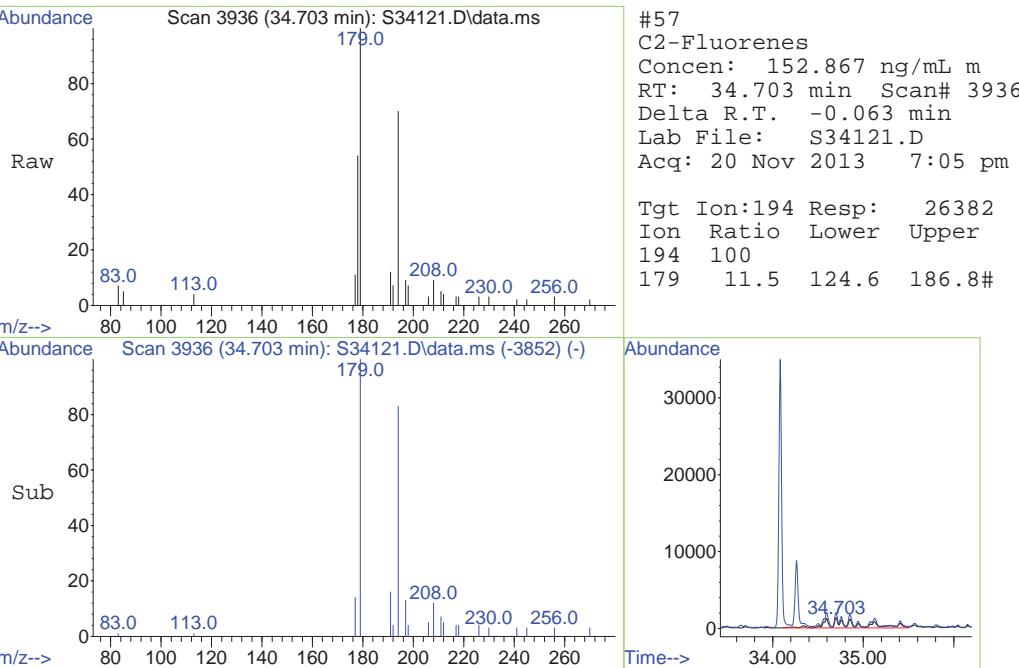


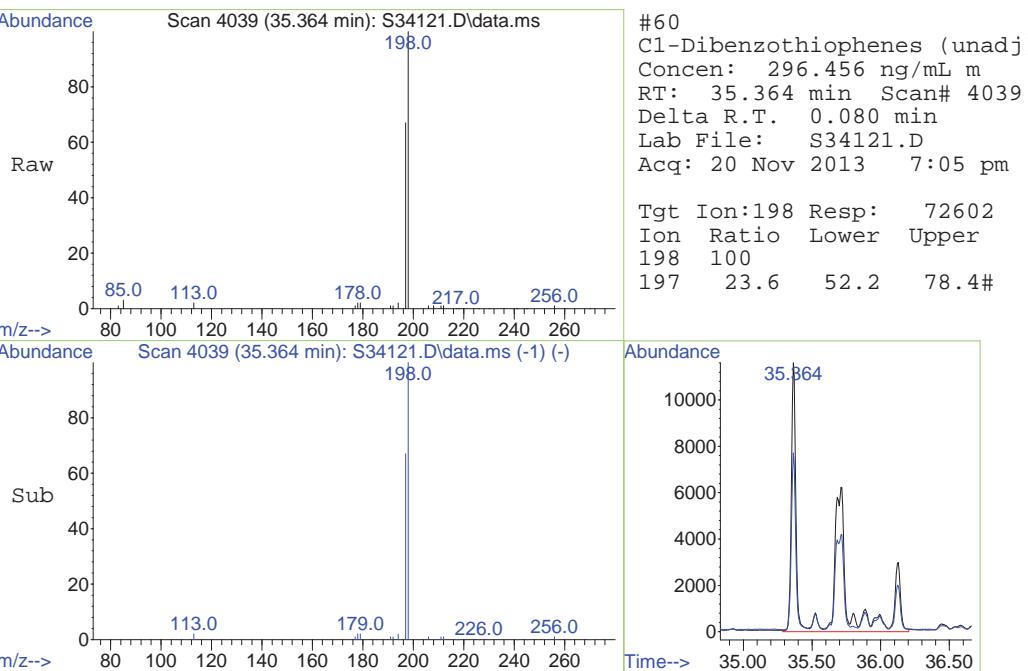
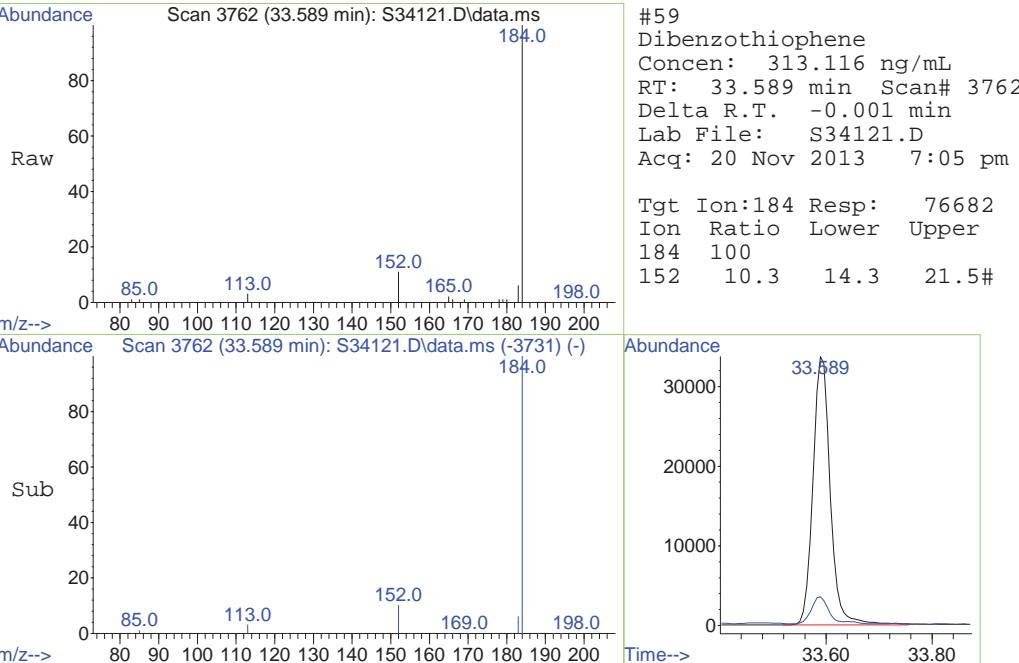


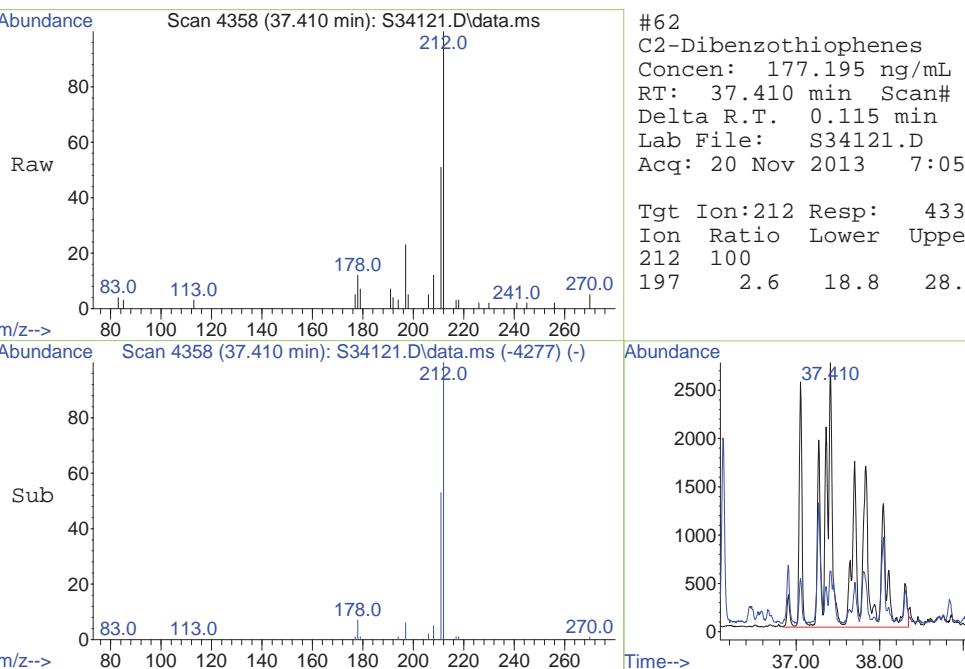
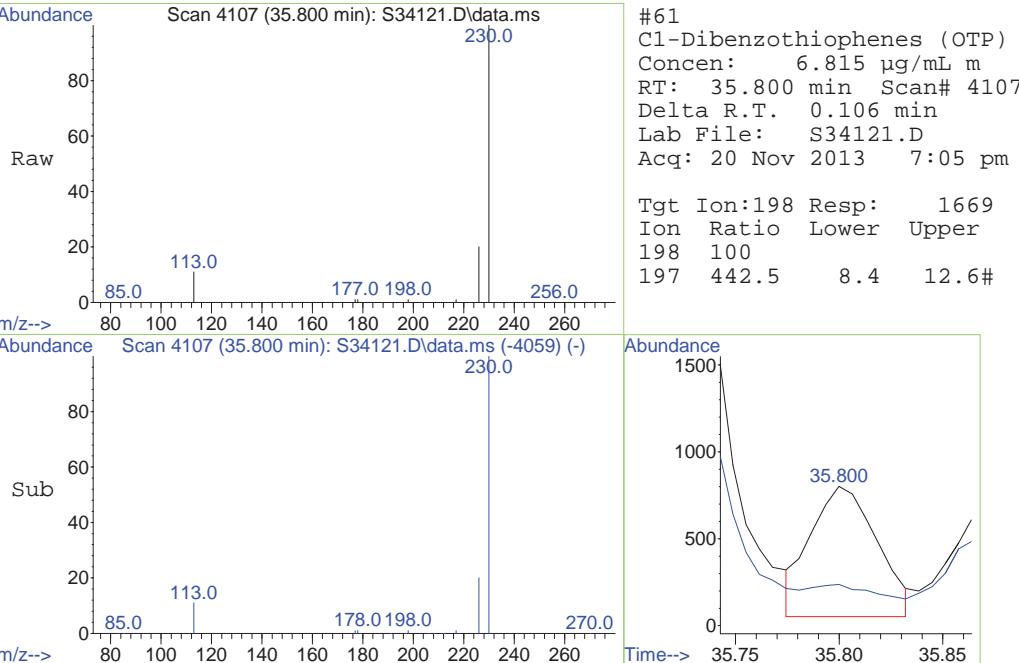


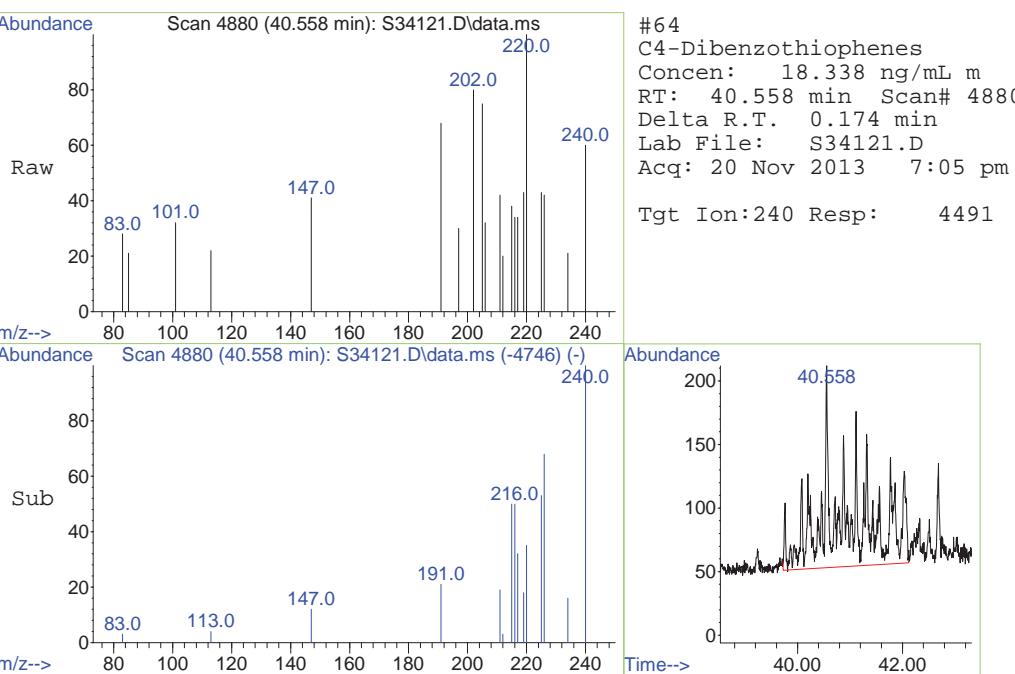
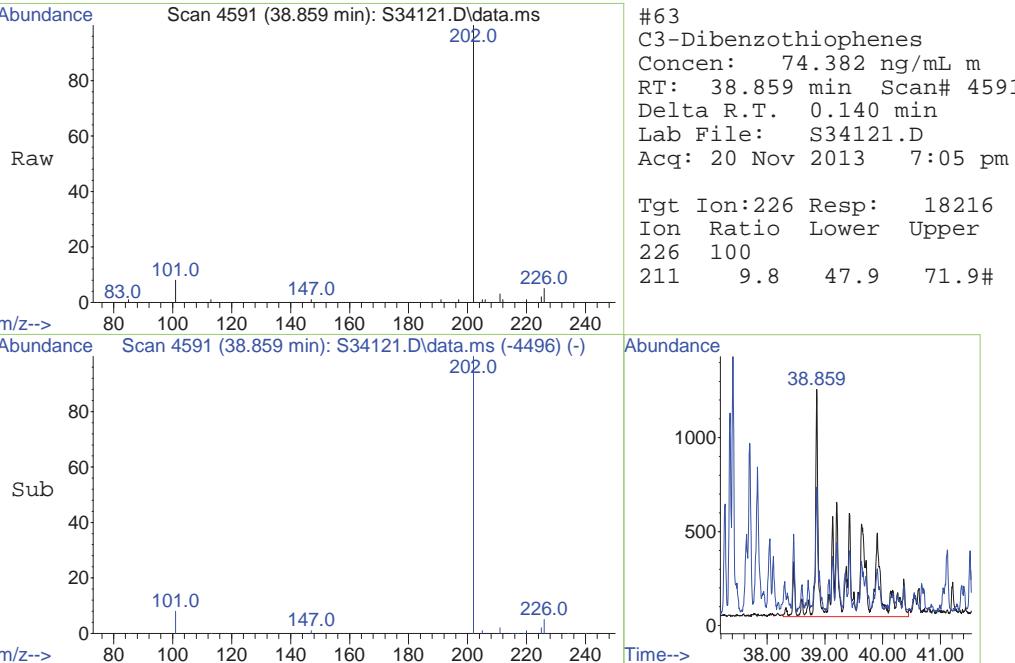


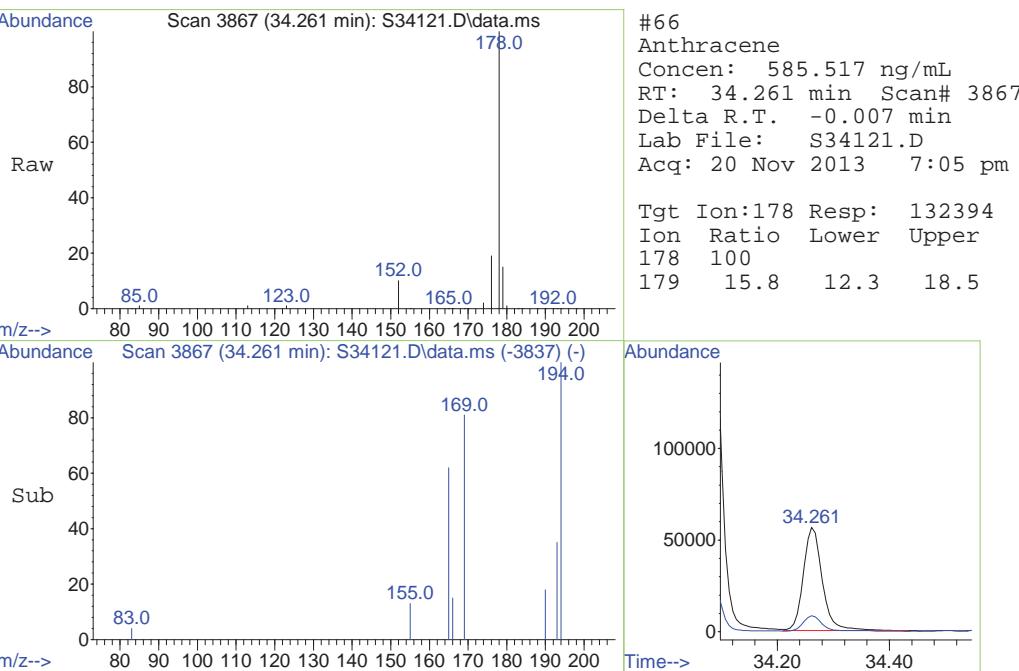
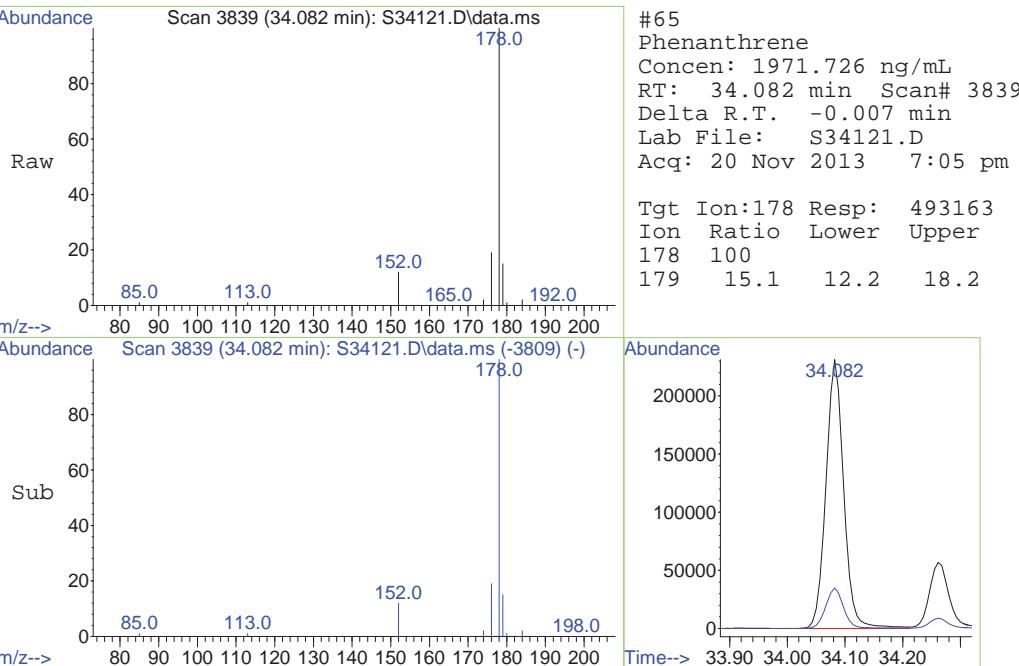


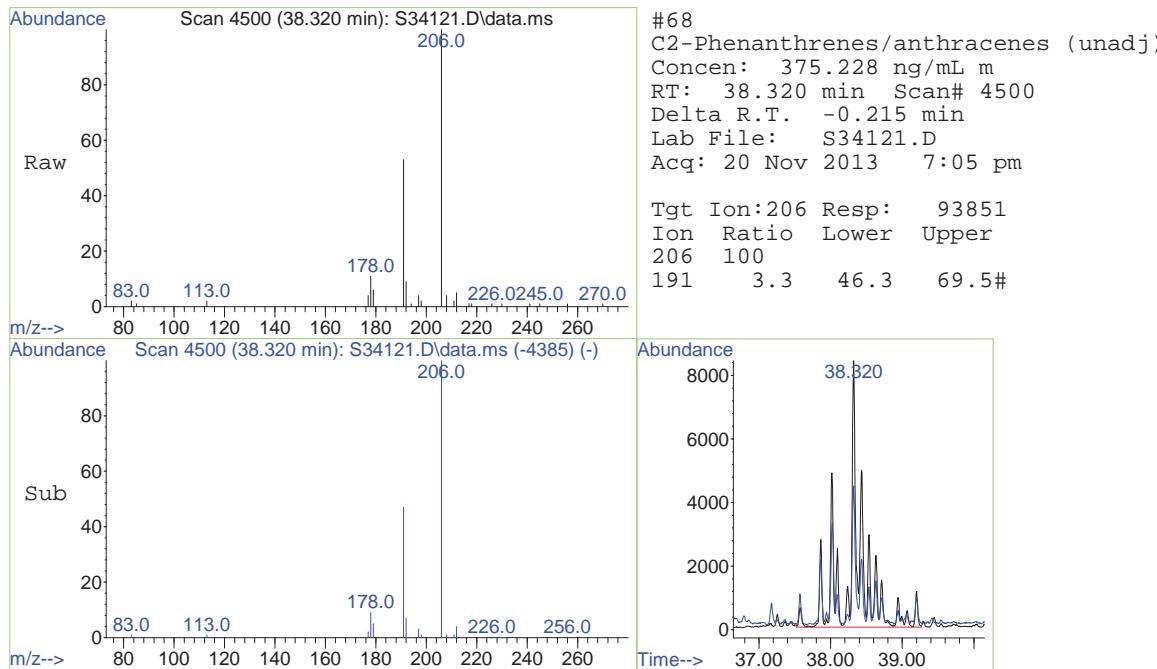
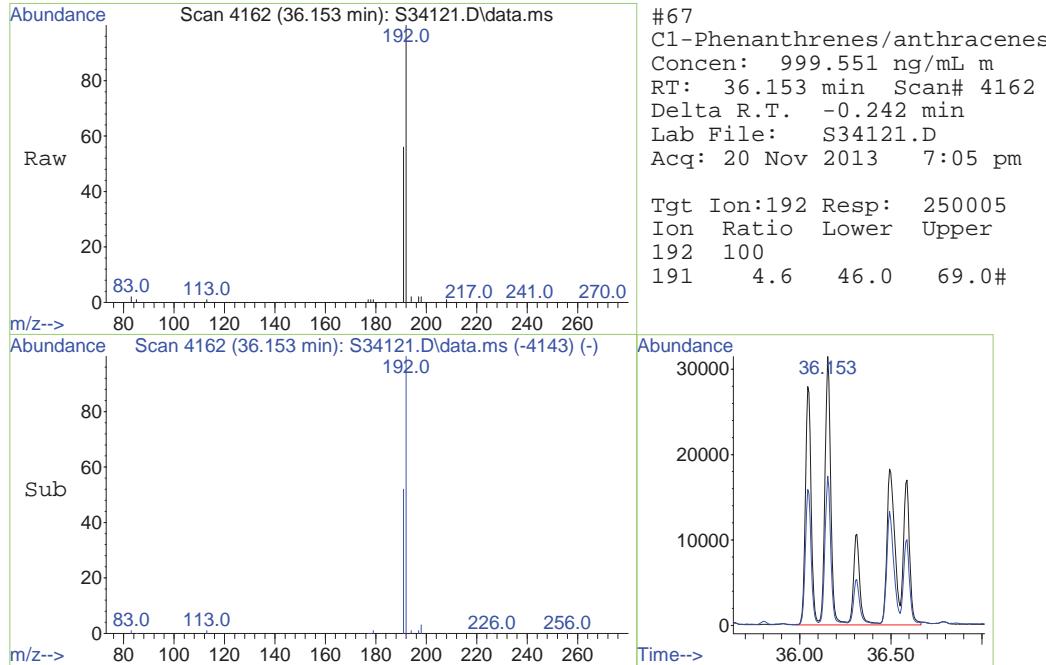




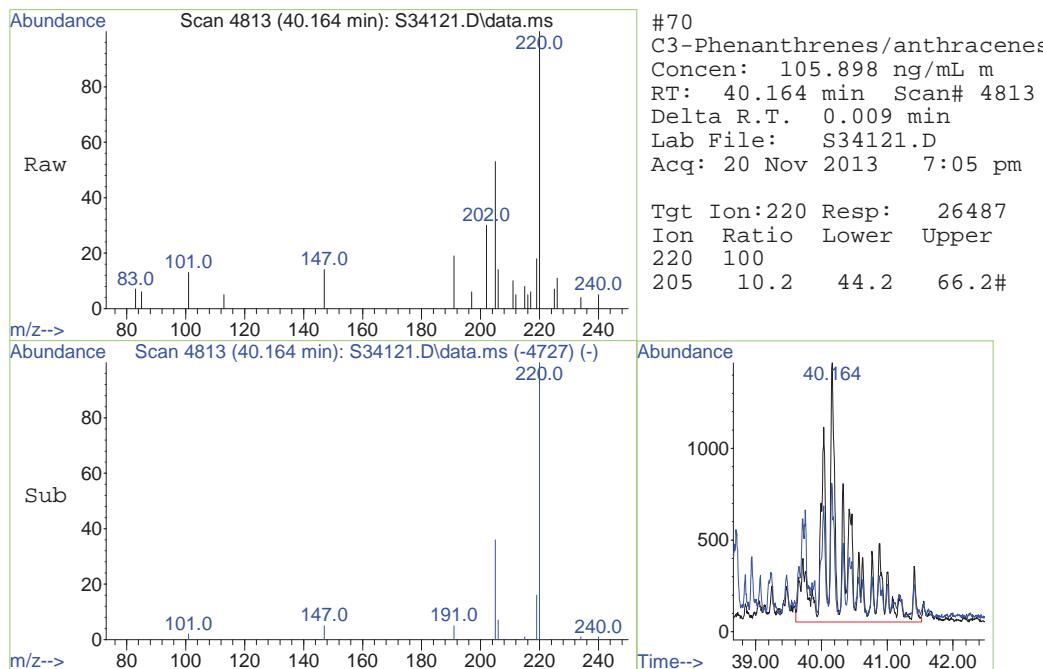
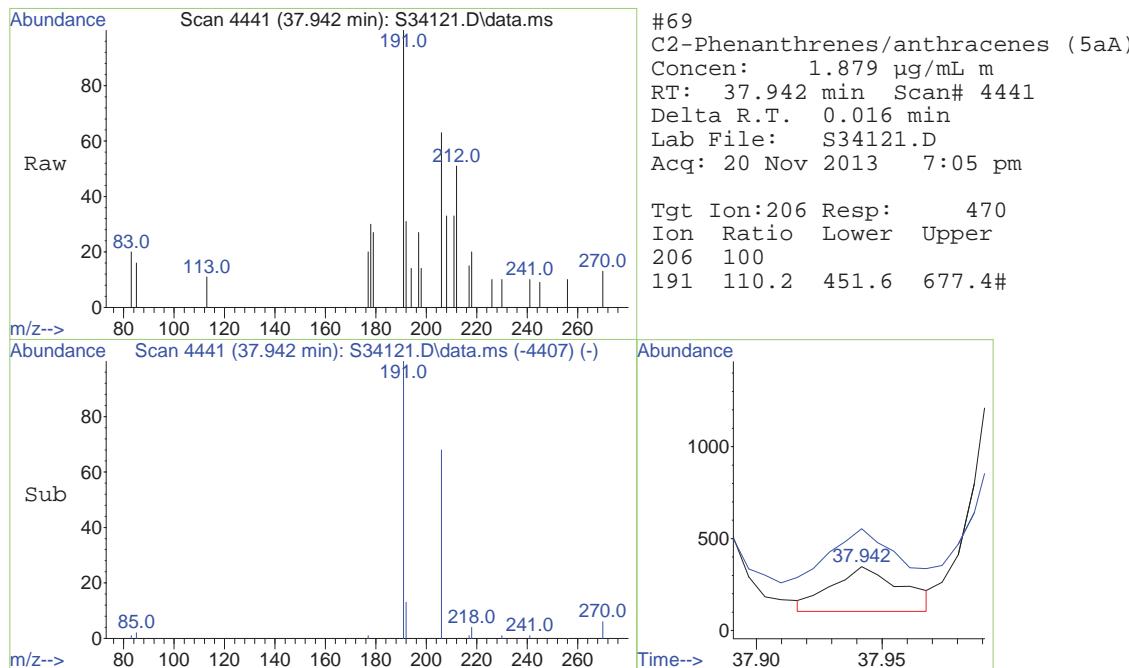


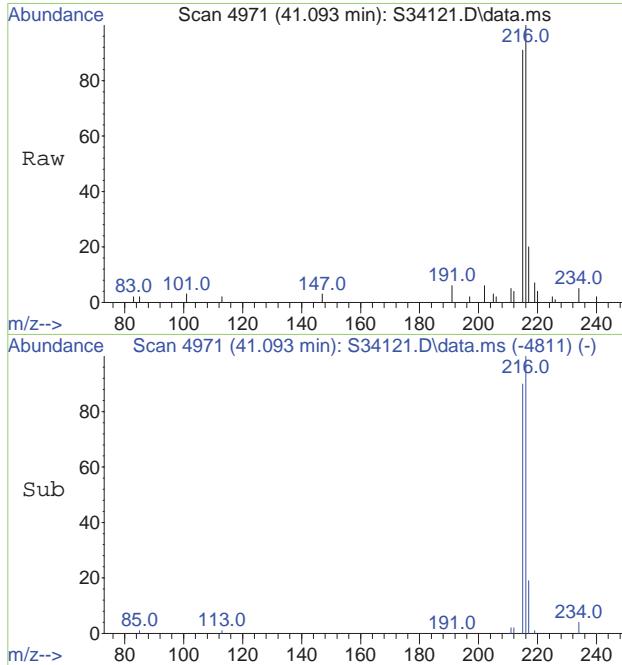






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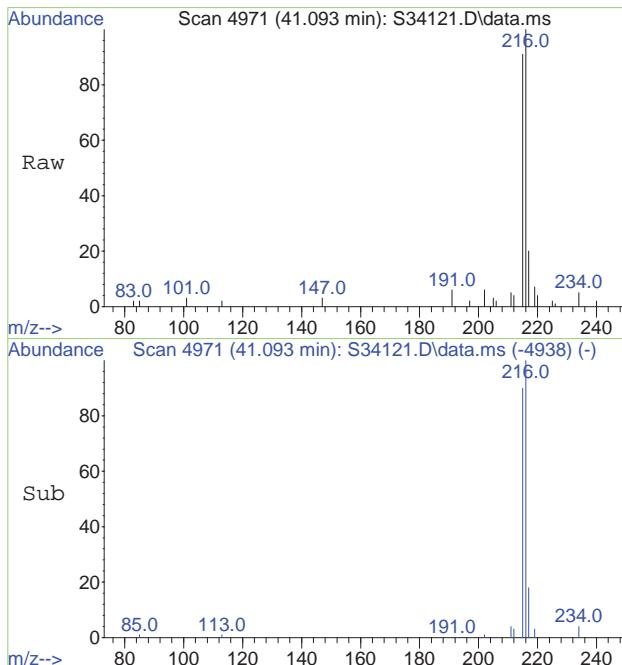


#71  
C4-Phenanthrenes/anthracenes  
Concen: 28.691 ng/mL m  
RT: 41.093 min Scan# 4971  
Delta R.T. -0.530 min  
Lab File: S34121.D  
Acq: 20 Nov 2013 7:05 pm  
  
Tgt Ion: 234 Resp: 7176  
Ion Ratio Lower Upper  
234 100  
219 0.0 49.4 74.0#

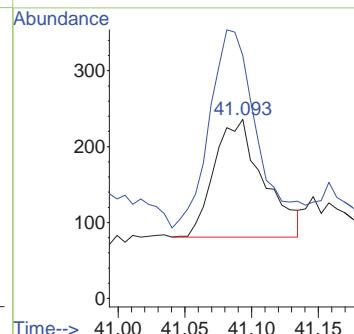
100 of 133

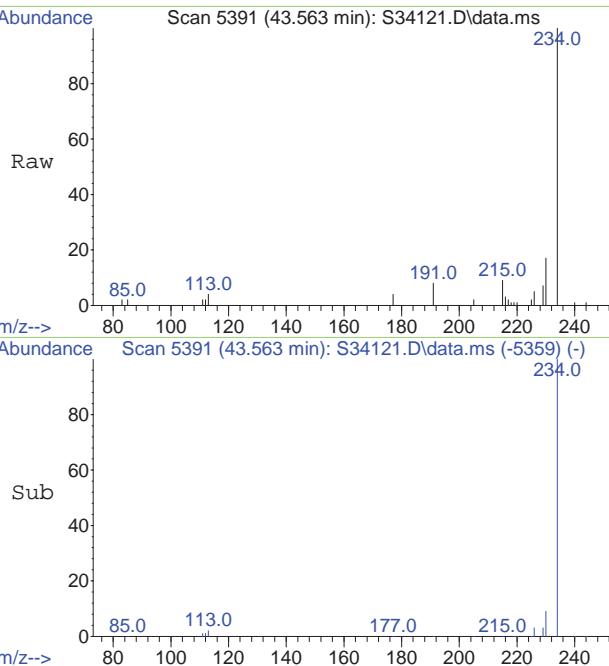
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7

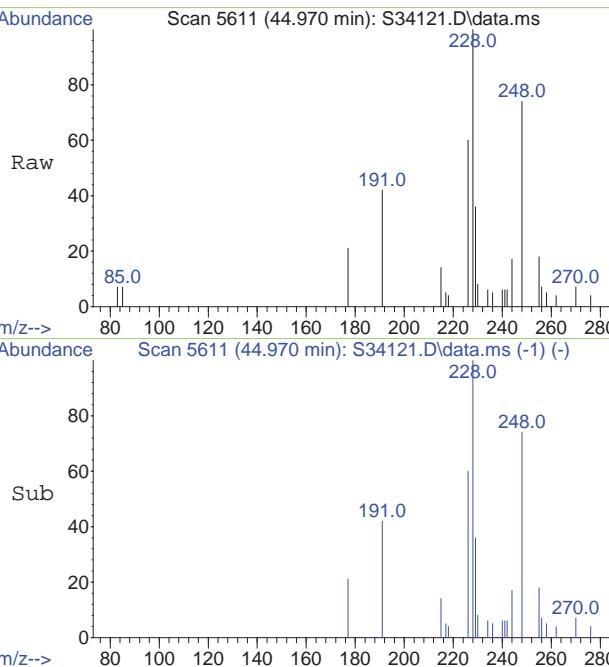


#72  
Retene  
Concen: 11.230 µg/mL  
RT: 41.093 min Scan# 4971  
Delta R.T. -0.006 min  
Lab File: S34121.D  
Acq: 20 Nov 2013 7:05 pm  
  
Tgt Ion: 234 Resp: 396  
Ion Ratio Lower Upper  
234 100  
219 162.1 131.3 196.9

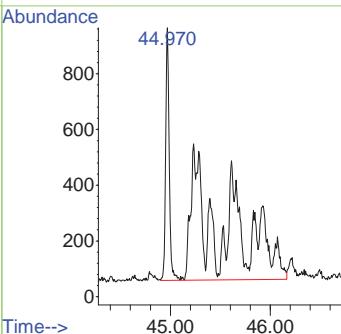
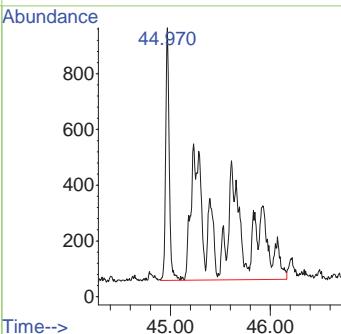


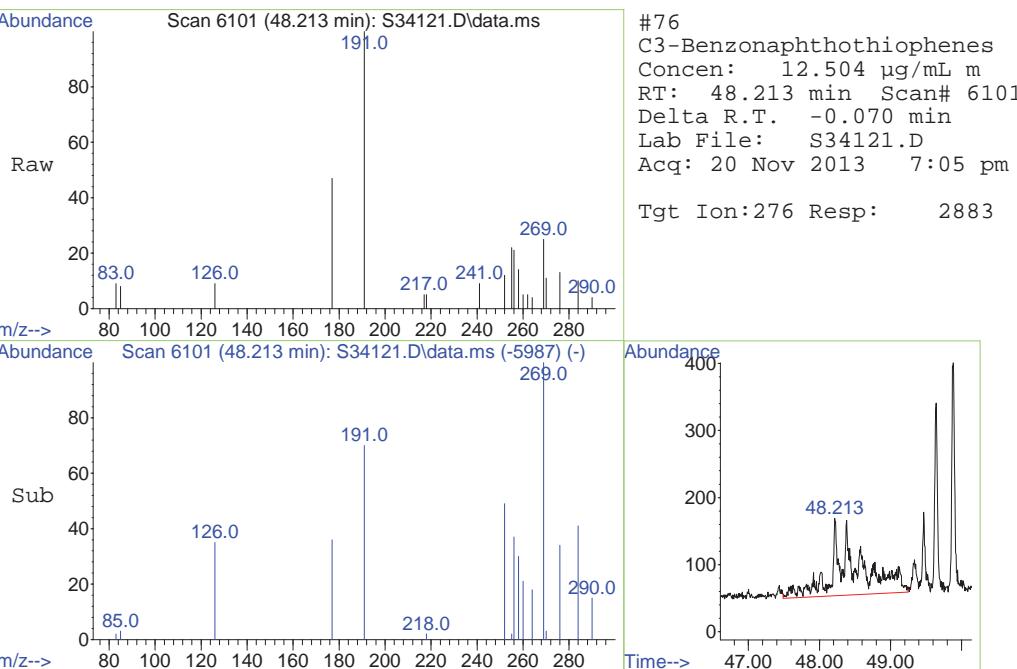
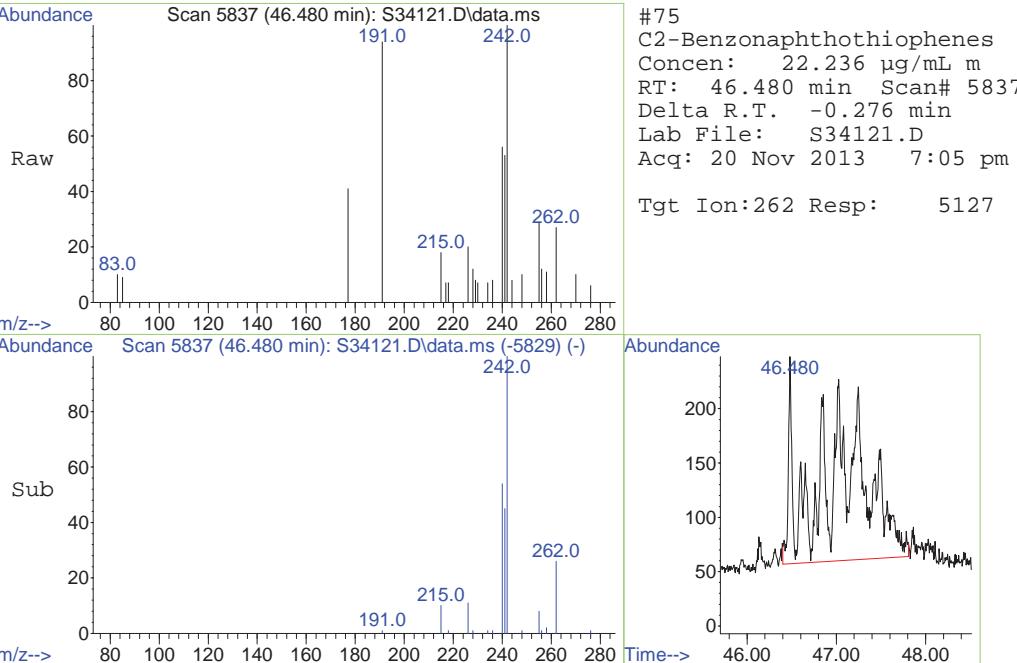


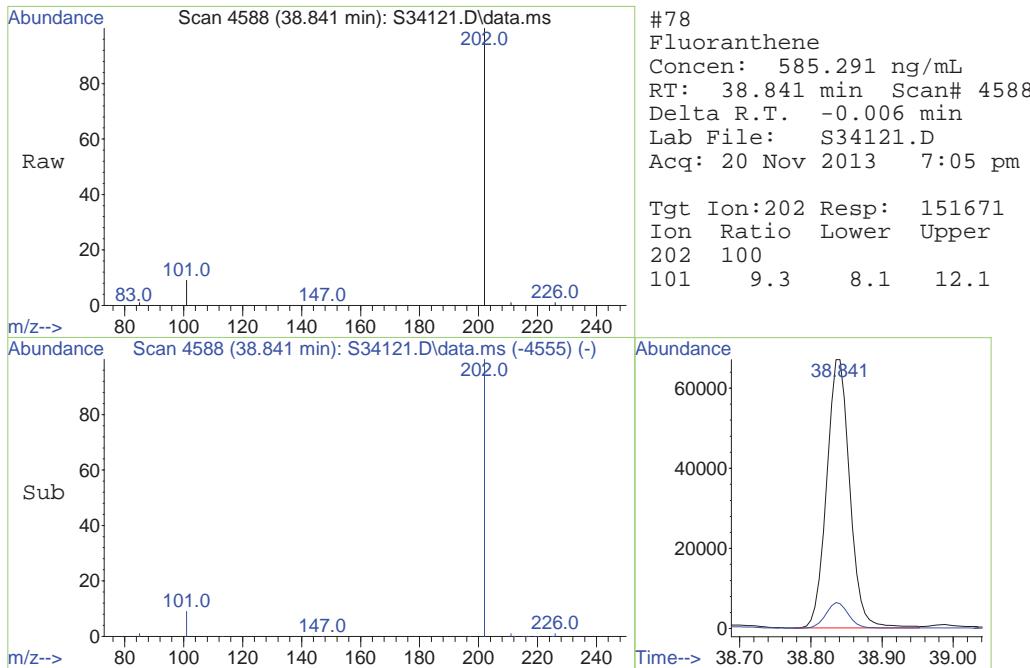
#73  
Benzo(b)naphtho(2,1-d)thiophene  
Concen: 51.519 µg/mL  
RT: 43.563 min Scan# 5391  
Delta R.T. -0.012 min  
Lab File: S34121.D  
Acq: 20 Nov 2013 7:05 pm  
Tgt Ion:234 Resp: 11879

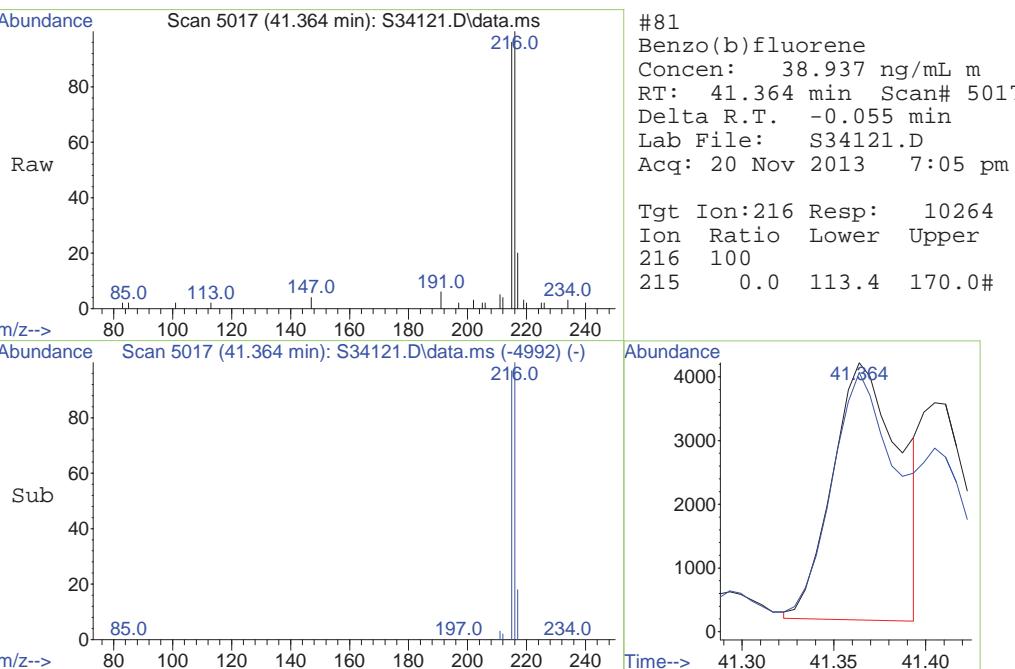
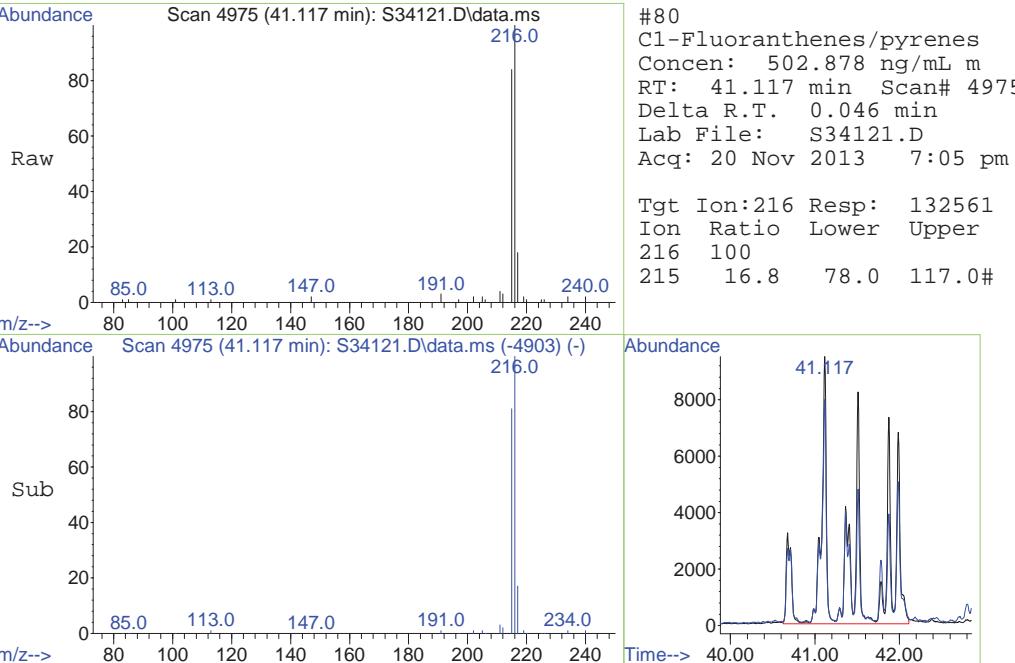


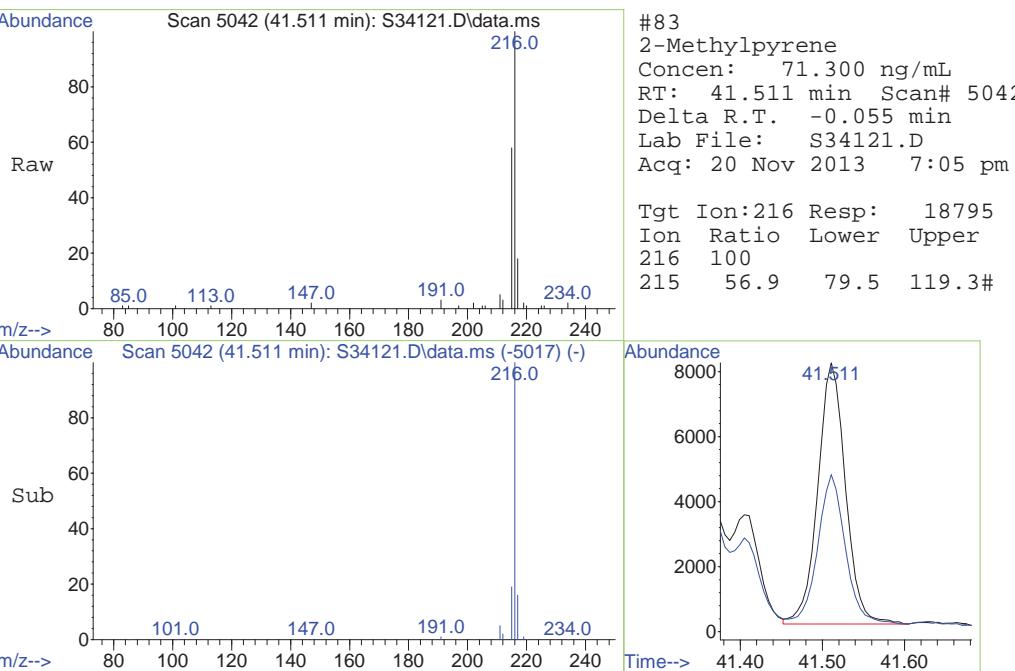
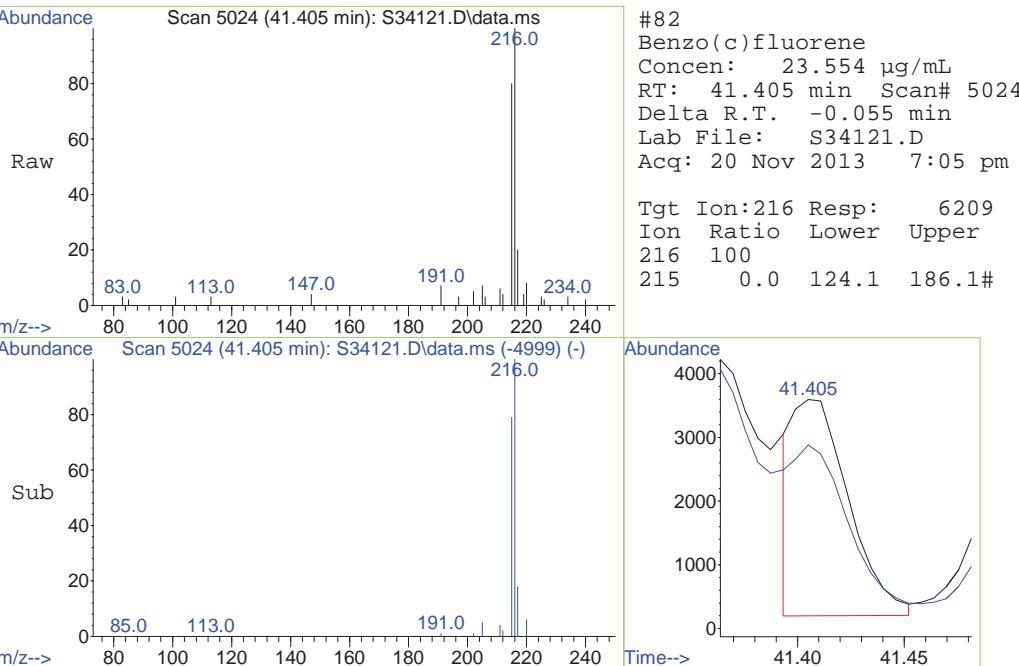
#74  
Cl-Benzonaphthothiophenes  
Concen: 55.392 µg/mL m  
RT: 44.970 min Scan# 5611  
Delta R.T. 0.232 min  
Lab File: S34121.D  
Acq: 20 Nov 2013 7:05 pm  
Tgt Ion:248 Resp: 12772

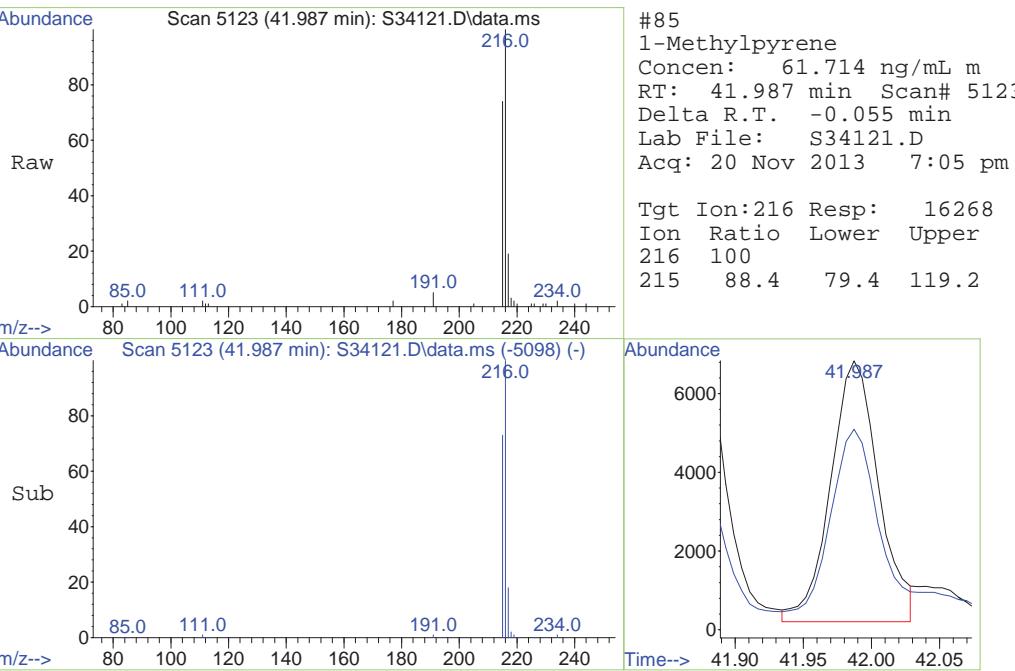
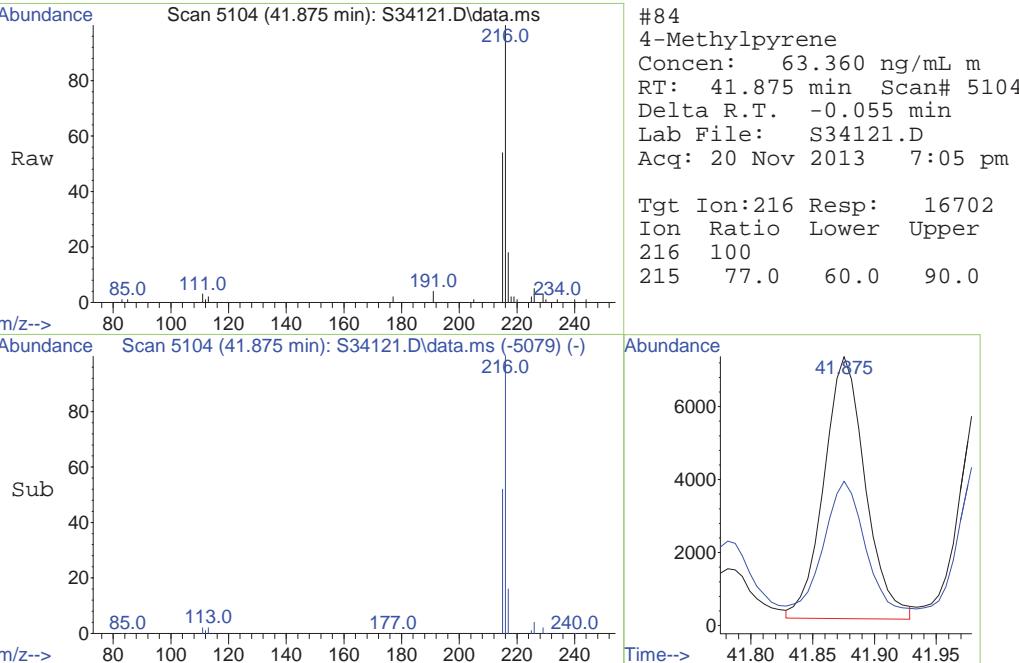


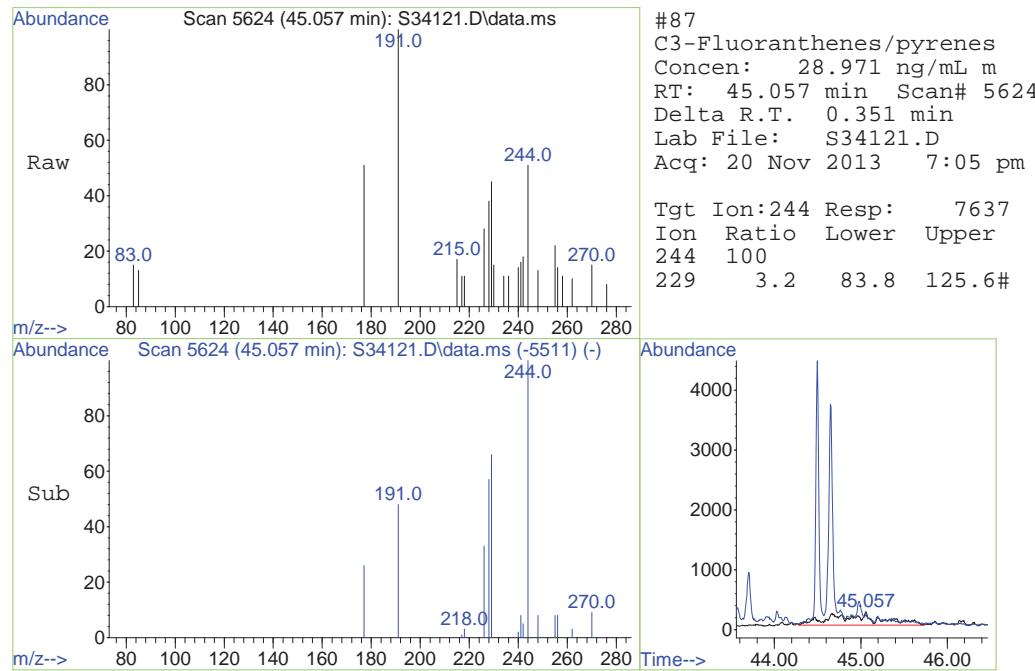
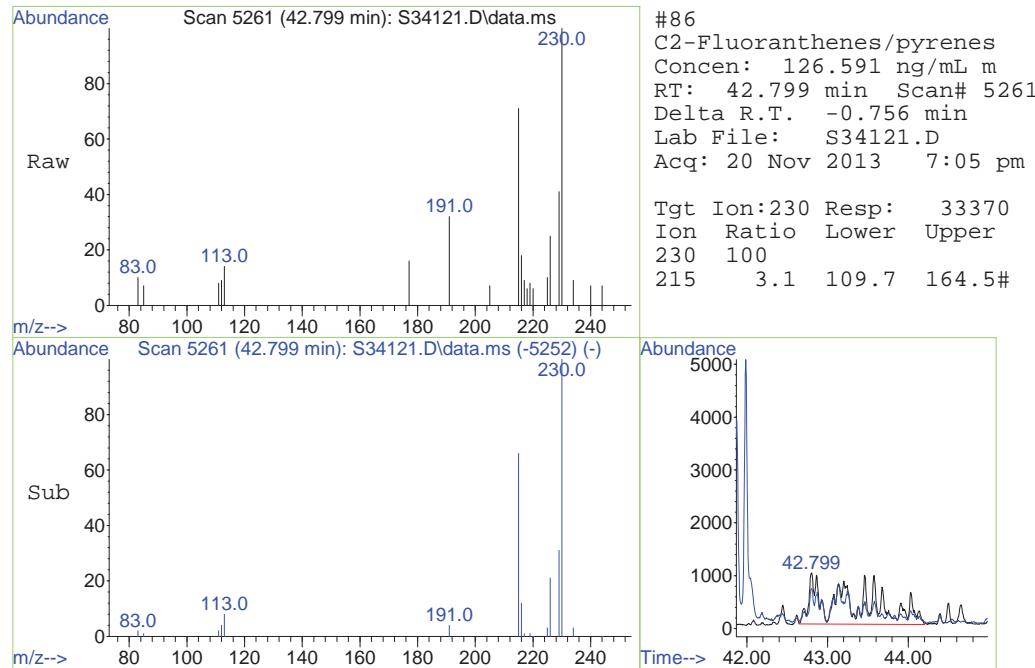


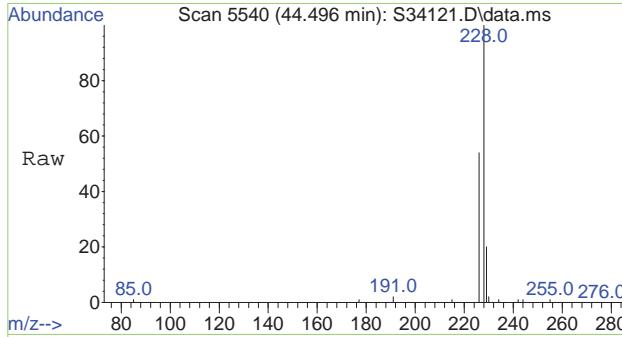






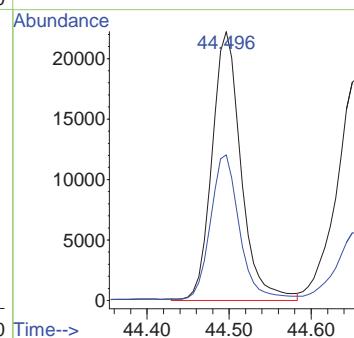
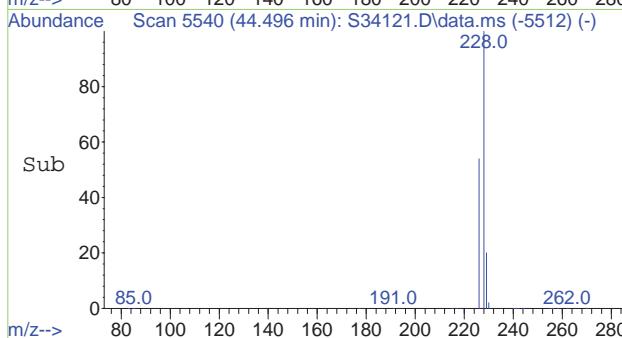






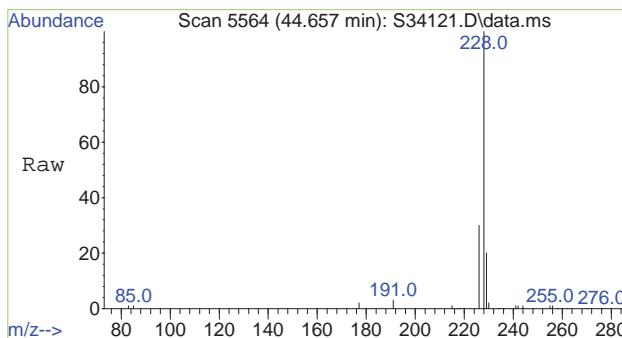
#88  
Benz(a)anthracene  
Concen: 244.165 ng/mL m  
RT: 44.496 min Scan# 5540  
Delta R.T. -0.014 min  
Lab File: S34121.D  
Acq: 20 Nov 2013 7:05 pm

Tgt Ion:	228	Resp:	54738
Ion Ratio	100	Lower	Upper
228	100	21.1	31.7
226	23.1		



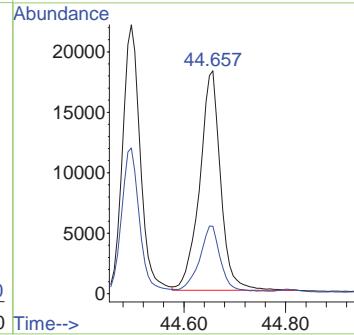
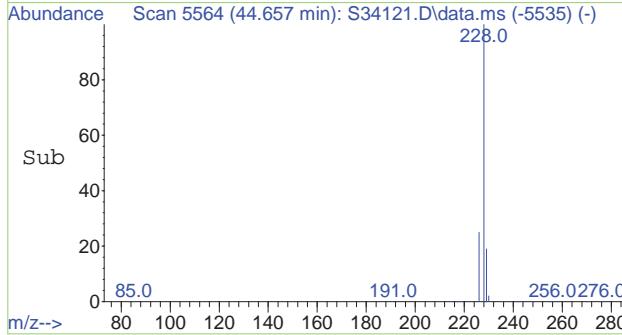
7.1.

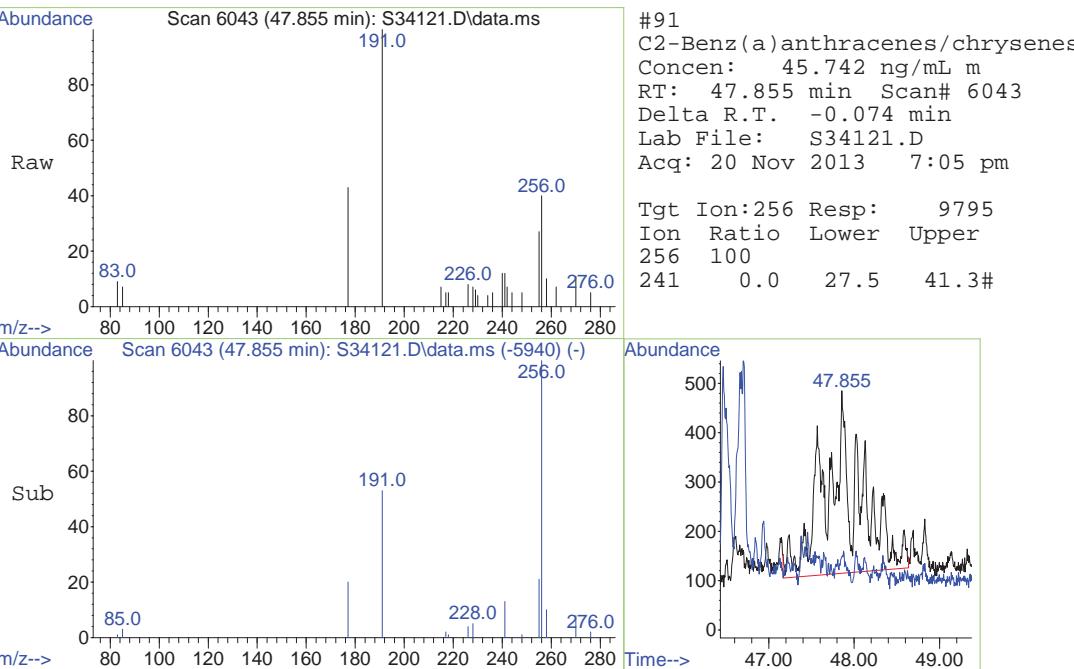
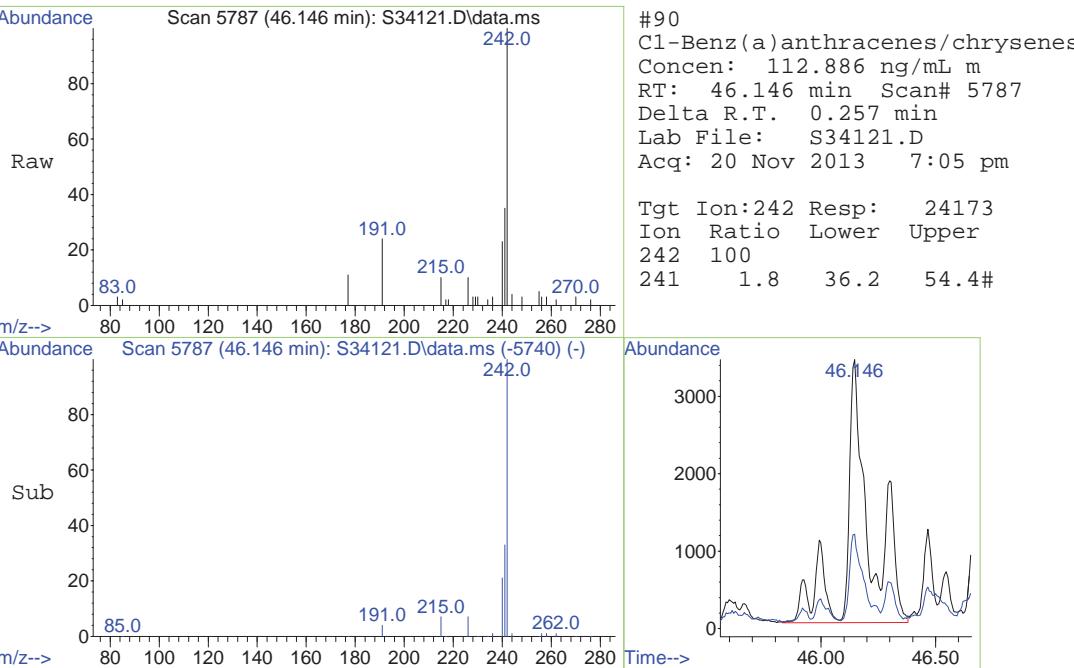
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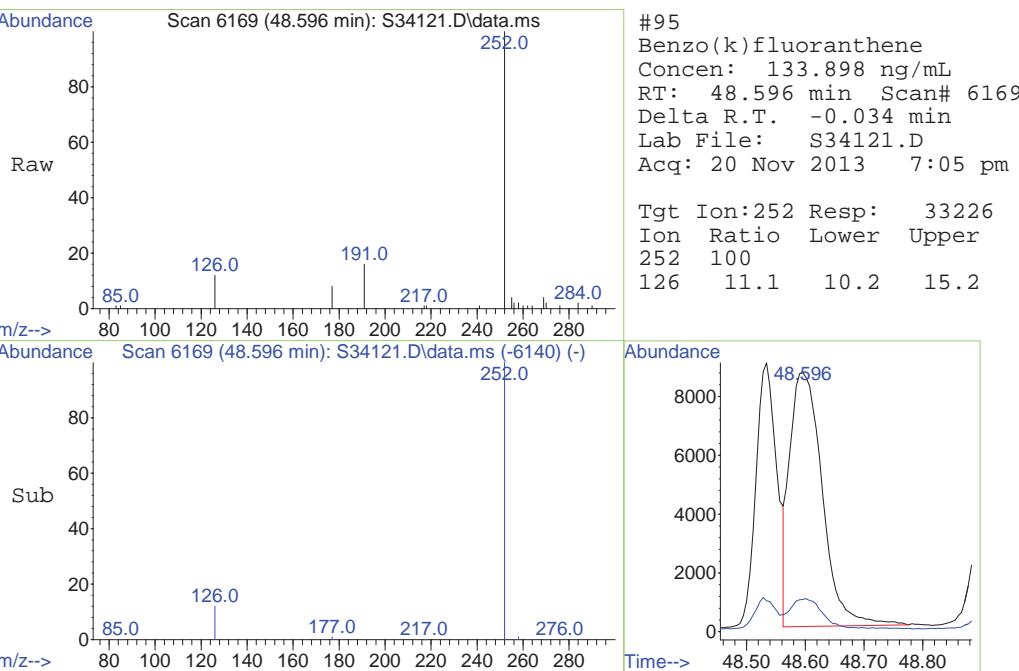
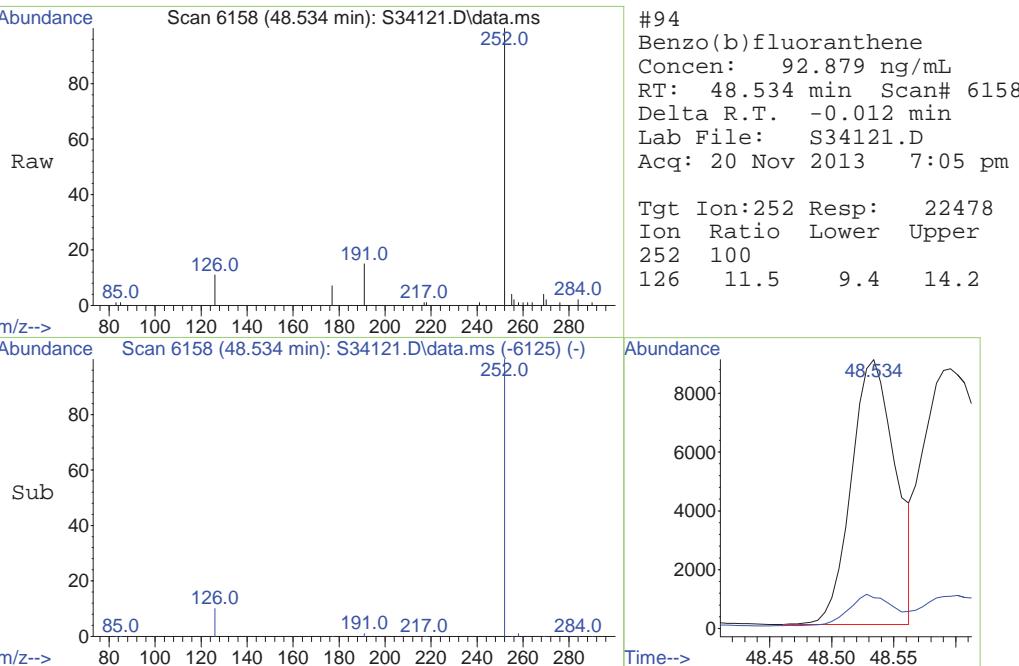


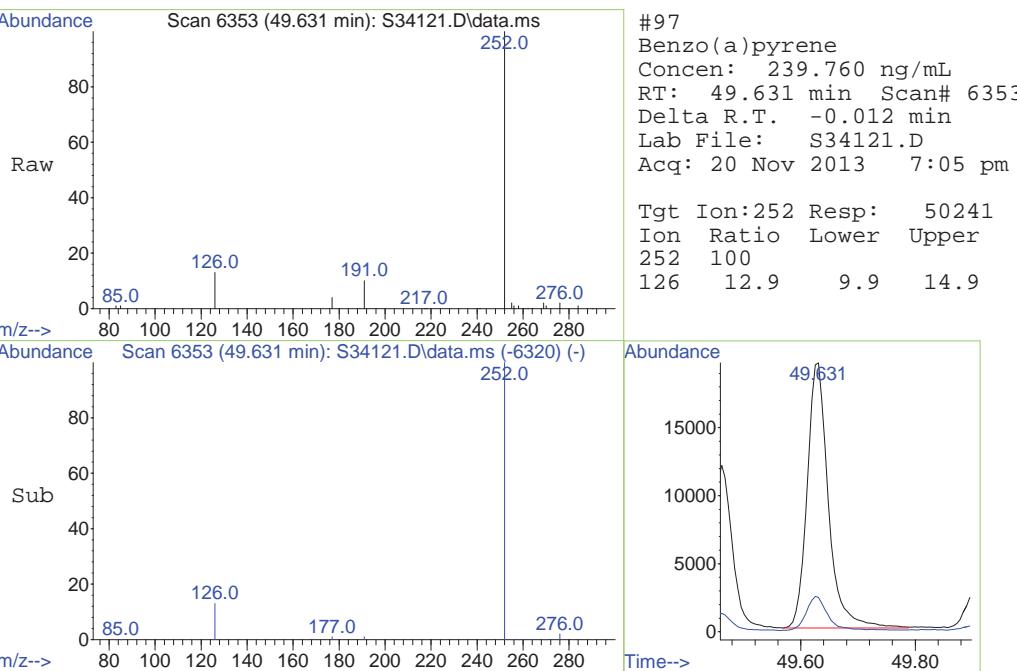
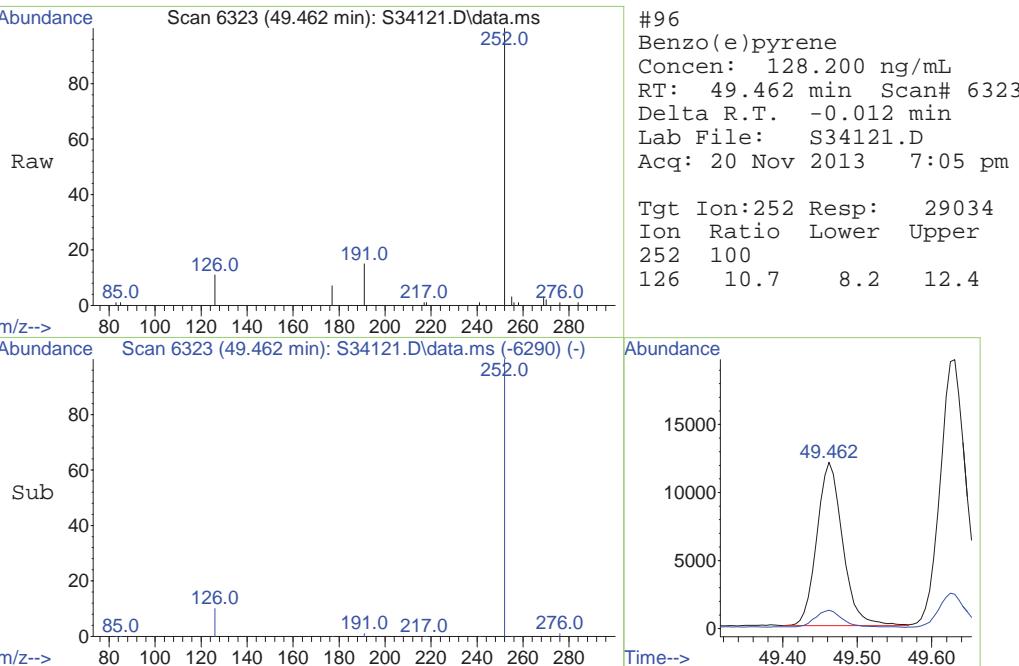
#89  
Chrysene  
Concen: 242.378 ng/mL  
RT: 44.657 min Scan# 5564  
Delta R.T. -0.007 min  
Lab File: S34121.D  
Acq: 20 Nov 2013 7:05 pm

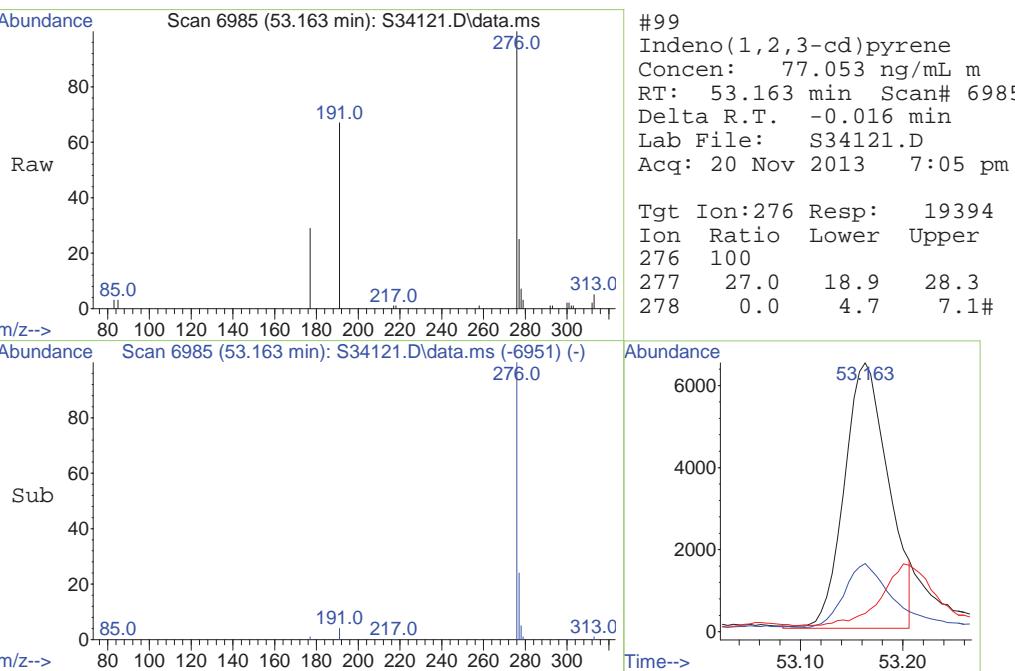
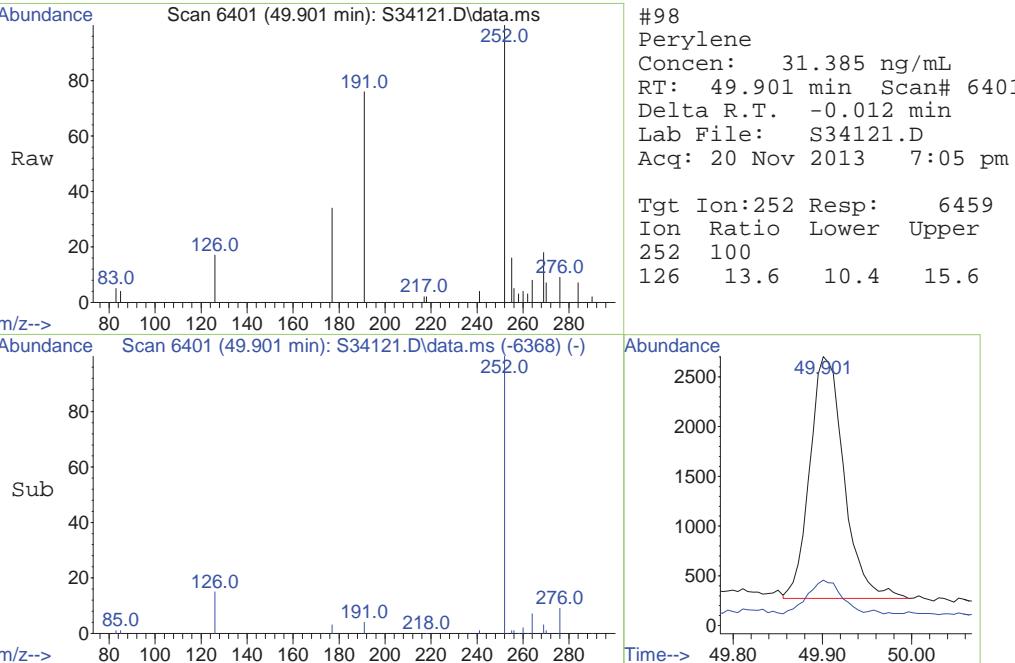
Tgt Ion:	228	Resp:	51902
Ion Ratio	100	Lower	Upper
228	100	23.7	35.5
226	29.5		

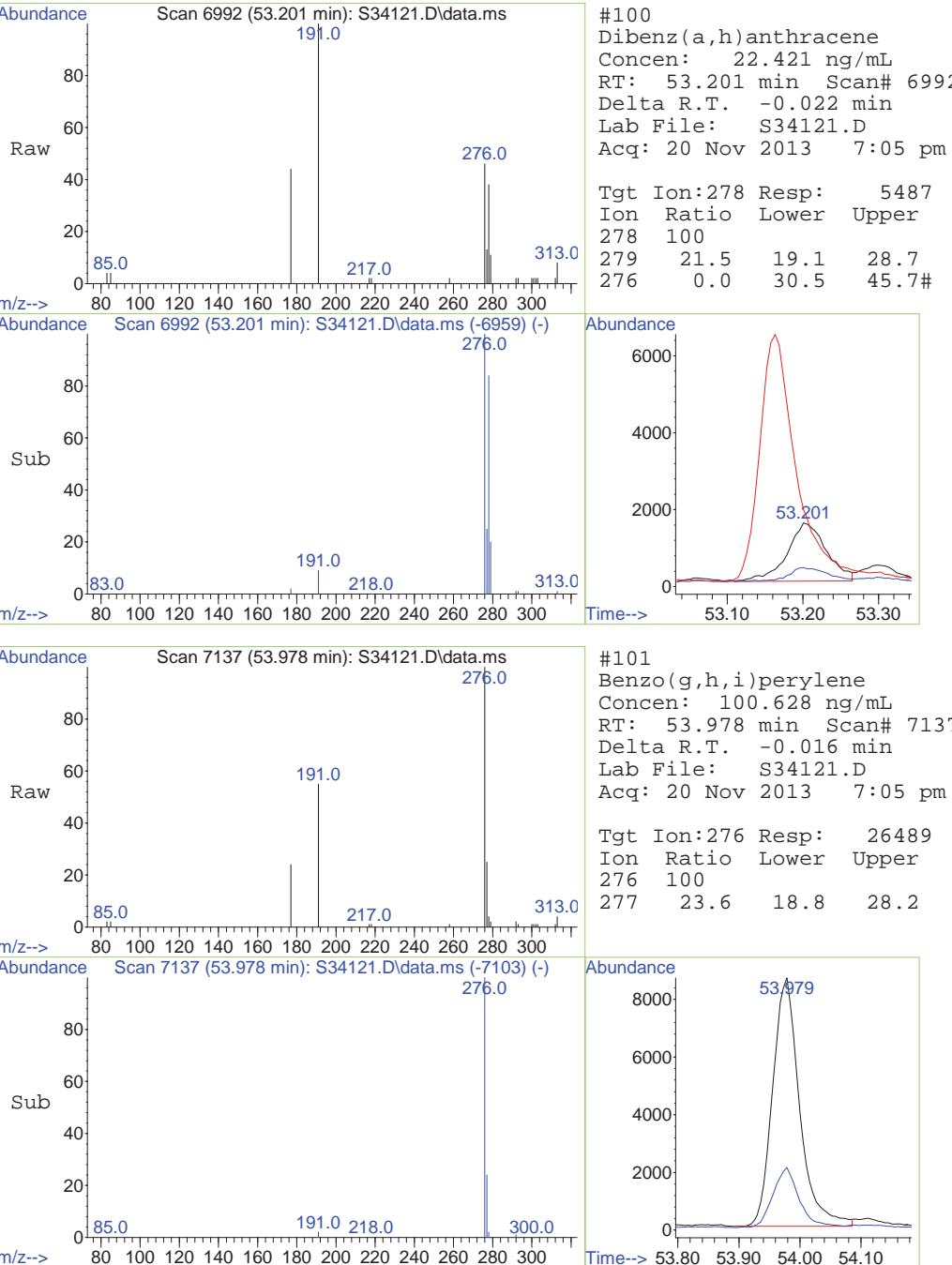


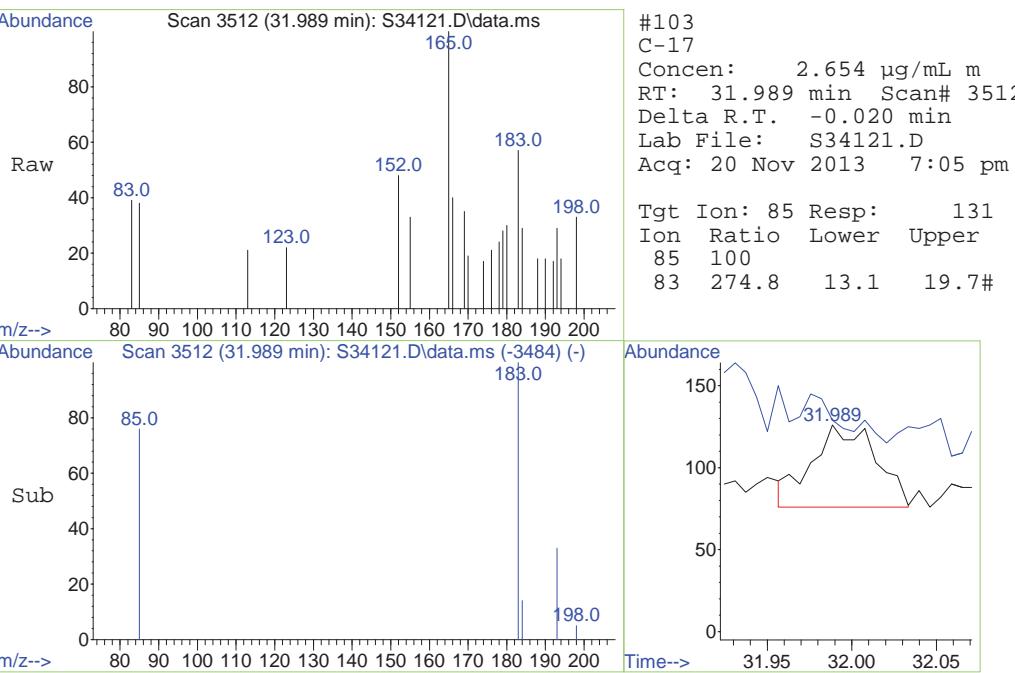
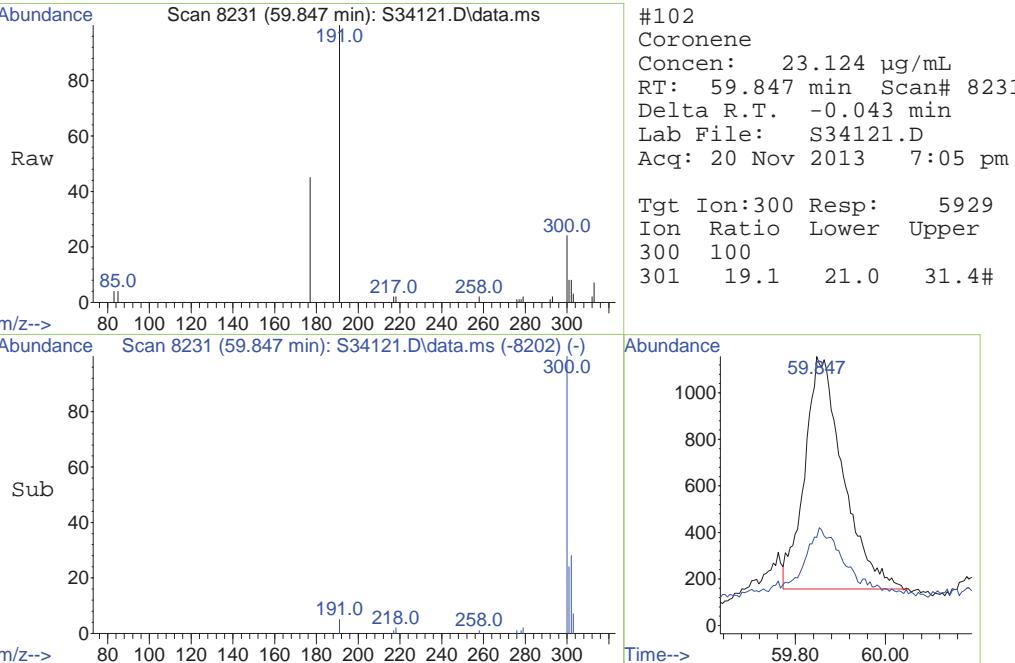


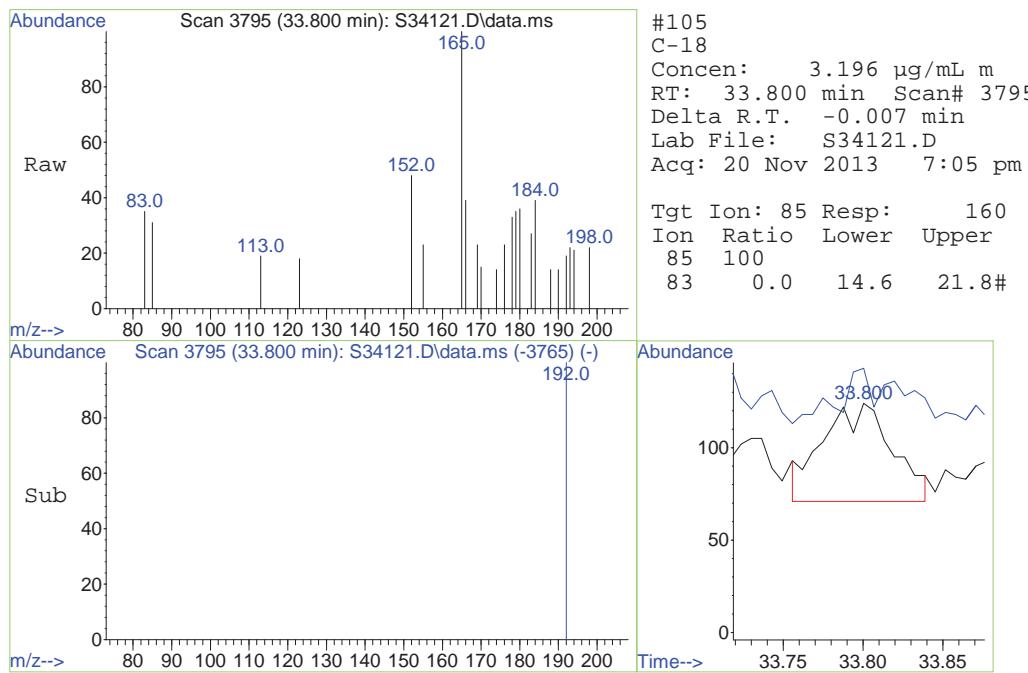
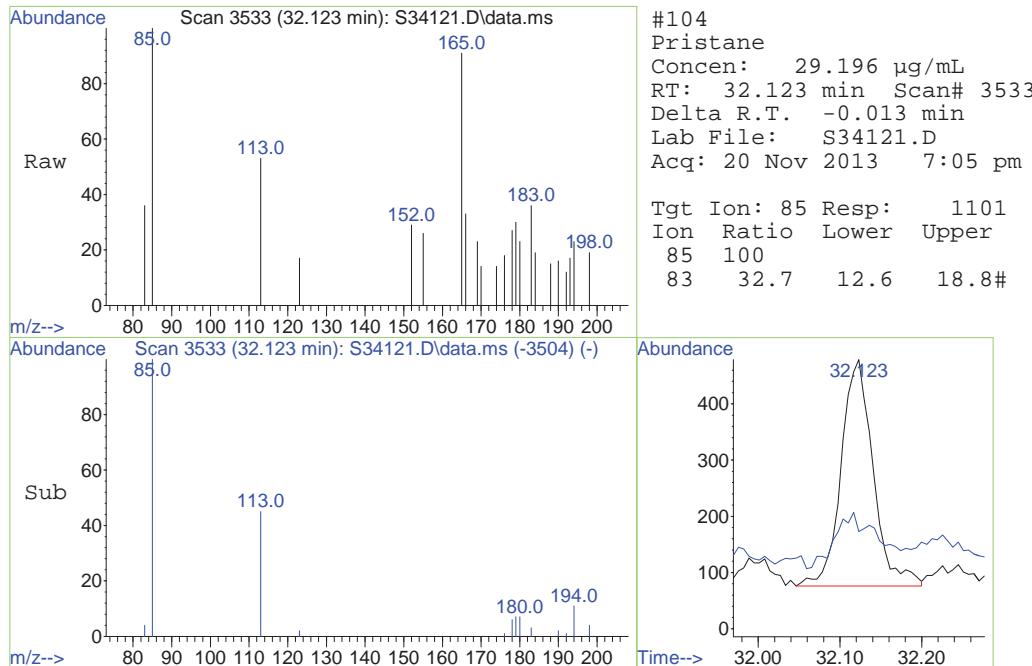


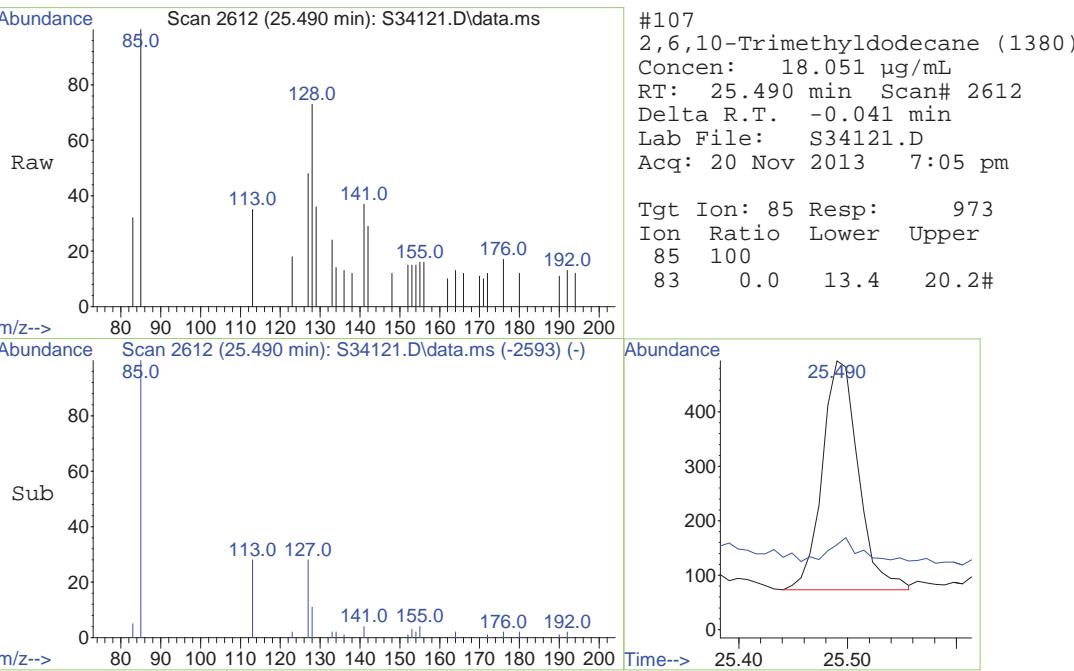
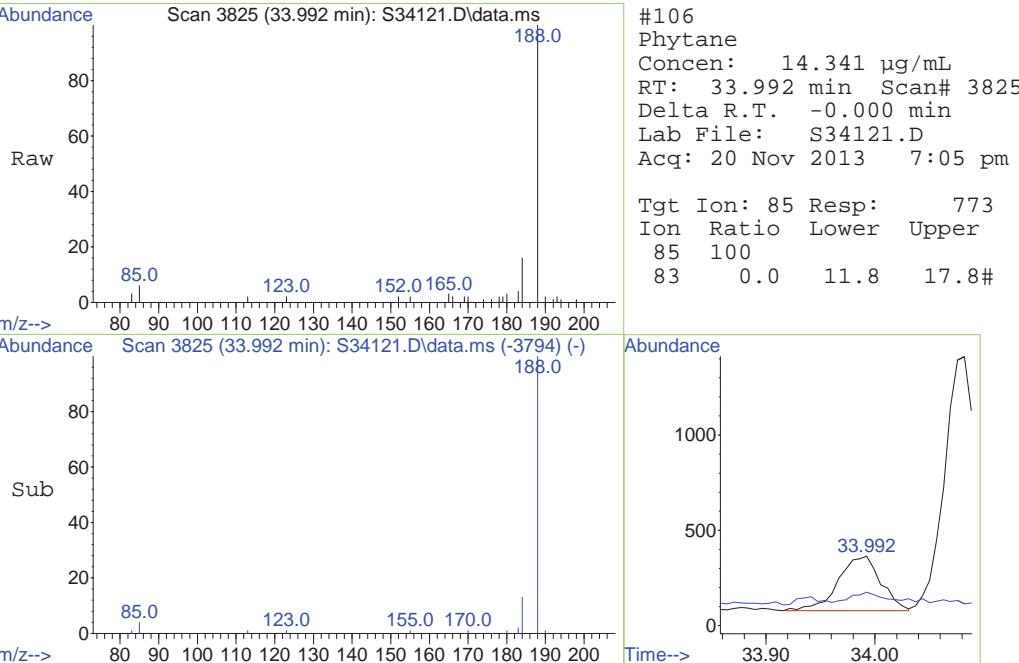


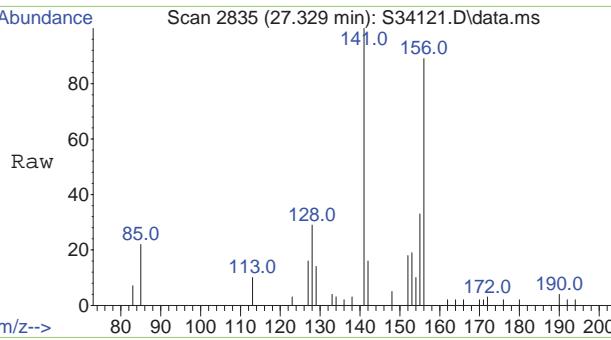






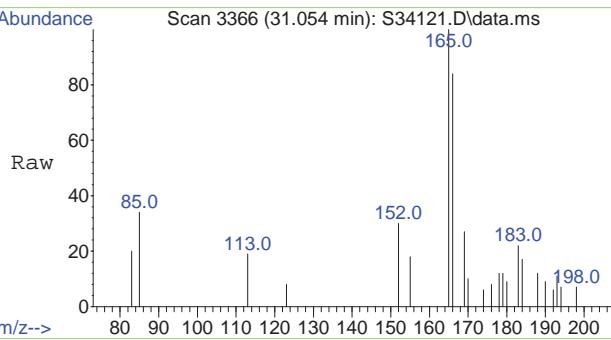
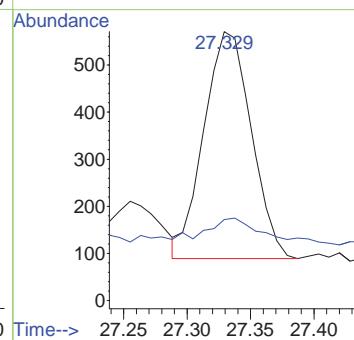
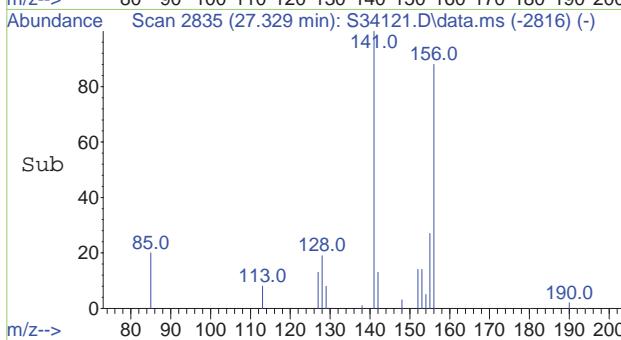






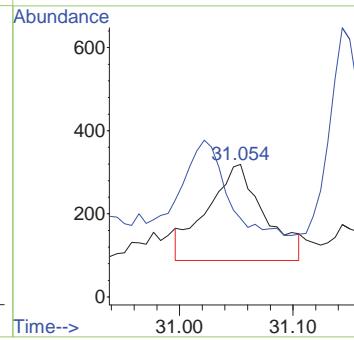
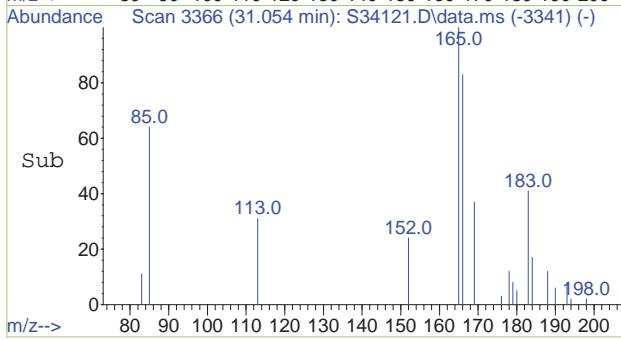
#108  
2,6,10-Trimethyltridecane (1470)  
Concen: 23.245 µg/mL  
RT: 27.329 min Scan# 2835  
Delta R.T. -0.041 min  
Lab File: S34121.D  
Acq: 20 Nov 2013 7:05 pm

Tgt Ion: 85 Resp: 1253  
Ion Ratio Lower Upper  
85 100  
83 17.6 13.8 20.8



#109  
2,6,10-Trimethylpentadecane (1650)  
Concen: 14.934 µg/mL m  
RT: 31.054 min Scan# 3366  
Delta R.T. -0.042 min  
Lab File: S34121.D  
Acq: 20 Nov 2013 7:05 pm

Tgt Ion: 85 Resp: 805  
Ion Ratio Lower Upper  
85 100  
83 0.0 20.7 31.1#



Data Path : Z:\1\data\S131120\  
 Data File : S34116.D  
 Acq On : 20 Nov 2013 1:04 pm  
 Operator : RUBENP  
 Sample : op35834-mb  
 Misc : op35834,mss1523,0.01,,,2,1  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Nov 20 16:15:04 2013  
 Quant Method : Z:\1\methods\S131114-MAHPAHEXT.M  
 Quant Title : PAHs & Alkylated PAHs by GC/MS/SIM  
 QLast Update : Fri Nov 15 10:17:28 2013  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<hr/>						
Internal Standards						
1) Acenaphthene-d10	28.166	164	85103	1000.00	ng/mL	0.00
<hr/>						
System Monitoring Compounds						
2) Toluene-d8	9.271	98	93898	1082.75	µg/mL	-0.03
Spiked Amount	1000.000		Recovery	=	108.28%	
3) Naphthalene-d8	21.218	136	146371	904.10	ng/mL	0.00
Spiked Amount	1000.000	Range 1 - 0	Recovery	=	90.41%#	
4) Phenanthrene-d10	33.999	188	136050	878.03	ng/mL	0.00
Spiked Amount	1000.000		Recovery	=	87.80%	
5) Perylene-d12	49.828	264	118424	875.27	ng/mL	0.00
Spiked Amount	1000.000		Recovery	=	87.53%	
<hr/>						
Target Compounds						
7) Benzene	6.844	78	4270	46.095	µg/mL#	1
61) C1-Dibenzothiophenes (...	35.819	198	14508	76.060	µg/mL	97
69) C2-Phenanthrenes/anthr...	38.058	206	7926	40.686	µg/mL	89

(#) = qualifier out of range (m) = manual integration (+) = signals summed

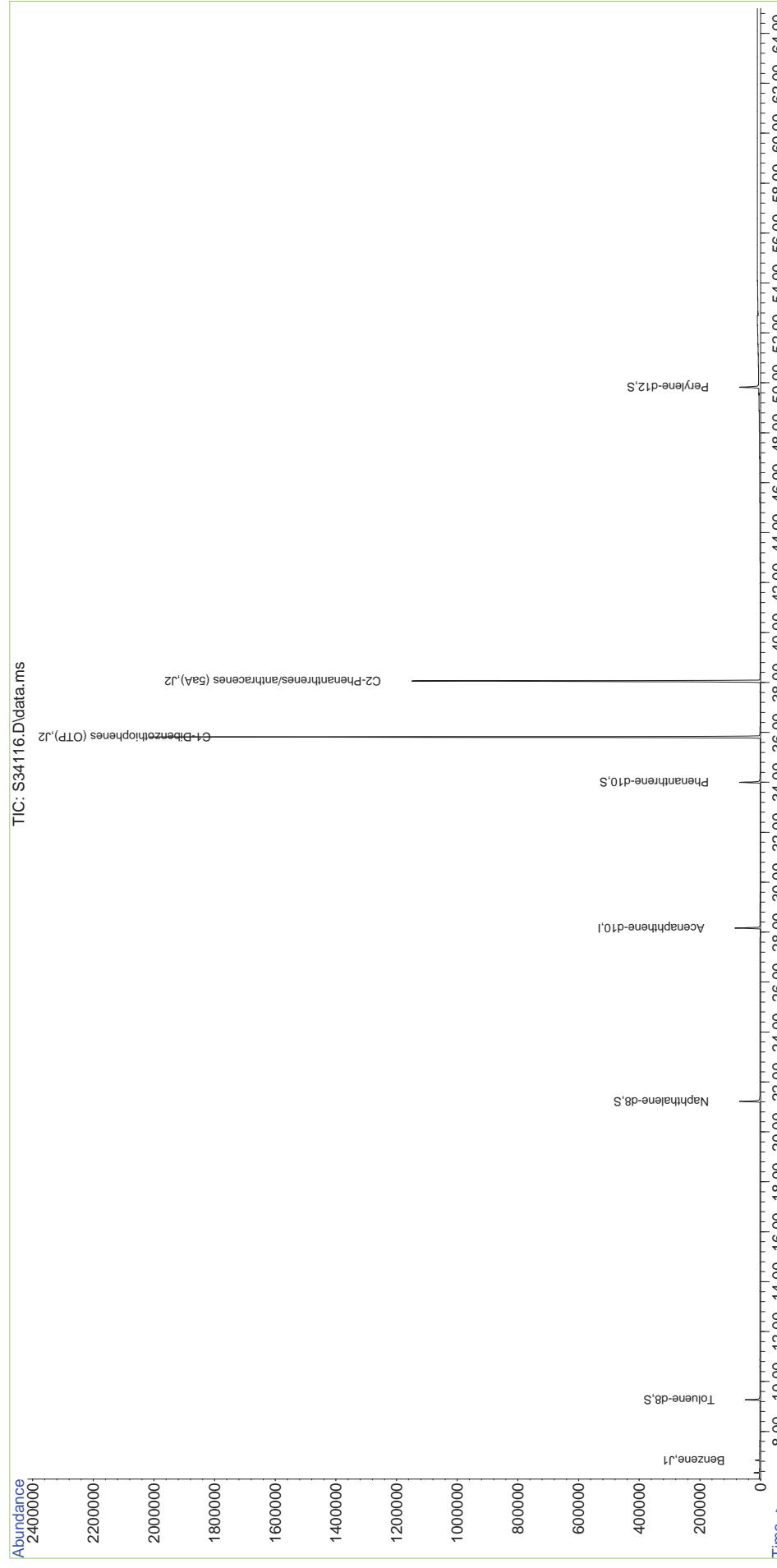
7.2.1

7

## Quantitation Report (QT Reviewed)

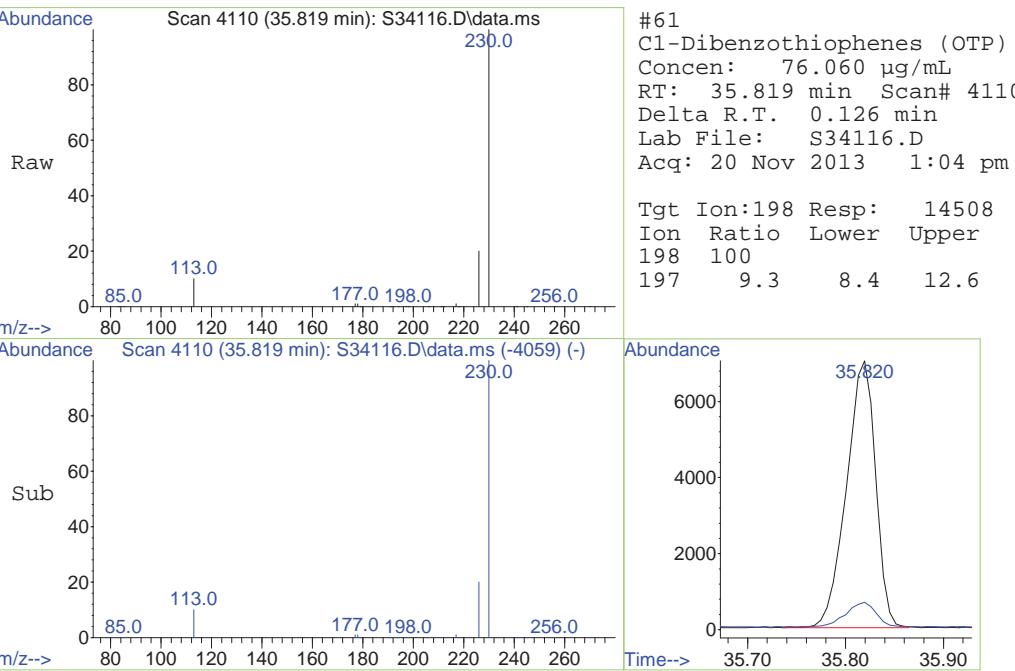
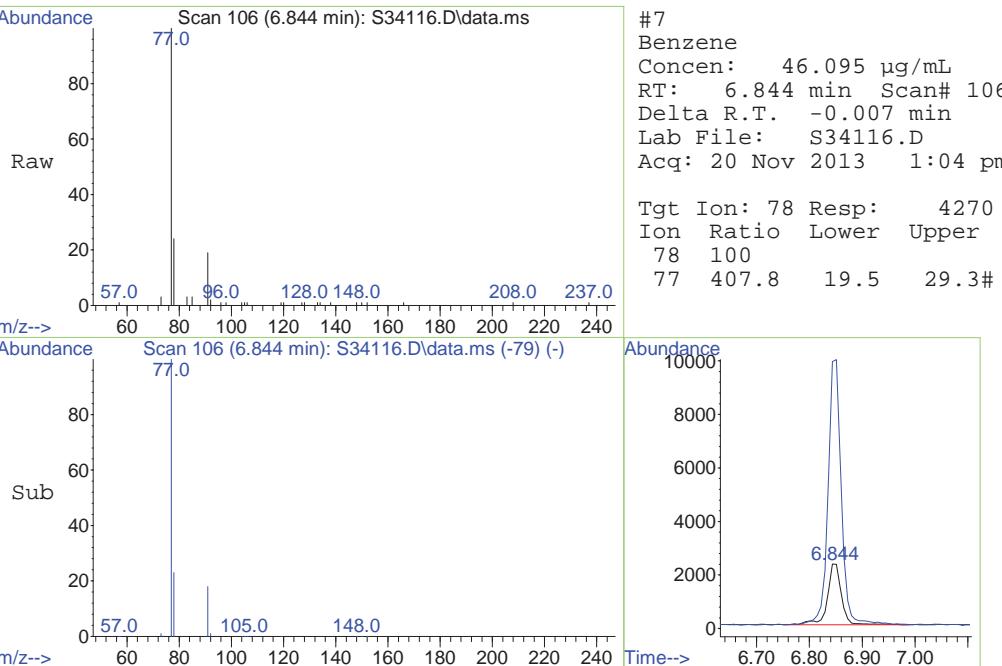
Data Path : Z:\1\data\S131120\  
 Data File : S34116.D  
 Acc On : 20 Nov 2013 1:04 pm  
 Operator : RUBENP  
 Sample : op35834-mb  
 Misc : op35834,mss1523.0.01,,2,1  
 ALS Vial : 3 Sample Multiplier: 1

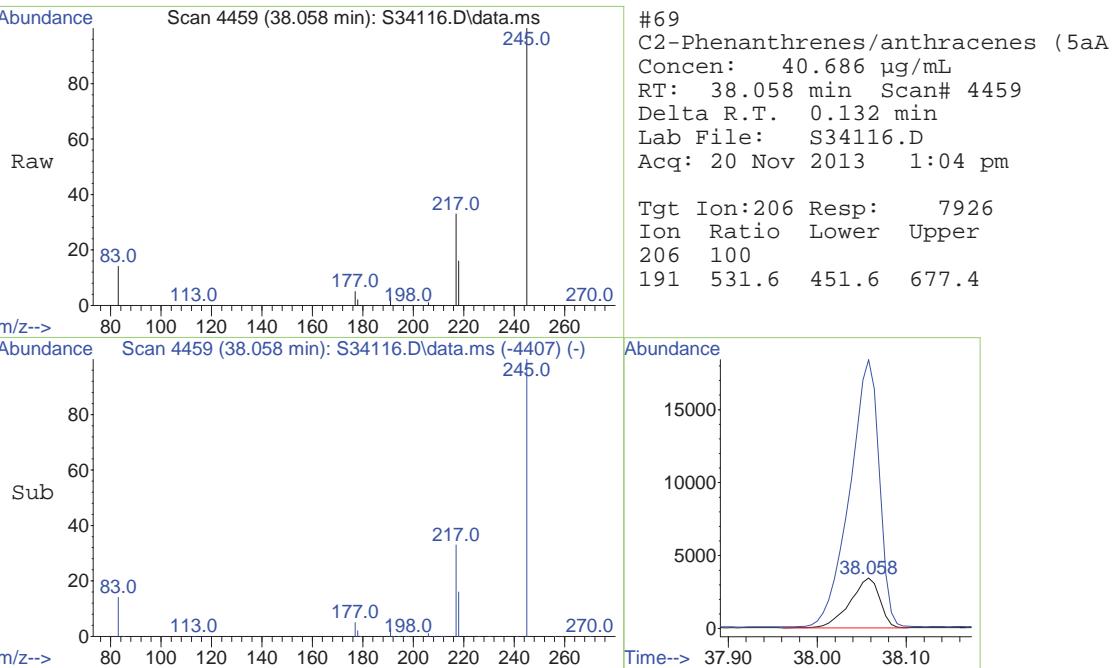
Quant Time: Nov 20 16:15:04 2013  
 Quant Method : Z:\1\methods\S131114-MAHPAHEXT.M  
 Quant Title : PAHs & Alkylated PAHs by GC/MS/SIM  
 QLast Update : Fri Nov 15 10:17:28 2013  
 Response via : Initial Calibration



.131114-MAHPAHEXT.M Thu Nov 21 12:24:42 2013

Page: 2





7.2.1

7



## GC Semi-volatiles

### QC Data Summaries



Includes the following where applicable:

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries
- Surrogate Recovery Summaries

**Method Blank Summary**

Page 1 of 1

**Job Number:** MC26103  
**Account:** METAMAW META Environmental, Inc.  
**Project:** Parsons, Halleck Street, NJ

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP35833-MB	BG42071.D	1	11/19/13	RP	11/19/13	OP35833	GBG1601

**The QC reported here applies to the following samples:****Method:** ASTM D3328-06

MC26103-1

CAS No.	Compound	Result	RL	Units	Q
	TPH (C8-C40)	ND	20000	mg/kg	

CAS No.	Surrogate Recoveries	Limits
84-15-1	o-Terphenyl	101% 40-140%

**Blank Spike Summary**

Page 1 of 1

**Job Number:** MC26103**Account:** METAMAW META Environmental, Inc.**Project:** Parsons, Halleck Street, NJ

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP35833-BS1	BG42072.D	1	11/19/13	RP	11/19/13	OP35833	GBG1601

**The QC reported here applies to the following samples:****Method:** ASTM D3328-06

MC26103-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH (C8-C40)		ND		40-140

CAS No.	Surrogate Recoveries	BSP	Limits
84-15-1	o-Terphenyl	101%	40-140%

8.2.1

8

\* = Outside of Control Limits.

**Duplicate Summary**

Page 1 of 1

**Job Number:** MC26103  
**Account:** METAMAW META Environmental, Inc.  
**Project:** Parsons, Halleck Street, NJ

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP35833-DUP1	BG42081.D	1	11/20/13	RP	11/19/13	OP35833	GBG1601
MC26103-1 <sup>a</sup>	BG42080.D	1	11/20/13	RP	11/19/13	OP35833	GBG1601

The QC reported here applies to the following samples:

Method: ASTM D3328-06

MC26103-1

CAS No.	Compound	MC26103-1 DUP					
		mg/kg	Q	mg/kg	Q	RPD	Limits
	TPH (C8-C40)	2510000		2320000		8	30

CAS No.	Surrogate Recoveries	DUP	MC26103-1 Limits		
84-15-1	o-Terphenyl		445% * <sup>b</sup>	430% * <sup>b</sup>	40-140%

(a) Sample extracted beyond the recommended holding time for soils.

(b) Outside control limits due to dilution.

8.3.1

8

\* = Outside of Control Limits.

**Semivolatile Surrogate Recovery Summary**

Page 1 of 1

**Job Number:** MC26103**Account:** METAMAW META Environmental, Inc.**Project:** Parsons, Halleck Street, NJ**Method:** ASTM D3328-06**Matrix:** SO**Samples and QC shown here apply to the above method**

<b>Lab Sample ID</b>	<b>Lab File ID</b>	<b>S1 <sup>a</sup></b>
MC26103-1	BG42080.D	430.0* <sup>b</sup>
OP35833-BS1	BG42072.D	101.0
OP35833-DUP1	BG42081.D	445.0* <sup>b</sup>
OP35833-MB	BG42071.D	101.0

<b>Surrogate Compounds</b>	<b>Recovery Limits</b>
<b>S1 = o-Terphenyl</b>	40-140%

(a) Recovery from GC signal #1

(b) Outside control limits due to dilution.

8.4.1

8



New England  
**ACCUTEST<sup>®</sup>**  
LABORATORIES

## GC Semi-volatiles

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Raw Data

---

Data Path : G:\1\DATA\BG131119\  
 Data File : BG42080.d  
 Signal(s) : FID1A.ch  
 Acq On : 20 Nov 2013 4:01 am  
 Operator : RubenP  
 Sample : MC26103-1  
 Misc : OP35833,GBG1601,0.0022,,,10,1  
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 22 10:19:51 2013  
 Quant Method : G:\1\METHODS\BG130925ALK-Front.m  
 Quant Title : n-C8 - n-C40 normal alkanes w/ isoprenoids  
 Response via : Initial Calibration

Compound	R.T.	Response	Conc	Units
<hr/>				
Internal Standards				
1) I 5a-Androstan e	31.102	136871228	50.000	µg/mL
<hr/>				
System Monitoring Compounds				
2) S o-Terphenyl	29.101	61777629	21.491	µg/mL
Spiked Amount	25.000	Recovery	=	85.96%
<hr/>				
Target Compounds				
41) H TPH (C8-C40)	30.880	1450898006	551.605	µg/mL
<hr/>				
SemiQuant Compounds - Not Calibrated on this Instrument				
<hr/>				

(f)=RT Delta &gt; 1/2 Window

(m)=manual int.

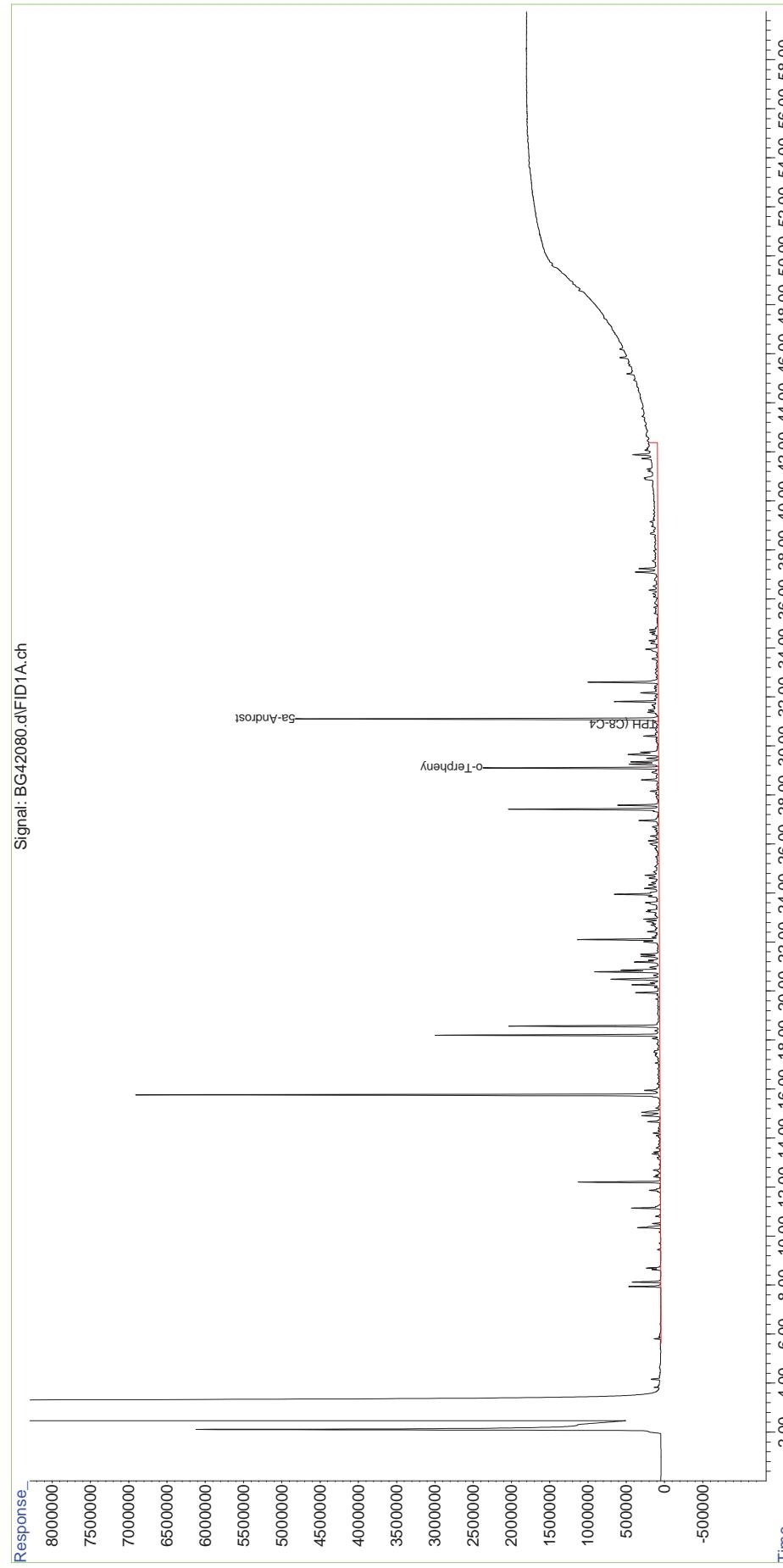
9.1.1

6

## Quantitation Report (QT Reviewed)

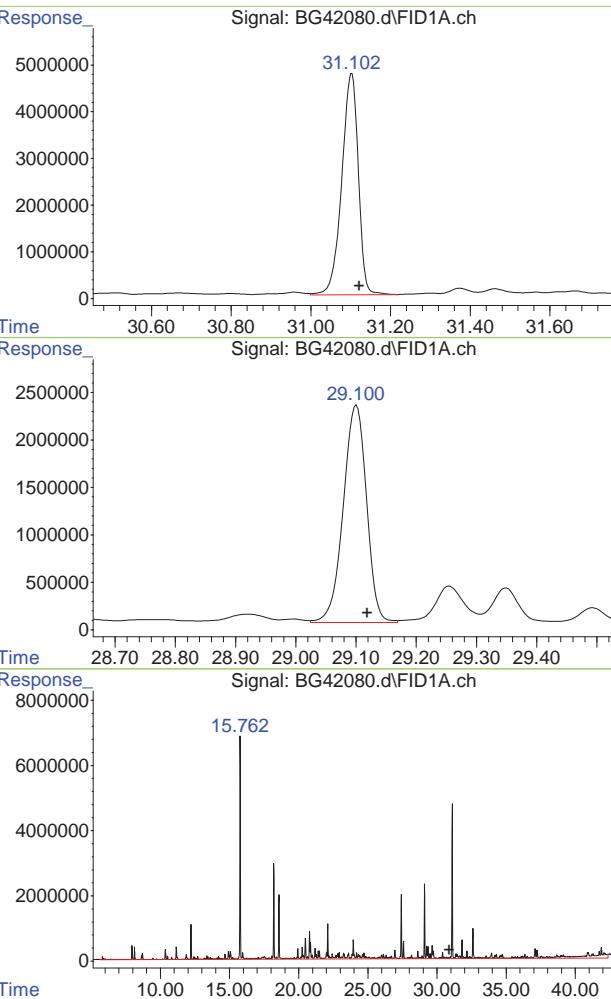
Data Path : G:\1\DATA\BG131119\  
 Data File : BG42080.d  
 Signal(s) : FID1A.ch  
 Acq On : 20 Nov 2013 4:01 am  
 Operator : RubenP  
 Sample : MC26103-1  
 Misc : OP35833,GBG1601,0.0022,,10,1  
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: Nov 22 10:19:51 2013  
 Quant Method : G:\1\METHODS\BG130925ALK-Front.m  
 Quant Title : n-C8 - n-C40 normal alkanes w/ isoprenoids  
 Response via : Initial Calibration



G130925ALK-Front.m Fri Nov 22 10:20:11 2013

Page: 2



#1 5a-Androstan

R.T.: 31.102 min  
Delta R.T.: -0.019 min  
Response: 136871228  
Conc: 50.00 µg/mL

#2 o-Terphenyl

R.T.: 29.101 min  
Delta R.T.: -0.018 min  
Response: 61777629  
Conc: 21.49 µg/mL

#41 TPH (C8-C40)

R.T.: 30.880 min  
Delta R.T.: 0.000 min  
Response: 1450898006  
Conc: 551.61 µg/mL

9.1.1

6

Data Path : G:\1\DATA\BG131119\  
 Data File : BG42071.d  
 Signal(s) : FID1A.ch  
 Acq On : 19 Nov 2013 6:03 pm  
 Operator : RubenP  
 Sample : op35833-MB  
 Misc : OP35833,GBG1601,0.01,,,2,1  
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Nov 20 13:50:21 2013  
 Quant Method : G:\1\METHODS\BG130925ALK-Front.m  
 Quant Title : n-C8 - n-C40 normal alkanes w/ isoprenoids  
 Response via : Initial Calibration

Compound	R.T.	Response	Conc Units
<hr/>			
Internal Standards			
1) I 5a-Androstan e	31.132	127120354	50.000 µg/mL
<hr/>			
System Monitoring Compounds			
2) S o-Terphenyl	29.127	67666792	25.345 µg/mL
Spiked Amount	25.000	Recovery	= 101.38%
<hr/>			
Target Compounds			
41) H TPH (C8-C40)	30.880	75556542	30.929 µg/mL
<hr/>			
SemiQuant Compounds - Not Calibrated on this Instrument			
<hr/>			

(f)=RT Delta &gt; 1/2 Window

(m)=manual int.

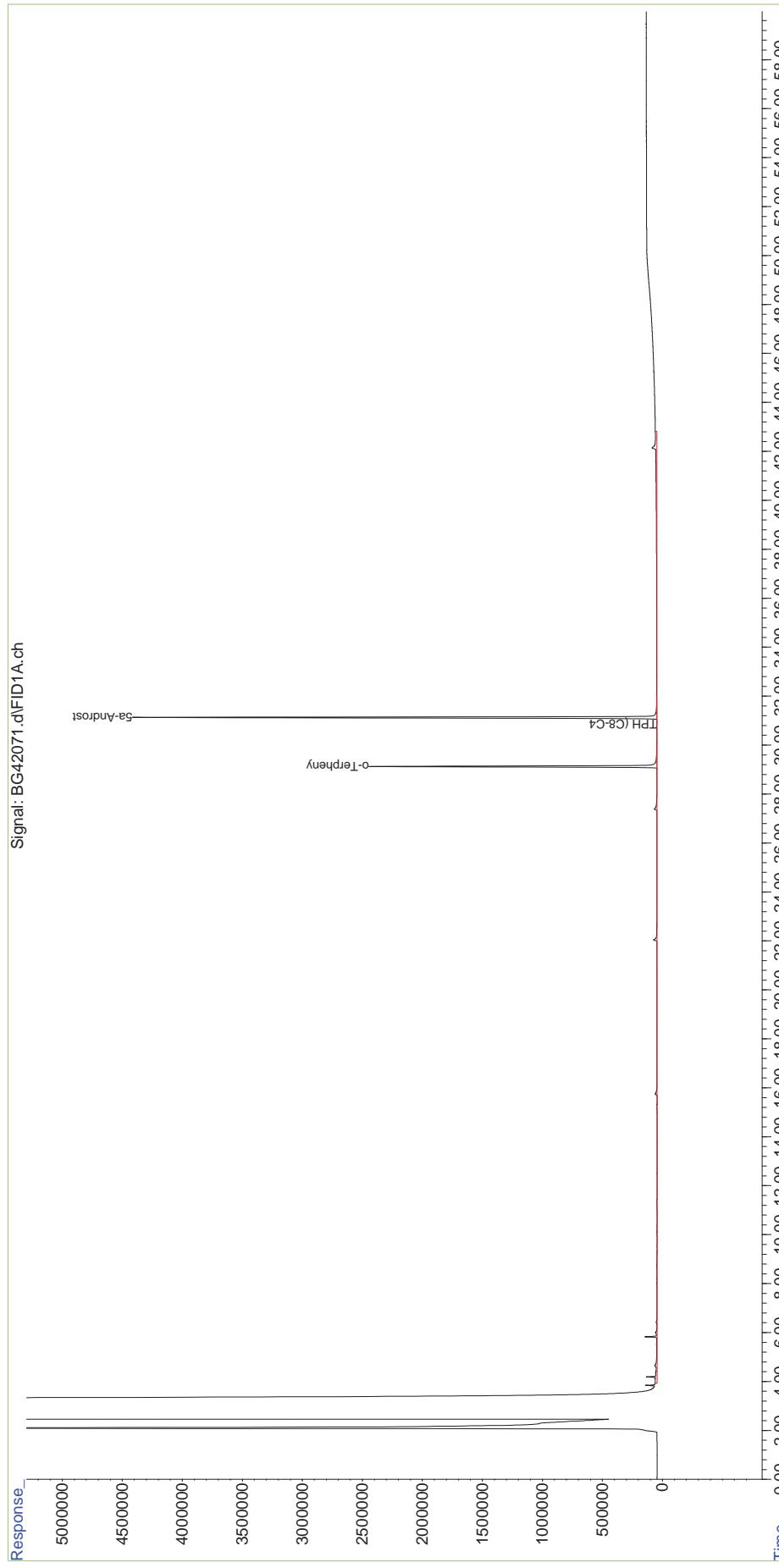
9.2.1

9

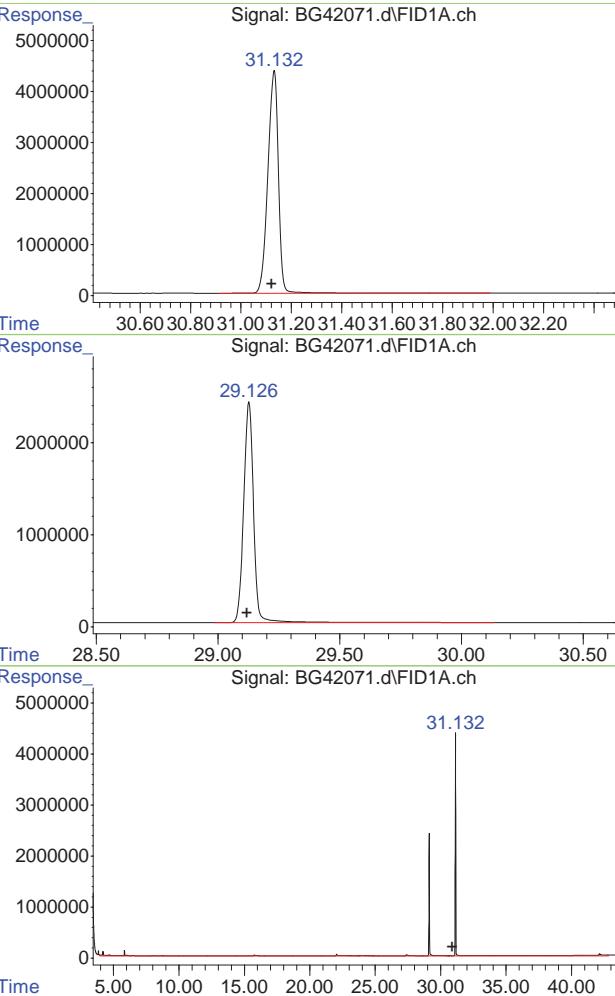
## Quantitation Report (QT Reviewed)

Data Path : G:\1\DATA\BG131119\  
 Data File : BG42071.d  
 Signal(s) : FID1A.ch  
 Acq On : 19 Nov 2013 6:03 pm  
 Operator : RubenP  
 Sample : OP35833-MB  
 Misc : OP35833,GBG1601,0.01,,2,1  
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Nov 20 13:50:21 2013  
 Quant Method : G:\1\METHODS\BG130925ALK-Front.m  
 Quant Title : n-C8 - n-C40 normal alkanes w/ isoprenoids  
 Response via : Initial Calibration



G130925ALK-Front.m Thu Nov 21 11:20:03 2013



#1 5a-Androstan

R.T.: 31.132 min  
 Delta R.T.: 0.011 min  
 Response: 127120354  
 Conc: 50.00  $\mu\text{g/mL}$

#2 o-Terphenyl

R.T.: 29.127 min  
 Delta R.T.: 0.009 min  
 Response: 67666792  
 Conc: 25.35  $\mu\text{g/mL}$

#41 TPH (C8-C40)

R.T.: 30.880 min  
 Delta R.T.: 0.000 min  
 Response: 75556542  
 Conc: 30.93  $\mu\text{g/mL}$