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Alameda County Environmental Health

ADDITIONAL SITE CHARACTERIZATION RESULTS 224 RICKENBACKER CIRCLE LIVERMORE, CALIFORNIA

SUBMITTED

TO

MR. ROBERT STRONG
SAN RAMON, CALIFORNIA

PREPARED

 \mathbf{BY}

ENGEO INCORPORATED

PROJECT NO. 7584.100.101

APRIL 8, 2008 REVISED JULY 23, 2008



Project No. **7584.100.101**

April 8, 2008 Revised July 23, 2008

Mr. Robert Strong 500 Bollinger Canyon Way, Suite A4 San Ramon, CA 94583

Subject:

224 Rickenbacker Circle Livermore, California

ADDITIONAL SITE CHARACTERIZATION RESULTS

Reference:

ENGEO Inc.; Revised Work Plan for Additional Site Characterization; 224 Rickenbacker Circle, Livermore, California; Revised July 9, 2007;

Project No. 7584.1.001.01.

Dear Mr. Strong:

ENGEO Incorporated is pleased to present our findings for additional site characterization completed for 224 Rickenbacker Circle (Property) in Livermore, California (Figure 1). The report summarizes previous investigative activities along with work conducted December 2007 through March 2008 at the site. A copy of this report should be provided to the Alameda County Health Service Department for their review.

We are pleased to be of continued service to you on this project. If you have any questions, please contact us.

Very truly yours,

ENGEO Incorporated

kk/sm/jb

Shawn Munger, CHG, REAL

No. HG 413



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1.0 BACKGROUND

1.1 Site Description

The Property is a rectangular-shaped, level parcel of land, located at 224 Rickenbacker Circle in Livermore, California (Figure 1). The site is currently occupied by a single structure and paved parking areas. A landscaped area is located north of the building, along Rickenbacker Circle. The property is located within a commercial area. The Livermore Water Reclamation Plant is located approximately 750 feet to the west and the Livermore Municipal Airport is located approximately ½ mile west of the property. Residential development is located approximately 750 feet to the south and 1,500 feet to the east.

1.2 Site History

The Property was formerly operated as a dry cleaning facility that utilized a tetrachloroethene (PCE)-based machine. According to the property owner, approximately 10 years ago the PCE-based machine was replaced by an Exxon DF2000 clean solvent machine and subsequently a silicon-based machine. All equipment was removed from the building in October 2005. Based on a site reconnaissance, a former boiler room was located in the southeastern corner of the building and a conventional washing machine pad with a grated drain was observed just north of the boiler room. A concrete patch was visible on the floor, as indicated on Figure 2, which is the assumed sanitary sewer alignment. A sanitary sewer cleanout was visible between the building and Rickenbacker Circle.

In October 2005, JMK Environmental Solutions, Inc. advanced three soil borings to a depth of approximately 35 feet below the ground surface and recovered soil samples from each boring. Analytical results of the soil samples indicated the presence of PCE to the maximum depth

explored in the two borings nearest the dry cleaning machine location. Based on review of the laboratory results for the soil samples, several samples exhibited concentrations of PCE in excess of the San Francisco Bay Regional Water Quality Control Board (SFRWQCB) Environmental Screening Levels (ESLs) for vapor intrusion. Groundwater was not encountered during the investigation, and therefore, no groundwater samples were collected.

A copy of the report prepared by JMK Environmental Solutions, Inc. was submitted to the Alameda County Health Services Agency along with a request for Site/Case Closure. Alameda County issued a letter dated July 6, 2006, in response to the request for case closure, requesting a work plan to delineate the extents of contamination at the Property.

In February 2007, ENGEO submitted an interim site characterization report to Alameda County Environmental Health (ACEH). The report included results of a soil vapor assessment as requested by ACEH as the first part of additional site characterization. Results from the soil vapor assessment indicated elevated concentrations of volatile organic compounds (VOCs). Based on the results of the survey, five boring locations were approved by ACEH to delineate the extents of the PCE-impacted soil and possible groundwater impact. The boring locations were selected based on previous soil data presented by JMK Environmental, results of soil gas survey, and to address potential uncertainty in the direction of groundwater flow.

Five borings were advanced under the supervision of ENGEO in March 2007, and soil and grab groundwater samples were recovered from each location. VOCs and total petroleum hydrocarbons as gasoline and diesel were reported in several soil samples. The reported concentrations for VOCs were below the San Francisco Regional Water Quality Control Board's environmental screening levels (ESLs) for commercial soil to indoor air (Table E-1) and screening levels for commercial land use where groundwater is a current or potential drinking water resource (Table F-1a). Additionally, the reported concentrations for TPH as diesel and

motor oil did not exceed the ESLs for commercial direct exposure (Table K-2), and would not be expected to impact commercial use of the property.

Analytical results from the grab groundwater samples indicated the presence of VOCs and total petroleum hydrocarbons as diesel. PCE was reported at concentrations exceeding the Maximum Contaminant Level (MCL) of $5\,\mu g/L$, as established by the California Department of Health Services, in several of the grab groundwater samples. All other VOCs were reported at concentrations below their respective MCLs. Additionally, the reported concentration of TPH as diesel did not exceed the ESLs for groundwater as a potential drinking water source.

1.3 Geology

The Property is located within the Livermore-Amador Valley, an east-west trending, inland structural basin located in northeastern Alameda County. The valley covers approximately 42,000 acres and is surrounded primarily by north-south trending faults and hills of the Diablo Range.

The valley is partially filled with recent alluvial fan, stream and lake deposits of Pleistocene-Holocene Age. The alluvial fans range in thickness from a few feet along the margins to nearly 800 feet in the west-central portion of the valley. The Livermore Formation (Pleistocene age), found below the majority of the alluvium in the groundwater basin, consists of beds of clayey gravels and sands, silts, and clays that are unconsolidated to semi-consolidated and estimated to be 4,000 feet thick in the southern and western portion of the basin.

Native soil encountered during advancement of the deep boring near the southwestern corner of the Property consisted of six primary zones: (1) Silty clay/sandy clay from surface to approximately 27 feet below the ground surface (bgs); (2) Silty sand with some subrounded fine

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gravel from approximately 27 feet to approximately 42 feet bgs; (3) Silty clay from approximately 42 feet bgs to approximately 60 feet bgs; (4) Gravel grading to sand from approximately 60 feet bgs to approximately 72 feet bgs; (5) Silty clay from approximately 72 feet bgs to approximately 75 feet bgs; and (6) Gravel grading to sand from approximately 75 feet bgs to the bottom of the boring at 97.5 feet bgs.

1.4 <u>Hydrogeology</u>

A review of available hydrogeology information from the Zone 7 Water Agency indicates that although multiple aquifers have been identified in the Main Basin alluvium, wells have been classified generally as being in one of two aquifer zones, separated by a relatively continuous silty clay aquitard up to approximately 50 feet thick.

<u>Upper Aquifer Zone</u> - The upper aquifer zone consists of alluvial materials, including primarily sandy gravel and sandy clayey gravels. These gravels are usually encountered underneath the surficial clays (typically 20 to 40 feet below ground surface [bgs]) to about 80 to 150 feet bgs. Groundwater in this zone has been classified as generally unconfined.

<u>Lower Aquifer Zone</u> - All sediments encountered below the clay aquitard in the center portion of the basin have been known collectively as the Lower Aquifer Zone. The aquifer materials consist of semi-confined to confined, leaky, coarse-grained, water-bearing units interbedded with relatively impermeable, fine-grained units.

As described above, several unconfined water bearing zones appear to be present at the subject site, separated by relatively impermeable clay layers.



1.5 Scope of Services

Based on data gaps identified as part of the conceptual site model prepared during the work plan preparation, the scope of work provided by ENGEO for the additional site characterization consisted the following:

- Notifying Underground Service Alert to identify known subsurface utilities.
- Advancement of one deep boring with soil and groundwater sampling (Figure 2).
- Advancement of seven additional soil gas probes (Figure 2).
- Installation of three groundwater monitoring wells.
- Recovery of soil samples from each of the three monitoring well locations.
- Field screening of the soil samples for organic vapors during drilling activities using a photoionization detector.
- Analysis of the selected soil samples and groundwater samples for Volatile Organic Compounds (VOCs) and total petroleum hydrocarbons as gasoline, diesel, and motor oil.
- Preparation of a Utility Corridor Assessment.
- Preparation of this final report documenting the field and laboratory activities.

2.0 FIELD ACTIVITIES, LABORATORY TESTING AND RESULTS

2.1 Deep Soil Boring

As requested by ACEH, one deep soil boring, 1-B1, was advanced in the southwestern portion of the site (Figure 2). Field activities were conducted on November 27 and 28, 2007, using a Mini-Sonic sampling rig. The boring was advanced to approximately 100 feet below the ground surface and logged in the field by an ENGEO Staff Geologist. The boring log is presented in Appendix A.

2.1.1 Soil Sampling

The Mini-Sonic was used to recover soil samples in 1½-inch-diameter sample cores in clear acrylic tubes. The acrylic sampling tubes were then cut to collect the soil sample at the desired depth. During sampling, retrieved soils were screened for organic vapors using a portable photoionization detector. All sampling equipment was regularly decontaminated using appropriate controls and protocols. Soil samples for analyses were sealed with Teflon sheets secured by tight fitting plastic end caps and tape, and were labeled to indicate a unique sample number, sample location, and time and date collected.

The soil samples from Boring 1-B1 were analyzed for the following:

- Volatile Organic Compounds (VOCs) by EPA Method 8260B.
- Total Petroleum Hydrocarbons (TPH) as gasoline, diesel and motor oil by EPA Method 8015B.

The laboratory analytical reports prepared by TestAmerica are included in Appendix B. A summary of the soil analytical results are presented in Tables 1 and 2.

PCE was reported above laboratory detection limits within Boring 1-B1 at depths of 10 feet, 20 feet and 50 feet below the ground surface. Reported concentrations range from below laboratory detection limits to 0.079 milligrams per kilogram (mg/kg). The environmental screening level (ESL) as established by the San Francisco Regional Water Quality Control Board (SFRWQCB) for commercial/industrial land use deep soil where groundwater is a potential source of drinking water is 0.70 mg/kg. No other VOCs were reported above laboratory detection limits in the soil samples recovered from the deep boring.

Total petroleum hydrocarbons as diesel were reported in the soil samples recovered from depths of 20 and 50 feet below the ground surface at concentrations of 17 mg/kg and 1.1 mg/kg, respectively. The reported concentrations for TPH as diesel do not exceed either residential or commercial ESLs. TPH gasoline and motor oil were not reported above laboratory detection limits in any of the soil samples recovered from the deep boring.

2.1.2 Groundwater Sampling

Groundwater samples were recovered from Boring 1-B1 at discrete depths from zones identified as water-bearing layers from the deep boring using a modified hydropunch tool. The identified depths were 35 feet, 70 feet and 90 feet below the ground surface. The grab-groundwater samples were recovered using a dedicated polyethylene tube equipped with a check valve. Following recovery, the groundwater samples were decanted into appropriate laboratory glassware. The samples were labeled to indicate a unique sample number, sample location, and time and date collected. All soil and groundwater samples were placed in an ice-cooled chest for delivery to TestAmerica in Pleasanton, California.

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The groundwater samples were analyzed for the following:

Volatile Organic Compounds (VOCs) by EPA Method 8260B.

• Total Petroleum Hydrocarbons (TPH) as gasoline, diesel and motor oil by EPA Method

8015B.

The laboratory analytical reports prepared by TestAmerica Laboratories, Inc. are included in

Appendix B. A summary of the groundwater analytical results are presented in Table 3.

VOCs detected in the groundwater samples include benzene, ethylbenzene, toluene and xylenes.

Reported concentrations for the constituents in the groundwater sample recovered from 70 feet

below the ground surface was 2.7 micrograms per liter (ug/L), 1.2 ug/L, 2.3 ug/L and 1.0 ug/L

respectively. The reported concentrations were compared to the groundwater concentrations for

evaluation of potential vapor intrusion concerns (Table E-1) ESL and the Maximum

Contaminant Level (MCL) of as established by the California Department of Health Services.

None of the detected VOCs were found to exceed the established screening levels. VOCs were

not reported in the other two groundwater samples recovered from the deep boring. No

chlorinated VOCs were reported for the groundwater samples recovered from the deep boring.

Total Petroleum Hydrocarbons (TPH) as gasoline and diesel was reported in the groundwater

sample recovered from 70 feet below the ground surface at concentrations of 98 ug/L and

190 ug/L respectively. The reported concentrations are below the ESLs for groundwater as a

potential source of drinking water. TPH as gasoline and diesel were not reported above

laboratory detection limits for the two other samples recovered from the deep boring. TPH as

motor oil was not reported above laboratory detection limits in any of the groundwater samples.

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2.2 Soil Vapor Sampling

Seven additional soil vapor probes were advanced on December 17, 2007. The soil vapor samples were collected using protocols and procedures consistent with "Advisory-Active Soil Gas Investigations dated January 13, 2003" (Los Angeles Regional Water Quality Control Board and DTSC). Significant rainfall, as defined in the above listed document as ½ inch or greater, was not observed within 48 hours prior to sampling. The GeoprobeTM direct-push soil vapor probes utilized a hydraulic hammer to drive a 2-inch-diameter rod to approximately 5 feet below the ground surface, and a bentonite seal was applied between the drive rod and ground surface. A disposable drive tip at the end of the rod was fitted with disposable poly tubing to which a laboratory-assembled sampling manifold was attached.

One purge and one 6-liter sample canister (Summa) were securely attached to the manifolds. Each manifold consisted of a flow regulator and moisture filter that was used solely for one soil vapor probe. Each probe was allowed to equilibrate for approximately 30 minutes following installation. A leak test was performed for approximately 10 minutes by bringing the sampling manifold up to vacuum and observing any loss in pressure. After the 30-minute equilibration period, qualitative soil permeability observations were completed, and the entire sampling train was purged using seven purge volumes. A purge volume test was conducted during previous soil gas sampling activities at the site, and seven purge volumes was shown to produce the highest resultant concentrations of target analytes. Qualitative soil permeability observations were conducted by observing the changes in the canister pressures during purging. The volume of the sampling system was approximately 600 milliliters (mL) and was equipped with a flow regulator set for 200 milliliter per minute (mL/min). Based on the number of purge volumes and the flow rate, each location was purged for approximately 3.15 minutes. Following purging, samples were collected in the 6-liter Summa canisters. Samples times varied per location due to

differences in the lithology at each location. Table 5 summarizes the Summa canister pressures prior to and after sampling, as well as upon arrival to the laboratory.

For quality assurance purposes, a leak detection compound, rubbing alcohol (2-Propanol), was applied to gauze strips attached to system connections and seals along the sampling manifold. Additional gauze strips were placed around the bentonite seal at the ground surface, as well as at the top of the probe. Probe rods were decontaminated with a non-phosphate detergent and a triple wash between each soil boring location.

The collected samples were returned to Torrent Laboratory, Inc., a fixed-base laboratory in Milpitas, California under documented chain-of-custody. The samples were analyzed using EPA Method TO-15 for volatile organic compounds (VOCs).

The laboratory analytical reports as prepared by Torrent Laboratory, Inc. are presented in Appendix C. These values have been compared to the ESLs for evaluation of potential indoor air impacts. Volatile organic compounds were detected in all soil vapor samples collected; however, the detected concentrations were below ESLs. PCE in particular was reported at concentrations ranging from below laboratory detection limits to 64 ug/m³, which are well below the ESL (1,400 ug/m³) for commercial vapor intrusion. Analytical reports prepared by Torrent Laboratory, Inc. are included as Appendix C. A summary of the groundwater analytical results are presented in Table 3.

2.3 Monitoring Well Installation and Sampling

Three groundwater monitoring wells were installed at locations MW-1 through MW-3 (Figure 2). The selected location of the wells was based upon discussions with ACEH. The well

borings were advanced using a truck-mounted drill rig equipped with 8½-inch-diameter hollow-stem augers on December 18, 2007.

The monitoring wells consist of 2-inch-diameter PVC casing with flush joints. The wells were constructed with approximately 25 feet of screened casing (0.01-inch slot width) and an appropriate length of solid PVC well casing (2-inch-diameter Schedule 40 PVC). The total depth of the monitoring wells is approximately 35 feet. A No. 2/16-sand-filter pack was placed from the base of the well to approximately one foot above the top of the screened interval. A ±12-inch-thick bentonite seal was placed at the top of the filter pack. The remaining annular space was backfilled with a cement-bentonite grout seal. The well was completed with flush-mounted monument. The top of the well casings were surveyed and secured with a locking waterproof cap. The drill cuttings were placed within sealed 55-gallon drums, and upon review of the laboratory analyses, a disposal plan for the soil cuttings will be developed.

On January 16, 2008, the wells were developed through purging to maximize the hydraulic connection between the well and the surrounding aquifer material and remove fine sediment to produce relatively non-turbid groundwater. Approximately 20 well volumes of water were removed during the development process. The purged water was placed on-site in Department of Transportation approved drums.

2.3.1 Soil Sampling

During monitoring well installation, soil samples were recovered at 5-foot intervals to the base of the boring and screened in the field with a photoionization detector (PID). Soil samples for laboratory analysis were recovered from locations exhibiting significant PID readings and at the top of the saturated zone.

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The soil samples were analyzed for the following:

• Volatile Organic Compounds (VOCs) by EPA Method 8260B.

• Total Petroleum Hydrocarbons (TPH) as gasoline, diesel and motor oil by EPA Method

8015B.

The laboratory analytical reports prepared by TestAmerica are included in Appendix B. A

summary of the soil analytical results are presented in Tables 1 and 2.

PCE was reported above laboratory detection limits in the soil samples recovered from MW-1 at

depths of 5½ and 10½ feet below the ground surface at concentrations of 0.081 mg/kg and

0.068 mg/kg respectively. The environmental screening level (ESL) as established by the

San Francisco Regional Water Quality Control Board (SFRWQCB) for commercial/industrial

land use where groundwater is a potential source of drinking water is 0.70 mg/kg for both

shallow and deep soils. No other VOCs were reported above laboratory detection limits in the

soil samples recovered from the monitoring well locations.

Total petroleum hydrocarbons as diesel was reported in the soil sample recovered from MW-3 at

a depth of 26 feet below the ground surface, at a concentration of 2.2 mg/kg. The reported

concentrations for TPH as diesel do not exceed the ESLs for commercial/industrial land use.

TPH diesel was not reported above laboratory detection limits in the other soil samples recovered

from the monitoring well locations and TPH as gasoline and motor oil were not reported above

laboratory detection limits in any of the soil samples recovered from the monitoring well

locations.

2.3.2 Groundwater Sampling

On January 28, 2008, the three wells were purged using a 12-volt submersible pump. Water

quality parameters including dissolved oxygen (DO), pH, and ORP were measured using a

hand-held Hanna 9828 Multi Parameter Water Quality Meter with flow through cell. The

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samples were decanted into pre-cleaned laboratory glassware and cooled in an ice chest until delivery under documented chain-of-custody to TestAmerica in Pleasanton, California. Data summary sheets prepared during the groundwater sampling are included as Appendix D.

The groundwater samples were analyzed for the following:

- Total Petroleum Hydrocarbons (TPH) as diesel and motor oil by EPA Method 8015B.
- Volatile Organic Compounds (VOCs) by EPA Method 8260B.

The laboratory analytical reports prepared by TestAmerica Laboratories, Inc. are included in Appendix B. A summary of the groundwater analytical results are presented in Table 3 and a table reporting the depths to groundwater is presented in Table 6.

PCE was reported above laboratory detection limits in the groundwater samples recovered from MW-1 and MW-2. Reported concentrations for the wells were 0.80 ug/L and 0.98 ug/L respectively. The reported concentrations were compared to the groundwater concentrations for evaluation of potential vapor intrusion concerns (Table E-1) ESL and the MCL (5 ug/L). The reported concentrations did not exceed the screening levels. No other VOCs were reported above laboratory detection limits.

The monitoring wells were surveyed by a PLS Surveys Inc. in February 2008. A copy of the report prepared by PLS Surveys Inc. is included as Appendix D. The initial monitoring of the three monitoring wells currently on Site indicates that the groundwater flow direction is to the northwest. Additionally, based on the initial depth to groundwater readings of the onsite wells, the groundwater gradient beneath the Site has been determined to be approximately 0.006 feet per foot (ft/ft). A groundwater elevation contour map is included as Figure 3.



3.0 UTILITY CORRIDOR ASSESSMENT

A utility corridor assessment was conducted in accordance with the December 15, 2004 DTSC Guidance for the Evaluation and Mitigation of Subsurface Vapor Intrusion to Indoor Air. As part of the assessment, facility records including copies of the utility maps, building "as-built' drawings, building construction specifications, and municipal utility corridor maps were reviewed as available.

Information available at the City of Livermore indicated that the storm drain, sanitary sewer and drinking water mains extend the length of Rickenbacker Circle, near the center of the street section. Plans indicate that a 24-inch storm drain is located approximately 10 feet bgs, an 8-inch sanitary sewer line is located approximately 6 feet bgs, and an 8-inch water main is located approximately 4 feet bgs. Based on a review of available 'as-built' plans, all utilities enter the site from Rickenbacker Circle. No information was readily available pertaining to the depth to gas and electrical lines; however, a contact in the City of Livermore Engineering Department indicated that as a general rule, no more that 24 inches of cover are recommended for gas and electrical lines.

The depth to groundwater at the site has been measured at approximately 25 feet below the ground surface. Based on the depth to groundwater, it is unlikely that the utility corridors will act as pathways for groundwater contaminant migration.

Soil vapor data collected from the locations near utility trenches originating within the building (SG-7, SG-10, SG-11 and SG-12) show decreasing concentrations of PCE as the distance from the former dry cleaning machine increases. Based on the analytical results and the location of the utility trenches at the Property, it appears that concentrations of constituents in the soil vapor decrease as the distance from the former dry cleaning machine increases; it is therefore unlikely that impacted soil vapor would migrate through on-site utility trenches to the street at concentrations that would adversely impact downgradient sites.



4.0 SITE CONCEPTUAL MODEL UPDATE

4.1 Current Soil Vapor, Soil and Groundwater Concentrations

4.1.1 Soil Vapor

Soil vapor samples were recovered on two separate occasions, in January 2007 and December 2007. The January 2007 samples were recovered by TEG-Northern California using the "syringe" method and analyzed by means of an on-site mobile laboratory using EPA method 8260. The December 2007 soil vapor sampling was conducted using conventional methods outlined in the "Advisory-Active Soil Gas Investigations document dated January 13, 2003" (Los Angeles Regional Water Quality Control Board and DTSC). The December 2007 samples were recovered in 6-liter Summa canisters with TO-15 analysis. A summary of soil vapor analytical results are presented in Table 4. Analytical results are compared to the San Francisco Regional Water Quality Control Board's (SFRWQCB's) environmental screening levels (ESLs¹) for evaluation of potential indoor air impacts (Table E). A number of soil vapor samples exceeded the respective commercial ESLs

The soil vapor results are further discussed in section 5.1.

4.1.2 Soil Samples

Several VOCs have been detected in soil samples recovered from the Property including acetone, cis 1,2 dichloroethene, trans-1,2-dichloroethene, trichloroethene, and PCE. Acetone was detected in

¹ San Francisco Bay Regional Water Quality Control Board; November 2007, Interim Final, Screening for Environmental Concerns at Site with Contaminated Soil and Groundwater, Volume 1, Table E, Indoor Air and Soil Gas, Commercial/Industrial Land Use Exposure Scenario.



sample S-3 at a depth of 4 feet below the ground surface (bgs) at a concentration of 0.049 mg/kg. In sample S-3 at 2 and 4 feet bgs, cis-1,2-dichloroethene was detected at concentrations of 0.054 mg/kg to 0.061 mg/kg and trans-1,2-dichloroethene was detected at concentrations of 0.015 mg/kg to 0.0065 mg/kg, respectively. PCE was detected in sample S-3 at depths of 4 and 8 feet bgs at concentrations of 0.013 mg/kg and 0.0066 mg/kg respectively. Additionally, PCE was detected at concentrations ranging from 0.012 to 0.079 mg/kg in sample S-3 at 4, 8, and 10 feet bgs, and in boring MWB1 at depths of 5½ and 10½ ft bgs at concentrations of 0.081 and 0.068 mg/kg respectively. PCE was reported in the deep boring, 1-B1, at depths of 10, 20 and 50 ft bgs at concentrations ranging from 0.014 to 0.079 mg/kg. The reported VOC concentrations were not found to exceed established ESLs for commercial/industrial land use where groundwater is a potential source of drinking water. No VOCs were reported above laboratory detection limits for samples S-1, S-2, S-3, S-4 and S-5 recovered from the saturated zone at each sample location.

Total Petroleum Hydrocarbons (TPH) as diesel were reported at concentrations ranging from 1 to 17 mg/kg in soil samples recovered at sampling locations S-2, S-3, S-5, 1-B1 and MWB3@26. Two soil samples were recovered from the near the drain in the former boiler room, and reported concentrations were found to range from below laboratory detection limits to 190 mg/kg (P-1 at 5 feet). TPH as gasoline was reported at a concentration of 0.33 mg/kg at sample location S-3 at a depth of 2 feet bgs. TPH as motor oil was not reported above laboratory detection limits in any of the samples. The reported concentration of TPH as diesel recovered from 5 feet bgs at P-1, 190 mg/kg, exceeds the established ESL for commercial/industrial land use where groundwater is a potential source of drinking water. No other constituents were reported at concentrations exceeding the established ESLs. Analytical results for the shallow soil samples (recovered from 0 to 10 feet below the ground surface) are presented in Table 1 and deep soil samples (recovered from greater than 10 feet) are presented in Table 2.



4.1.3 Groundwater Samples

VOCs detected in the groundwater samples include benzene, ethylbenzene, cis-1,2-dichloroethene, trichloroethene, toluene, xylenes and PCE. Benzene, ethylbenzene and xylenes were reported in the groundwater sample recovered from the deep boring at 70 ft bgs (DB-1(GW@70)), at concentrations of 2.7 ug/L, 1.2 ug/L and 1.0 ug/L respectively. Cis-1,2-dichloroethene was reported in samples S-3 and S-5 at concentrations of 1.6 micrograms per liter (µg/L) and 0.54 µg/L. Additionally, trichloroethene was reported in samples S-3 and S-5 at concentrations of 2 µg/L and 2.2 µg/L respectively. Analytical results indicated the presence of toluene in samples S-3 through S-5, and DB-1(GW@70) at concentrations ranging from 0.86 µg/L to 2.3 µg/L. PCE was reported in samples S-2 through S-5, at concentrations ranging from 1.8 µg/L (sample S-2) to 36 µg/L (sample S-5). PCE was reported in the groundwater samples recovered from monitoring wells MW-1 and MW-2 at concentrations of 0.8 µg/L and 0.98 µg/L respectively. The reported concentrations were compared to the groundwater concentrations for evaluation of potential vapor intrusion concerns (Table E-1) ESL and the MCLs. Concentrations of PCE in the grab samples recovered from S-3, S-4 and S-5 were found to exceed the established MCL of 5 µg/L, but are below the ESL of 420 µg/L. No other constituents were reported above the MCL or ESL (Table E) screening levels.

Total Petroleum Hydrocarbons (TPH) as diesel was reported in the groundwater samples recovered at sample location S-4 and DB-1(GW@70) at concentrations of 70 μ g/L and 190 μ g/L. TPH as gasoline was reported in the groundwater samples recovered from DB-1(GW@70) at a concentration of 98 μ g/L. TPH as motor oil was not reported above laboratory detection limits in any of the groundwater samples. No reported concentrations of TPH were found to exceed the established MCLs or ESLs (Table E-1). Groundwater analytical results are summarized in Table 3.



4.2 Sensitive Receptors

As part of the Soil and Groundwater Sampling Results report, a request was submitted to Zone 7 Water Agency to provide a ½-mile radius well search. Zone 7 provided a map depicting well locations (Figure 2), and two wells were identified approximately 360 feet southeast of the Property. In preparation of this work plan, additional information regarding the two wells was requested from Zone 7. According to Mr. Wyman Hong, the two wells located southeast of the Property are no longer in existence. Well 3S/2E-7C1 (open blue triangle) was destroyed on December 7, 1983, and it was an agricultural well with a 14-inch-diameter casing and 280 feet deep. Well 3S/2E-7D1 (yellow cross) cannot be located and most was likely destroyed by construction; this well was documented in a San Francisco Water Department report dated 1912. No other information regarding these two wells is available.

The nearest residence is located approximately 950 feet south of the Property, and Rancho Las Positas Elementary School is located approximately 1300 feet southeast of the site at 401 E Jack London Boulevard. Lake Del Valle, the local reservoir used for drinking water supply, is located approximately 6 miles southeast of the Property.

As described below, three monitoring wells were installed at the Property. Based on the initial depth to water readings, the groundwater flow direction at the Property appears to be to the northwest. Additionally, based on the initial readings of the onsite wells, the groundwater gradient beneath the Site has been determined to be approximately 0.006 feet per foot (ft/ft).

4.3 Exposure Pathways

A flowchart was created to evaluate potential exposure pathways, as shown on Figure 4. The site is currently occupied by a single structure located in a commercial area. Impacts from the former dry

cleaning facility are present in shallow soil below the building, in groundwater below and to the southwest of the building, and soil vapor throughout portions of the site. The primary chemical of

concern, PCE, was stored and used onsite as part of the previous dry cleaning operations. It is

suspected that surface spills of PCE and discharge of PCE-impacted wastewater were the release

mechanisms that allowed PCE to enter the subsurface soil beneath the site. Through leaching, PCE

has migrated from soil to groundwater beneath the site. Through volatilization of the impacts in soil

and groundwater, soil vapor has been also impacted.

According to the well search performed by Zone 7 Water Agency, there are no domestic wells

within ½-mile of the site. This information was confirmed by a windshield survey completed by

ENGEO on May 22, 2007. Given the lack of nearby domestic wells and the limited impacts to

groundwater at the site, it is unlikely that the ingestion, direct contact, or inhalation exposure

pathways are complete for groundwater.

Inhalation of impacted vapors created from volatilization within the building was identified as a

complete exposure pathway. We believe the potential for exposure to vapor inhalation is limited

due to (1) industrial use of the site, (2) an existing passive ventilation system present in the

structure, and (3) the relatively low concentrations reported from the December 2007 soil gas

sampling.

The nearest surface water body identified is Lake Del Valle, the local reservoir used for drinking

water supply, located approximately 6 miles southeast of the site. Due to the large distance

between the site and Lake Del Valle, the potential migration of impacts to surface water or

sediments was found to be incomplete.

7584.100.101

April 8, 2008

Revised July 23, 2008

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5.0 DISCUSSION

5.1 Soil Vapor Analysis

The previous field soil vapor screening assessment conducted in January 2007 found significantly higher soil vapor results than the recent fixed based laboratory analyses. A specific reason for the noted discrepancy between the two soil vapor data sets is unknown. TEG used a "syringe" method for the recovery of the soil vapor samples with on-site mobile laboratory analysis using EPA 8260 methodology. It is also conceivable that syringes used for sample collection may have not been properly purged prior to sampling and may exhibit cross-contamination from other off-site sources. In addition, the EPA 8260 method used for mobile analysis was not specifically developed for air analysis. Discrepancies between 8260 and TO-15 analysis have been demonstrated in the past.

Review of the analytical results prepared by the fixed-base laboratory (Torrent) indicate that the air samples were diluted prior to analytical analysis. Discussions with the laboratory indicate that the use of a dilution does not bias the final analytical results, only raises the minimum reporting limit. A copy of the correspondence with the laboratory can be found in Appendix C.

Additional evaluation of soil vapor at the site is recommended to address the discrepancies between the January 2007 and December 2007 soil vapor sampling events.

5.2 Deep Boring

Review of soil and groundwater data from the deep boring (1-B1) found no significant constituents of concern. Total petroleum hydrocarbons and BTEX were reported at concentrations below ESLs/MCLs in the groundwater samples recovered at 70 feet; however,



given the lack of TPH/BTEX in the shallow groundwater samples, we would attribute the TPH/BTEX to off-site source(s).

5.3 Groundwater Monitoring Well Sampling

Low level PCE was reported for wells MW-1 and MW-2 at 0.8 ug/l and 0.98 ug/l, respectively. Based on groundwater elevation measurements, well MW-3 is located at the downgradient area of the site, with no PCE or other VOCs reported. This data suggest the PCE impacts in groundwater are limited in extent.



6.0 CONCLUSIONS

Several VOCs and total petroleum hydrocarbon were detected in the soil at the Property at concentrations below the environmental screening levels established by the SFRWQCB for commercial soil to indoor air (Table E-1) and for commercial land use where groundwater is a current or potential drinking water resource (Table F-1a). Soil impacts are generally limited to the area in the vicinity of the former dry cleaning machine and to a depth of approximately 20 feet below the ground surface.

Several VOCs were detected in the groundwater beneath the Property. Reported concentrations of PCE exceed the Maximum Contaminant Level of 5 ug/L as established by the California Department of Health Services for grab groundwater samples collected in March 2007; however, groundwater samples recovered from the monitoring wells in January 2008 did not report concentrations of PCE exceeding the established MCLs. Based on the depth to groundwater measurements recorded in January 2008, the direction of groundwater flow is to the northwest at a gradient of approximately 0.006 ft/ft. Analytical results from the sample location located northwest of the former dry cleaning machine (S-2) reported PCE concentrations in the groundwater less than the established MCL. In addition, the sample collected from the downgradient well MW-3 found no detectable PCE or VOCs, suggesting a plume of limited extent.

Initial screening level soil gas sampling conducted in January 2007 found significant levels of VOCs in soil gas; however, subsequent soil gas recovery in December 2007, using preferred sampling methodologies and test methods, found no significant soil gas impacts. While the soil gas data appears fundamentally in conflict, we believe the Summa canister recovery and TO-15 methodology to be representative of site conditions.



7.0 LIMITATIONS

We performed our professional services in accordance with generally accepted environmental engineering principles and practices currently employed in Northern California at the time of this report. No other warranty is expressed or implied.

We limited our investigation to the authorized scope of work. Our investigation is not intended to be comprehensive, to identify all potential concerns, or to guarantee that no additional environmental contamination beyond that described in this report exists at the site.

The findings in this report are valid as of the time of the investigation; however, changes in subsurface conditions can occur over time, whether due to natural processes or human activity on the Property or on surrounding properties. ENGEO Incorporated has prepared this report for the exclusive use of Mr. Robert Strong. It is recognized and agreed that ENGEO has assumed responsibility only for undertaking the study for the client. The responsibility for disclosures or reports to a third party and for remedial or mitigative action shall be solely that of the Client.



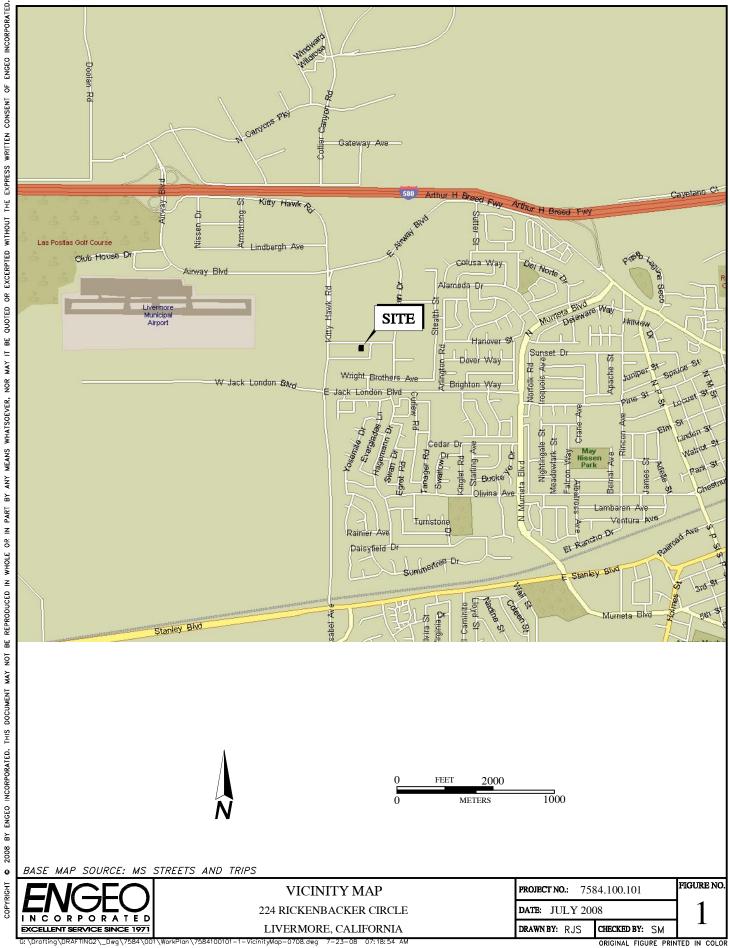
LIST OF FIGURES

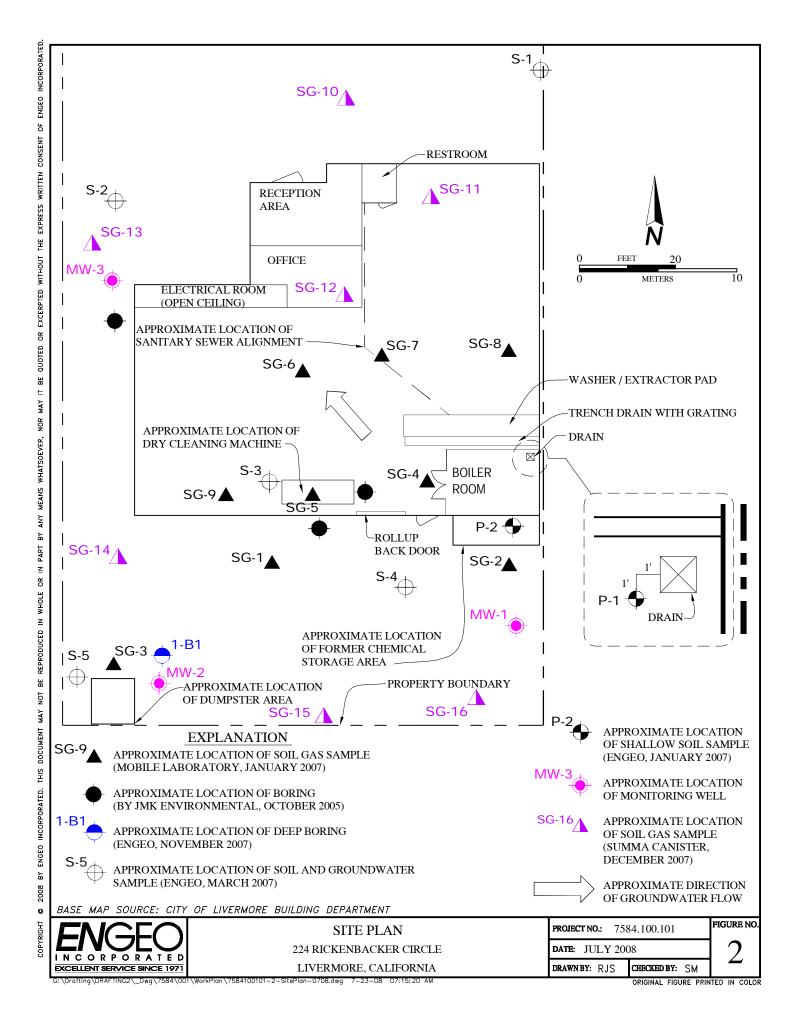
Figure 1 Vicinity Map

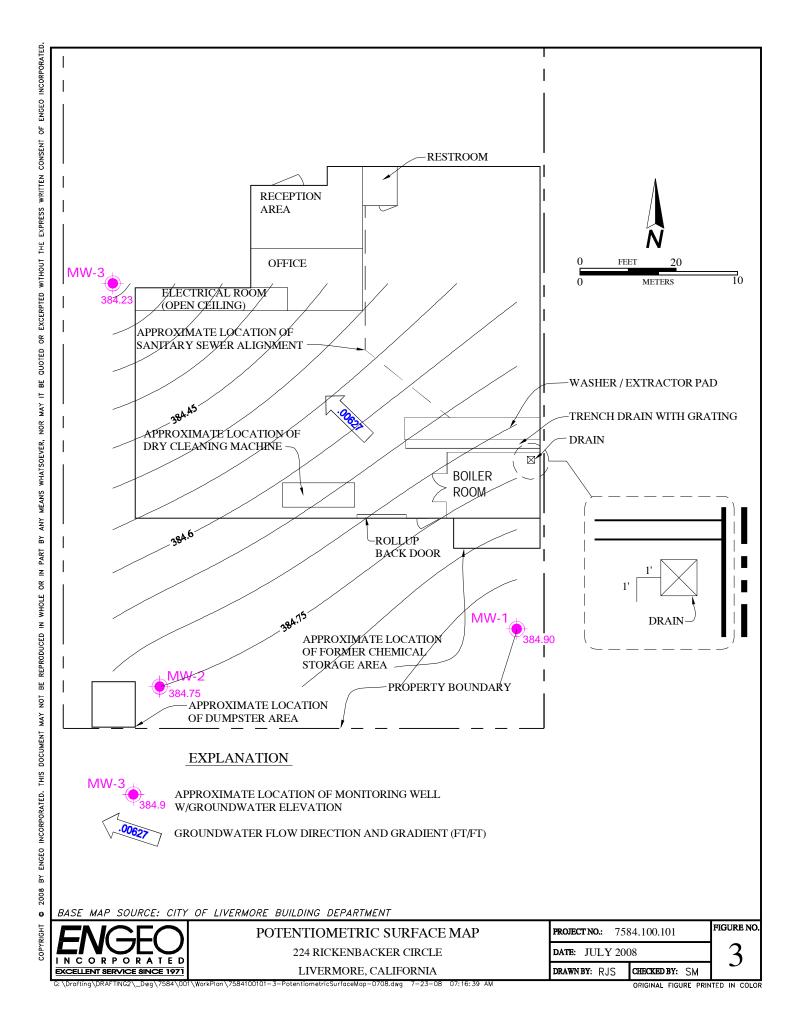
Figure 2 Site Plan

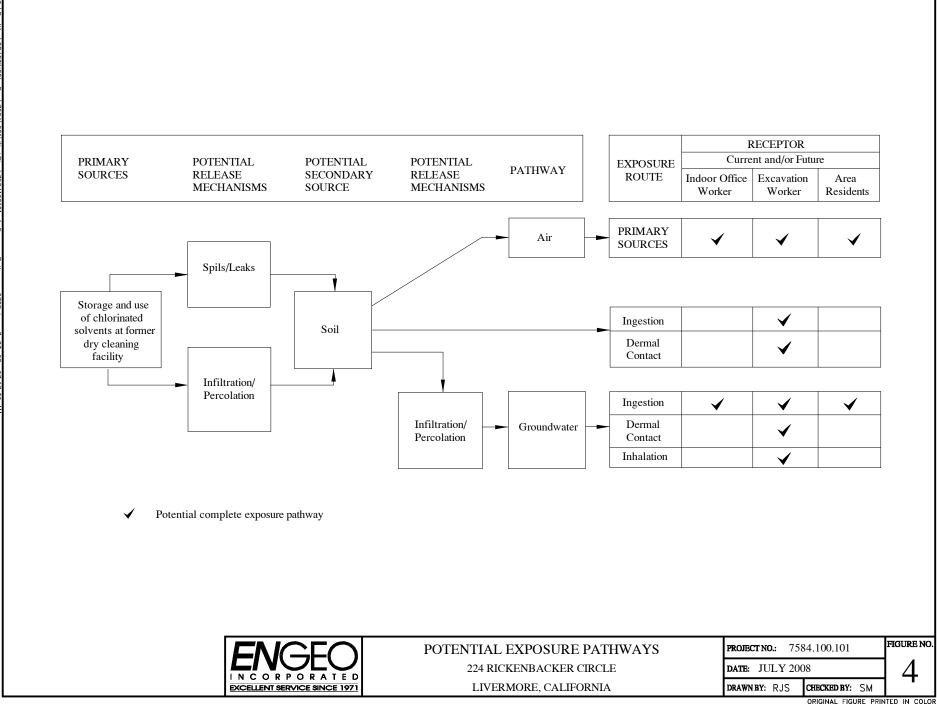
Figure 3 Potentiometric Surface Map

Figure 4 Potential Exposure Pathways











TABLES

Table 1	Shallow Soil Data
Table 2	Subsurface Soil Data
Table 3	Groundwater Data
Table 4	Soil Vapor Data
Table 5	Summa Canister Pressures
Table 6	Monitoring Well Depth to Groundwater

Table 1. Shallow Soil Data

	SFRWQCB	P-1	P-1	P-2	P-2	S-3@2'	S-3@4'	S-3@8'	MWB1@5-1/2'
		Soil	Soil	Soil	Soil	Soil	Soil	Soil	Soil
	ESL	1 ft	5 ft	1 ft	5 ft	2 ft	4 ft	8 ft	5 1/2 ft
		1/22/2007	1/22/2007	1/22/2007	1/22/2007	3/1/2007	3/1/2007	3/1/2007	12/18/2007
	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg
TARGET ANALYTE									
	TABLE A (C/I)								
ACETONE	2.1	0.062	< 0.049	< 0.050	< 0.047	<0.048	0.049	<0.048	< 0.047
cis-1,2-DICHLOROETHENE	0.19	<0.0048	< 0.0049	< 0.0050	< 0.0047	0.054	0.061	<0.0048	< 0.0047
trans-1,2-DICHLOROETHENE	0.67	<0.0048	< 0.0049	< 0.0050	< 0.0047	0.0065	0.015	<0.0048	< 0.0047
TETRACHLOROETHENE	1	<0.0048	0.0055	<0.0050	< 0.0047	<0.0048	0.012	0.079	0.081
TRICHLOROETHENE	0.46	<0.0048	< 0.0049	< 0.0050	< 0.0047	<0.0048	0.013	0.0066	< 0.0047
TPH-GASOLINE	83	<0.24	< 0.23	<0.24	<0.25	0.33	<0.23	<0.24	<0.23
TPH-MIDDLE DISTILLATES	83	2.6	190	2.9	< 0.99	4.5	1.0	< 0.96	<1.0
TPH-RESIDUAL FUELS	2,500	<48	1000	<49	<50	<50	<47	<46	<50

Table 2. Subsurface Soil Data

	SFRWQCB	S-1@24'	S-2@26'	S-3@10'	S-3@27'	S-4@25'	S-5@30'	S@10'	S@20'	S@30'	S@40'	S@50'	S@60'
		Soil	Soil	Soil	Soil	Soil	Soil						
	ESL	24 ft	26 ft	10 ft	27 ft	25 ft	30 ft	10 ft	20 ft	30 ft	40 ft	50 ft	60 ft
		3/2/2007	3/2/2007	3/2/2007	3/2/2007	3/2/2007	3/2/2007	11/27/2007	11/27/2007	11/27/2007	11/27/2007	11/27/2007	11/27/2007
	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg
TARGET ANALYTE													
	TABLE C (C/I)												
TETRACHLOROETHENE	0.70	<0.0045	< 0.0049	0.023	< 0.0047	< 0.0049	<0.0048	0.079	0.017	< 0.0049	< 0.0050	0.014	<0.0050
TPH-MIDDLE DISTILLATES	83	< 0.96	11	13	< 0.99	<0.98	1.0	<1.0	17	< 0.99	< 0.99	1.1	<0.99

Table 2. Subsurface Soil Data

_							
	SFRWQCB	S@70'	S@80'	S@90'	MWB1@10-1/2'	MWB2@25-1/2'	MW3@26'
		Soil	Soil	Soil	Soil	Soil	Soil
	ESL	70 ft	80 ft	90 ft	10 1/2 ft	25 1/2 ft	26 ft
		11/27/2007	11/27/2007	11/27/2007	12/18/2007	12/18/2007	12/19/2007
	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg
TARGET ANALYTE							
	TABLE C (C/I)						
TETRACHLOROETHENE	0.70	< 0.0049	< 0.0049	<0.0048	0.068	< 0.0050	< 0.0046
TPH-MIDDLE DISTILLATES	83	<0.98	<1.0	< 0.99	<1.0	<1.0	2.2

Table 3. Groundwater Data

		SFRWQCB	S-1	S-2	S-3	S-4	S-5	DB-1(GW@35')	DR-1/GW@70'\	DR-1/GW@00'\	MW-1	MW-2	MW-3
	1401							. ,	,	, ,			
	MCL	ESL	Water	Water	Water	Water	Water	Water	Water	Water	Water	Water	Water
			3/2/2007	3/2/2007	3/2/2007	3/2/2007	3/2/2007	11/28/2007	11/28/2007	11/28/2007	1/28/2008	1/28/2008	1/28/2008
	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
TARGET ANALYTE													
		TABLE E-1											
BENZENE	1	1,800	<1.0	<1.0	< 0.50	< 0.50	< 0.50	< 0.50	2.7	< 0.50	< 0.50	< 0.50	< 0.50
cis-1,2-DICHLOROETHENE	6	17000	<1.0	<1.0	1.6	< 0.50	0.54	< 0.50	< 0.50	<0.50	<0.50	< 0.50	< 0.50
ETHYLBENZENE	300	170,000	<1.0	<1.0	< 0.50	< 0.50	< 0.50	< 0.50	1.2	< 0.50	<0.50	< 0.50	< 0.50
TETRACHLOROETHENE	5	420	<1.0	1.8	27	16	36	< 0.50	< 0.50	<0.50	0.8	0.98	< 0.50
TRICHLOROETHENE	5	1,800	<1.0	<1.0	2.2	< 0.50	2	< 0.50	< 0.50	< 0.50	<0.50	< 0.50	< 0.50
TOLUENE	150	530000	<1.0	<1.0	0.86	0.96	1.8	< 0.50	2.3	< 0.50	<0.50	< 0.50	< 0.50
XYLENE(S)	1750	160,000	<2.0	<2.0	<1.0	<1.0	<1.0	<1.0	1.0	<1.0	<1.0	<1.0	<1.0
TPH - GASOLINE			<50	<50	<50	<50	<50	<50	98	<50	NA	NA	NA
TPH - MIDDLE DISTILLATES			<50	<50	<50	70	<50	<50	190	<50	NA	NA	NA

Bold Numbers - Exceed MCL or ESL Concentrations
-- - Screening level not provided
NA - Constituent not reported by laboratory

Table 4. Soil Vapor Data

	SFRWQCB	SG-1	SG-2	SG-3	SG-4	SG-5	SG-6	SG-7	SG-8	SG-9	SG-10	SG-11	SG-12	SG-12 (DUPLICATE)	SG-13	SG-14	SG-15	SG-16
		Soil Gas	Soil Gas	Soil Gas	Soil Gas	Soil Gas	Soil Gas	Soil Gas	Soil Gas	Soil Gas	Soil Gas	Soil Gas	Soil Gas	Soil Gas				
	ESL	5 ft	5 ft	5 ft	5 ft	5 ft	5 ft	5 ft	5 ft	5 ft	5 ft	5 ft	5 ft	5 ft				
		1/22/2007	1/22/2007	1/22/2007	1/22/2007	1/22/2007	1/22/2007	1/22/2007	1/22/2007	1/22/2007	12/17/2007	12/17/2007	12/17/2007	12/17/2007	12/17/2007	12/17/2007	12/17/2007	12/17/2007
	ug/m³	ug/m³	ug/m³	ug/m³	ug/m³	ug/m³	ug/m³	ug/m³	ug/m³	ug/m³	ug/m³	ug/m³	ug/m³	ug/m³	ug/m³	ug/m³	ug/m³	ug/m³
TARGET ANALYTE																		
VOCs	TABLE E (C/I)																	
1,1-DICHLOROETHENE	160	<100	<100	<100	<100	4700	<100	<100	<100	<100	<1.3	<1.3	<1.2	<1.2	< 0.79	<1.2	<1.2	<1.2
1,1-DIFLOUROETHANE		<100	<100	<100	<100	<100	<100	<100	<100	<100	NR	NR	NR	NR	NR	NR	NR	NR
1,2,4-TRIMETHYLBENZENE		NR	NR	NR	NR	NR	NR	NR	NR	NR	<1.4	<1.4	<1.4	<1.3	< 0.89	2.5	<1.3	<1.3
2-BUTANONE	2,900,000			NR	NR	NR	NR	NR	NR	NR	58	77	31	28	42	< 0.69	29	110
2-HEXANONE				NR	NR	NR	NR	NR	NR	NR	5.3	<1.4	<1.3	<1.3	< 0.86	1.7	<1.3	<1.3
4-ETHYL TOLUENE				NR	NR	NR	NR	NR	NR	NR	<1.2	<1.2	<1.2	<1.1	< 0.74	1.9	<1.1	<1.1
ACETONE	1,800,000			NR	NR	NR	NR	NR	NR	NR	100	180	70	62	110	< 0.90	200	180
BENZENE	280	<100	<100	<100	<100	<100	<100	<100	<100	<100	2.8	3.5	2.5	2.2	3.1	<1.4	4.0	6.6
CIS-1,2-DICHLOROETHENE	20,000	<100	<100	17,000	450	780,000 (50)	<100	470	<100	1,700	< 0.90	<0.88	<0.86	<0.82	< 0.55	<0.87	<0.81	7.9
ETHYLBENZENE	580,000	<100	<100	<100	<100	<100	<100	120	<100	<100	<0.51	<0.49	<0.48	<0.46	<0.31	1.7	<0.46	8.2
HEXANE			NR	NR	NR	NR	NR	NR	NR	NR	17	<2.8	<2.8	<2.7	27	<2.8	59	<2.7
TETRACHLOROETHENE	1,400	16,000	15,000	38,000	11,000	860,000 (50)	25,000	5,700	4,300	4,100	<2.1	64	10	8.7	<1.3	<2.0	<1.9	15
TOLUENE	180,000	<200	320	220	210	<200	250	550	270	270	31	25	16	14	48	3.3	68	30
TRANS-1,2-DICHLOROETHENE	41,000	<100	<100	<100	<100	140,000 (50)	<100	<100	<100	<100	<0.90	<0.88	<0.86	<0.82	<0.55	<0.87	<0.81	8.2
TRICHLOROETHENE	4,100	150	480	18,000	1,200	4,600,000 (50)	1,300	3,000	310	3,100	<0.86	<0.83	<0.82	<0.78	< 0.53	<0.83	<0.77	22
VINYL CHLORIDE	100	<100	<100	<100	<100	1,800	<100	<100	<100	<100	<0.40	< 0.39	< 0.39	< 0.37	<0.25	< 0.39	< 0.37	< 0.32
XYLENE(S)	58,000	<100	120	<100	<100	<100	<100	450	100	130	48	49	31.4	26.3	43.2	8	50	59

Bold Numbers - Exceed ESL Concentrations

-- - Screening level not provided NR - Constituent not reported by laboratory

Table 5. Summa Canister Pressures

Sample Location	Before Sampling (mm Hg)	After Sampling (mm Hg)
SG-10	-20	-7
SG-11	-30	-7
SG-12	-30	-6
SG-13	-30	-7
SG-14	-30	-7
SG-15	-30	-7
SG-16	-30	-7

Table 6. Monitoring Well Depth to Groundwater

	Date	Depth to Water from	Water Elev
	Date	Top of Casing (ft)	(msl)
MW-1	1/28/2008	25.10	384.90
MW-2	1/28/2008	25.23	384.75
MW-3	1/28/2008	25.25	384.23



APPENDIX A

Boring Logs



Well Installation Rickenbacker Assesment Livermore, California 7584.100.101

DATE DRILLED: 12/18/2007 HOLE DEPTH: Approx. 35 ft. HOLE DIAMETER: 8.0 in. LOGGED / REVIEWED BY: J. Preece / SM
DRILLING CONTRACTOR: Gregg Drilling & Testing
DRILLING METHOD: Hollow Stem Auger

	Live 7	rmc 758	ore, California 4.100.101	HOLE DIAMETER: 8.0 in. SURF ELEV (MSL): Approx. 410	ft.					Hollow Stem Auger Direct Push
Depth in Feet	Depth in Meters	Sample Type		DESCRIPTION	Log Symbol	Water Level	Blow Count/Foot	PID (mdd)	Unconfined Strength (tsf) *field approx	Well Construction
			Asphalt WELL GRADED SAND V moist, well graded	VITH GRAVEL (SW), gray, very dense,				17		Flush-Mounted Well Cover
_			SILTY CLAY (CL), dark b	rown, stiff, moist, slightly porous				10*		
_	- - -		SILTY CLAY (CH), dark g	gray, stiff, moist						
_	1 - -			_						
5 —	- - -		Grades to dark olive brow	n.						
-	-			•				0.4		
· · · · · -	_ 2 _									
_	- - -		SILTY GRAVEL (GM). lia	ht olive brown to light olive gray, loose,						
_	- - -		moist to dry, fine- to coars	se-grained sand, and fine gravel						
10 —	3							0.7		
***								0.7		
_	- - -									Bentonite Seal
-	- - 4 -									
-	-									0.01-inch slot width. No. 2/16 sand filter pack.
15 —	<u>-</u>							0.4		
-	_ _ 5							0.4		
_	<u> </u> -									
_	- - -				000					
_	- - - 6									
20 —		\prod								·. ·



Well Installation Rickenbacker Assesment Livermore, California 7584.100.101

LOG - ENVIRONMENTAL + WELL MONITOR WELLS.GPJ ENGEO INC.GDT 2/25/08

DATE DRILLED: 12/18/2007 HOLE DEPTH: Approx. 35 ft. HOLE DIAMETER: 8,0 in. LOGGED / REVIEWED BY: J. Preece / SM
DRILLING CONTRACTOR: Gregg Drilling & Testing
DRILLING METHOD: Hollow Stem Auger

7584.100.101 SURF ELEV (MSL): Approx. 410 ft. HAMMER TYPE: Direct Push Unconfined Strength (tsf) *field approx Blow Count/Foot Depth in Feet **DESCRIPTION** Well Construction og Symbol Water Level PID (ppm) SILTY CLAY (ML), light orange mottled with light gray, stiff, moist to wet, trace fine-grained sand SILTY SAND (SM), light orange mottled with light gray, stiff, moist to wet, trace fine-grained sand A SILTY CLAY (ML), light orange mottled with light gray, stiff, saturated SILTY SAND (SM), light orange mottled with light gray, stiff, saturated, trace fine-grained sand Bottom Plug Bottom of boring at approximately 35 feet.



Well Installation
Rickenbacker Assesment
Livermore, California
7584,100,101

DATE DRILLED: 12/18/2007 HOLE DEPTH: Approx. 35 ft. HOLE DIAMETER: 8.0 in. LOGGED / REVIEWED BY: R. Gandolfo / SM
DRILLING CONTRACTOR: Gregg Drilling & Testing
DRILLING METHOD: Hollow Stem Auger

Depth in Feet	Depth in Meters	Sample Type		DESCRIPTION	Log Symbol	Water Level	Blow Count/Foot	PID (ppm)	Unconfined Strength (tsf) *field approx	Well Construction
5 —	1 - 1		moist, well graded SILTY CLAY (CL), dark o stiff, moist, trace fine-grai	VITH GRAVEL (SW), gray, very dense, rangeish brown mottled with grayish broned sand	Vinio	N	ш	0.1	()	Flush-Mounted Well Cover
10 -	3		Grades to light brown with	orange and black mottling.				0.0		Bentonite Seal
15 —	5		SILTY GRAVEL (GM), ligi medium dense, moist, with subrounded fine gravel	nt olive brown and brownish gray, loose n fine- to medium-grained sand, and				0.0		0.01-inch slot width. No. 2/16 sand filter pack.



Well Installation Rickenbacker Assesment Livermore, California 7584.100.101

DATE DRILLED: 12/18/2007 HOLE DEPTH: Approx. 35 ft. HOLE DIAMETER: 8.0 in. LOGGED / REVIEWED BY: R. Gandolfo / SM
DRILLING CONTRACTOR: Gregg Drilling & Testing
DRILLING METHOD: Hollow Stem Auger

HAMMER TYPE: Direct Push 7584.100.101 SURF ELEV (MSL): Approx. 410 ft. Unconfined Strength (tsf) *field approx Blow Count/Foot Depth in Feet **DESCRIPTION** Well Construction Log Symbol Water Level PID (ppm) 0.1 SILTY CLAY (CL), brown mottled with orangeish gray, stiff, moist to saturated, trace fine-grained sand Grades to some fine grained sand and fine grained, subrounded gravel. 0.0 SILTY GRAVEL (GM), grayish brown, loose to medium dense, saturated, with fine- to medium-grained sand, and subrounded fine 0.0 LOG - ENVIRONMENTAL + WELL MONITOR WELLS.GPJ ENGEO INC.GDT 2/25/08 Bottom Plug 0.0 Bottom of boring at approximately 35 feet.



Well Installation Rickenbacker Assesment Livermore, California 7584.100.101

DATE DRILLED: 12/19/2007 HOLE DEPTH: Approx. 35 ft. HOLE DIAMETER: 8.0 in.

DRILLING CONTRACTOR: Gregg Drilling & Testing DRILLING METHOD: Hollow Stem Auger

LOGGED / REVIEWED BY: R. Gandolfo / SM

SURF ELEV (MSL): Approx. 410 ft. HAMMER TYPE: Direct Push

	1004	4.100.101	SURF ELEV (MSL): Approx. 410	J IL.		11/1	VIIVILIX	111 6.	Direct Push
Depth in Feet Depth in Meters	Sample Type		DESCRIPTION	Log Symbol	Water Level	Blow Count/Foot	PID (mdd)	Unconfined Strength (tsf) *field approx	
5 — 1		moist, well graded SILTY CLAY (CL), dark b	VITH GRAVEL (SW), gray, very dense, rown, stiff, moist, trace fine-grained sand				0.0		Flush-Mounted Well Cover Flush-Mounted Well Cover Flush-Mounted Flush-M
15 — 5		SILTY GRAVEL (GP), oliv moist to saturated, fine- to	re brown to grayish brown, medium dens o coarse-grained sand, and fine gravel				0.0		0.01-inch slot width. No. 2/16 sand filter pack.



Well Installation Rickenbacker Assesment Livermore, California 7584.100.101

LOG - ENVIRONMENTAL + WELL MONITOR WELLS.GPJ ENGEO INC.GDT 2/25/08

DATE DRILLED: 12/19/2007 HOLE DEPTH: Approx. 35 ft. HOLE DIAMETER: 8.0 in. LOGGED / REVIEWED BY: R. Gandolfo / SM
DRILLING CONTRACTOR: Gregg Drilling & Testing
DRILLING METHOD: Hollow Stem Auger

HAMMER TYPE: Direct Push 7584.100.101 SURF ELEV (MSL): Approx. 410 ft. Unconfined Strength (tsf) *field approx Blow Count/Foot Depth in Feet Sample Type **DESCRIPTION** Cog Symbol Well Construction Water Level PID (ppm) SILTY GRAVEL (GP), olive brown to grayish brown, medium dense, moist to saturated, fine- to coarse-grained sand, and fine gravel 0.0 0.0 SILTY CLAY (CL), dark olive brown mottled with orangeish brown, stiff, saturated, some fine-grained sand, and fine gravel 0.0 SILTY GRAVEL (GW), olive brown to grayish brown, medium dense, saturated, fine- to coarse-grained sand, and fine gravel Bottom plug 0.0 Bottom of boring at approximately 35 feet.



Environmental Assessment Rickenbacker Assement Livermore, California 07584.100.101

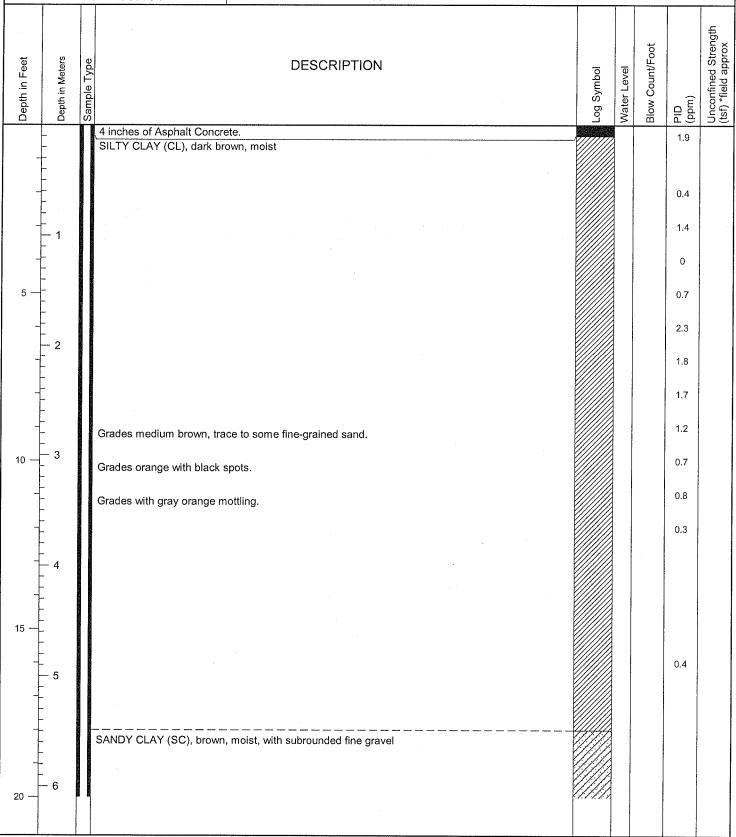
LOG - ENVIRONMENTAL 07584100101 - RNB ADDITIONAL STUDIES.GPJ ENGEO INC.GDT 12/7/07

DATE DRILLED: 11/27/2007 HOLE DEPTH: Approx. 97½ ft.

HOLE DIAMETER: 4.0 in. SURF ELEV (FT-MSL): Approx. 410 ft.

LOGGED / REVIEWED BY: R. Gandolfo / JJT

DRILLING CONTRACTOR: RSI
DRILLING METHOD: Rotosonic
HAMMER TYPE: N/A





Environmental Assessment Rickenbacker Assement Livermore, California 07584.100.101

LOG - ENVIRONMENTAL 07584100101 - RNB ADDITIONAL STUDIES.GPJ ENGEO INC.GDT 12/7/07

DATE DRILLED: 11/27/2007 HOLE DEPTH: Approx. 97½ ft. HOLE DIAMETER: 4.0 in.

SURF ELEV (FT-MSL): Approx. 410 ft.

LOGGED / REVIEWED BY: R. Gandolfo / JJT

DRILLING CONTRACTOR: RSI
DRILLING METHOD: Rotosonic
HAMMER TYPE: N/A

07	07584.100.101 SURF ELEV (FT-MSL): Approx. 410 ft. HAMMER TYPE:							1	
Depth in Feet Depth in Meters	Sample Type		DESCRIPTION		Log Symbol	Water Level	Blow Count/Foot	(mdd)	Unconfined Strength
-		4 inches of Asphalt Cond	crete.		<i>7/11/11/11</i>		····	1.9	
	Stellariti	SILTY CLAY (CL), dark t	brown, moist						
-	and a second							·	·
								0.4	
-	STATE OF STA							0.4	
F								1.4	
<u> </u>									
+								0	
-	A CONTRACTOR								
5 —	WORDEN OF THE PERSON NAMED IN COLUMN TWO IS NOT THE PERSON NAMED IN COLUMN TWO IS NAMED IN COLUMN TWO IS NOT THE PERSON NAMED IN COLUMN TWO IS NAMED IN COLUMN T							0.7	
-									
	Section of							2.3	
_ 2	(Marketin							1.0	
						1		1.8	
<u> </u>	editores supposed							1.7	
-	2000/0055 2000-00								
+		Grades medium brown to	race to some fine-grained sand.					1.2	
3		oracoo modium provin, a	rado to domo imo granica dana.						
10 —	Par Casago	Grades orange with black	spots.					0.7	
-									
		Grades with gray orange	mottling.					8.0	
								0.3	
	AMETERS							0.3	
-4									
-	STORY TO								
+	Selection.								
_						-			
15 —									
}	S CONTRA							0.4	
<u> </u>								0.4	
+								ĺ	
<u> </u>		CANDY OLAY (CO)							
<u> </u>	Security See	SANDY CLAY (SC), brow	n, moist, with subrounded fine gravel						
+	Service of the servic			·					
_ 6				k					
20 —				ĺ	7.11.22		1		



Environmental Assessment Rickenbacker Assement Livermore, California 07584.100.101

LOG - ENVIRONMENTAL 07584100101 - RNB ADDITIONAL STUDIES.GPJ ENGEO INC.GDT 12/7/07

DATE DRILLED: 11/27/2007 HOLE DEPTH: Approx. 97½ ft. HOLE DIAMETER: 4.0 in.

SURF ELEV (FT-MSL): Approx. 410 ft.

LOGGED / REVIEWED BY: R. Gandolfo / JJT DRILLING CONTRACTOR: RSI

DRILLING METHOD: Rotosonic
HAMMER TYPE: N/A

	07584.100.101 SURF ELEV (FT-MSL): Approx. 410 ft. HAMMER TYPE: N/A								,	
Depth in Feet	Depth in Meters	Sample Type		DESCRIPTION		Log Symbol	Water Level	Blow Count/Foot	(mdd)	Unconfined Strength (tsf) *field approx
25 — 30 — 35 — 40 —	7 10 11 12		Sandy Clay (continued) SILTY CLAY (CL), gray, m	oist, some fine-grained sand sh yellow, moist, trace clay, some subrounded fine grave ne to coarse gravel.					0.3	



Environmental Assessment Rickenbacker Assement Livermore, California

DATE DRILLED: 11/27/2007 HOLE DEPTH: Approx. 971/2 ft. HOLE DIAMETER: 4.0 in.

LOGGED / REVIEWED BY: R. Gandolfo / JJT

DRILLING CONTRACTOR: RSI DRILLING METHOD: Rotosonic

	07	55	34.100.101	SURF ELEV (FT-MSL): Approx. 410 ft.	HAMMER T	YPE: N/	А Т Т		1	
Depth in Feet	Depth in Meters	Sample Type		DESCRIPTION		Log Symbol	Water Level	Blow Count/Foot	PID (mdd)	Unconfined Strength
	13		Silty Sand (continued) SILTY CLAY (CL), gray, t						0.3	
45 — -	- - - - - - - - - - - - - -								0.5	
50 —			SILTY CLAY (CL), gray, r	noist, with fine-grained sand, some subrounded fine	gravel				0.5	
- - 55 — -			SILTY CLAY (CL), light be	rown mottled with black, moist, trace fine-grained sar	nd				To the second se	
-60 —	- - - - - 18		Grades with subrounded f	ine to coarse gravel.					-	



Environmental Assessment Rickenbacker Assement Livermore, California 07584.100.101

LOG - ENVIRONMENTAL 07584100101 - RNB ADDITIONAL STUDIES.GPJ ENGEO INC.GDT 12/7/07

DATE DRILLED: 11/27/2007 HOLE DEPTH: Approx. 97½ ft.

HOLE DIAMETER: 4.0 in. SURF ELEV (FT-MSL): Approx. 410 ft.

LOGGED / REVIEWED BY: R. Gandolfo / JJT

DRILLING CONTRACTOR: RSI
DRILLING METHOD: Rotosonic
HAMMER TYPE: N/A

07584.100	.101 SURF ELEV (FT-MSL): Approx. 410 ft.	HAMMER TYPE: N/A	
Depth in Feet Depth in Meters Sample Type	DESCRIPTION	Log Symbol Water Level	Blow Count/Foot PID (ppm) Unconfined Strength (tsf) *field approx
GRAVE subrour - 19 - 19 - 20 - 21 - 21 - 22 - 3ILTY C Grades - 23 - 38 - 38 - 48 - 58 - 58 - 68 - 75 - 23 - 68 - 68 - 75 - 23 - 75 - 75 - 75 - 75 - 75 - 75 - 75 - 75	trace to some clay. SP), brown, wet, medium-grained sand, clay, trace subrounded to rounded fine to trace gravel and some sand. L (GM), grayish brown, wet, subrounded fine gravel, with fine- to coarse-grained sand. L (GM), grayish brown, wet, subrounded fine gravel, with fine- to coarse-grained sand.	ed sand, with	a a a b a b a b a b a



Environmental Assessment Rickenbacker Assement Livermore, California 07584.100.101

LOG - ENVIRONMENTAL 07584100101 - RNB ADDITIONAL STUDIES.GPJ ENGEO INC.GDT 12/7/07

DATE DRILLED: 11/27/2007 HOLE DEPTH: Approx. 97½ ft. HOLE DIAMETER: 4.0 in.

SURF ELEV (FT-MSL): Approx. 410 ft.

LOGGED / REVIEWED BY: R. Gandolfo / JJT DRILLING CONTRACTOR; RSI

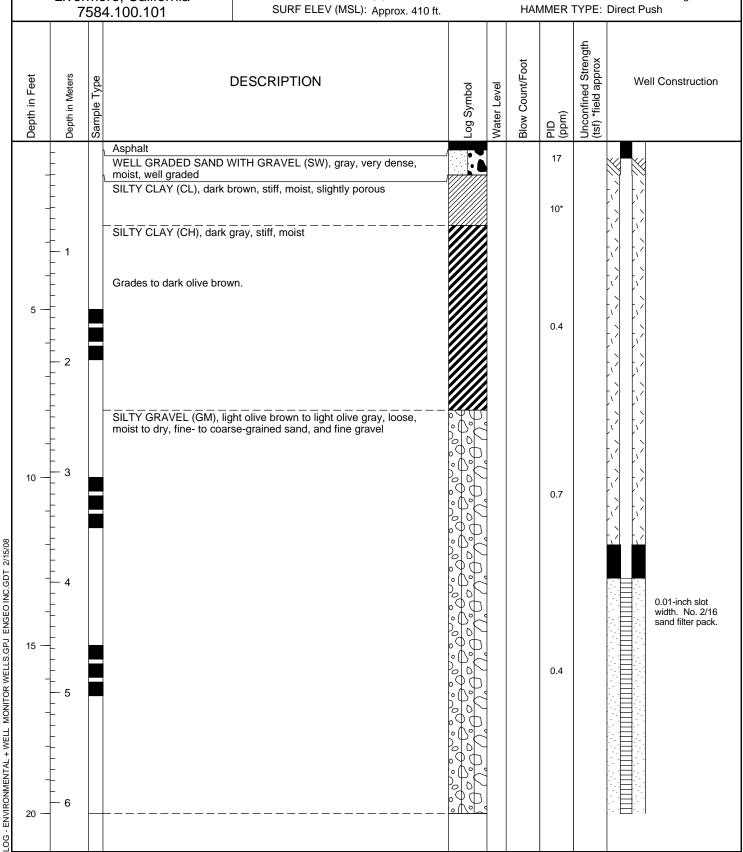
RILLING CONTRACTOR: RSI
DRILLING METHOD: Rotosonic
HAMMER TYPE: N/A

	07	758	34.100.101	SURF ELEV (FT-MSL): Approx. 410 ft. H	AMMER T	YPE: N/	A —		r	
Depth in Feet	Depth in Meters	Sample Type		DESCRIPTION		Log Symbol	Water Level	Blow Count/Foot	Old (mdd)	Unconfined Strength (tsf) *field approx
85 —	25			, subrounded coarse gravel, some silt, some clay					0	
90 —	28		Grades fine to coarse subro	unded gravel.					0.2	
			Bottom of boring at approxir	nately 97 1/2 feet. Ground water measured at approximately 4	42 feet.					



Well Installation Rickenbacker Assesment Livermore, California 7584.100.101

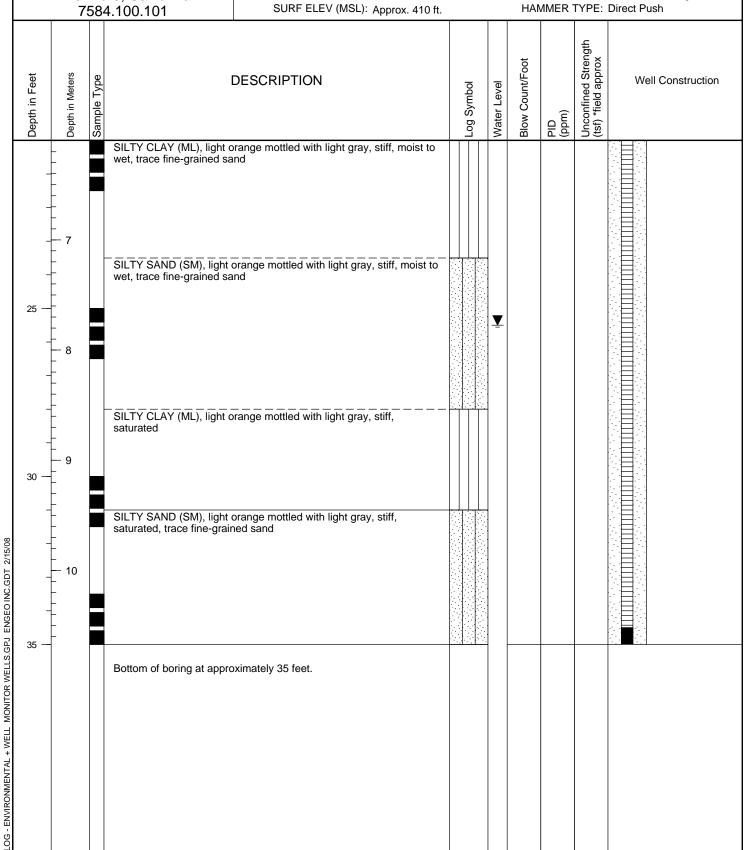
DATE DRILLED: 12/18/2007 HOLE DEPTH: Approx. 35 ft. HOLE DIAMETER: 8.0 in. LOGGED / REVIEWED BY: J. Preece / SM
DRILLING CONTRACTOR: Gregg Drilling & Testing
DRILLING METHOD: Hollow Stem Auger
HAMMER TYPE: Direct Push





Well Installation Rickenbacker Assesment Livermore, California 7584.100.101

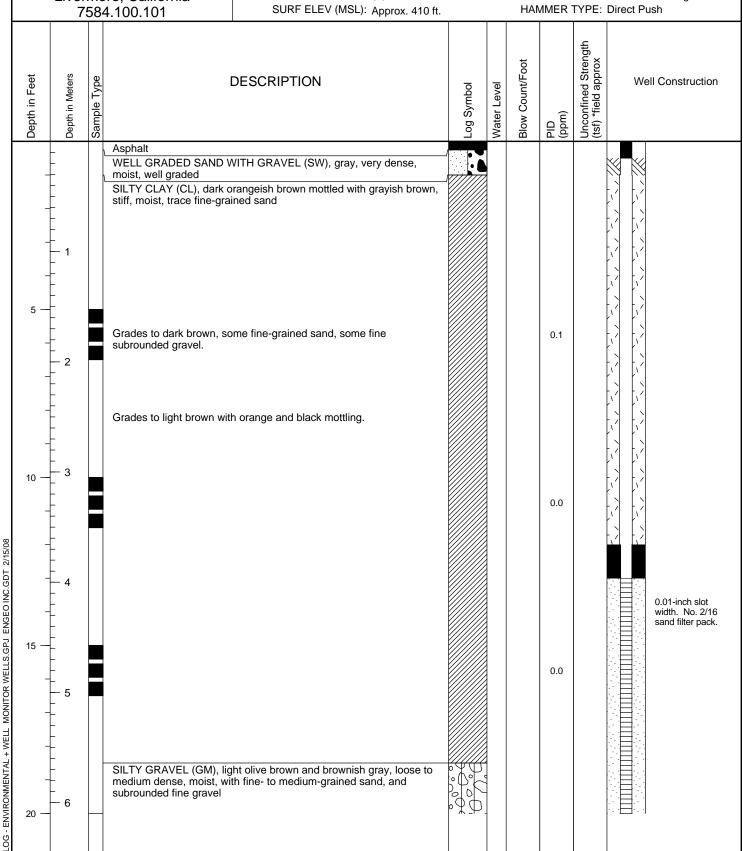
DATE DRILLED: 12/18/2007 HOLE DEPTH: Approx. 35 ft. HOLE DIAMETER: 8.0 in. LOGGED / REVIEWED BY: J. Preece / SM
DRILLING CONTRACTOR: Gregg Drilling & Testing
DRILLING METHOD: Hollow Stem Auger
HAMMER TYPE: Direct Push





Well Installation Rickenbacker Assesment Livermore, California 7584.100.101

DATE DRILLED: 12/18/2007 HOLE DEPTH: Approx. 35 ft. HOLE DIAMETER: 8.0 in. LOGGED / REVIEWED BY: R. Gandolfo / SM
DRILLING CONTRACTOR: Gregg Drilling & Testing
DRILLING METHOD: Hollow Stem Auger





Well Installation Rickenbacker Assesment Livermore, California 7584.100.101

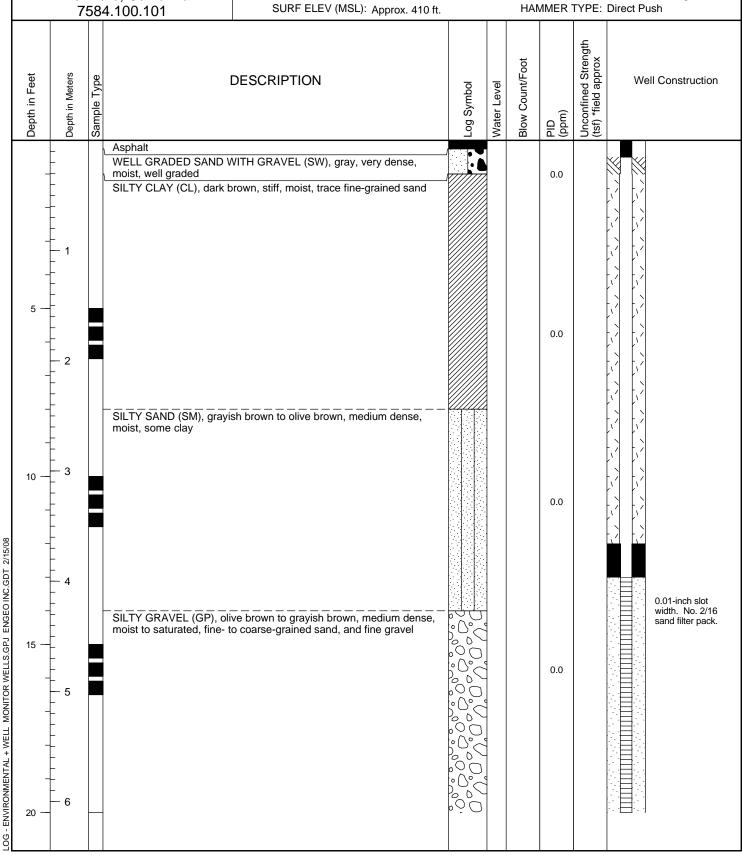
DATE DRILLED: 12/18/2007 HOLE DEPTH: Approx. 35 ft. HOLE DIAMETER: 8.0 in. SURF ELEV (MSL): Approx. 410 ft. LOGGED / REVIEWED BY: R. Gandolfo / SM
DRILLING CONTRACTOR: Gregg Drilling & Testing
DRILLING METHOD: Hollow Stem Auger
HAMMER TYPE: Direct Push

		7	584	1.100.101	SURF ELEV (MSL): Approx. 410 ft			ПАІ	VIIVIER	ITPE.	Direct Push
	Depth in Feet	Depth in Meters Sample Type		ן	DESCRIPTION	Log Symbol	Water Level	Blow Count/Foot	PID (mdd)	Unconfined Strength (tsf) *field approx	Well Construction
	25 —	- - - - - - - - - - - - - - - - - - -			mottled with orangeish gray, stiff, moist to ed sand		<u></u>		0.1		
3EO INC.GDT 2/15/08	30 —	- - - - - - - - - - - - - - - - - - -		SILTY GRAVEL (GM), grayish brown, loose to medium dense, saturated, with fine- to medium-grained sand, and subrounded fir gravel		10 (10 (10 (10 M)))		0.0			
LOG - ENVIRONMENTAL + WELL MONITOR WELLS.GPJ ENGEO INC.GDT 2/15/08	35 —			Bottom of boring at appro	ximately 35 feet.				0.0		



Well Installation Rickenbacker Assesment Livermore, California 7584 100 101

DATE DRILLED: 12/19/2007 HOLE DEPTH: Approx. 35 ft. HOLE DIAMETER: 8.0 in. LOGGED / REVIEWED BY: R. Gandolfo / SM
DRILLING CONTRACTOR: Gregg Drilling & Testing
DRILLING METHOD: Hollow Stem Auger





Well Installation Rickenbacker Assesment Livermore, California 7584.100.101

DATE DRILLED: 12/19/2007 HOLE DEPTH: Approx. 35 ft. HOLE DIAMETER: 8.0 in. SURF ELEV (MSL): Approx. 410 ft. LOGGED / REVIEWED BY: R. Gandolfo / SM
DRILLING CONTRACTOR: Gregg Drilling & Testing
DRILLING METHOD: Hollow Stem Auger
HAMMER TYPE: Direct Push

	7584.100.101 SURF ELEV (MSL): Approx. 410 ft.		t.			HAMMER		Direct Push		
Depth in Feet	Depth in Meters	Sample Type		DESCRIPTION	C C Log Symbol	Water Level	Blow Count/Foot	PID (mdd)	Unconfined Strength (tsf) *field approx	Well Construction
25 -			SILTY CLAY (CL), dark o stiff, saturated, some fine	re brown to grayish brown, medium dense, o coarse-grained sand, and fine gravel rive brown mottled with orangeish brown, and fine gravel regrained sand, and fine gravel ve brown to grayish brown, medium dense grained sand, and fine gravel				0.0		
35 —										



APPENDIX B

TESTAMERICA LABORATORIES, INC.

Laboratory Analytical Reports



ANALYTICAL REPORT

Job Number: 720-11927-1

Job Description: Rickenbacker

For: Engeo, Inc. 580 N Wilma Avenue Suite A Ripon, CA 95366-9502

Attention: Mr. Richard Gandolfo



Melissa Brewer
Project Manager I
melissa.brewer@testamericainc.com
12/04/2007

Job Narrative 720-J11927-1

Comments

No additional comments.

Receipt

All samples were received in good condition within temperature requirements.

GC/MS VOA

No analytical or quality issues were noted.

GC VOA

No analytical or quality issues were noted.

GC Semi VOA

No analytical or quality issues were noted.

Organic Prep
No analytical or quality issues were noted.

EXECUTIVE SUMMARY - Detections

Client: Engeo, Inc. Job Number: 720-11927-1

Lab Sample ID Analyte	Client Sample ID	Result / Qualifier	Reporting Limit	Units	Method	
720-11927-1	S @ 10'					
Tetrachloroethene		79	4.9	ug/Kg	8260B	
720-11927-2	S @ 20'					
Tetrachloroethene		17	5.0	ug/Kg	8260B	
Silica Gel Cleanup Diesel Range Orga		17	0.99	mg/Kg	8015B	
720-11927-5	S @ 50'					
Tetrachloroethene		14	4.9	ug/Kg	8260B	
Silica Gel Cleanup Diesel Range Orga		1.1	0.99	mg/Kg	8015B	

METHOD SUMMARY

Client: Engeo, Inc. Job Number: 720-11927-1

Description	Lab Location	Method	Preparation Method
Matrix: Solid			
Volatile Organic Compounds by GC/MS	TAL SF	SW846 8260B	
Volatile Organic Compounds by GC/MS (Low Level)	TAL SF	SW846 8260B	
Purge and Trap for Solids	TAL SF		SW846 5030B
Purge and Trap for Solids	TAL SF		SW846 5030B
Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)	TAL SF	SW846 8015B	
Ultrasonic Extraction	TAL SF		SW846 3550B

Lab References:

TAL SF = TestAmerica San Francisco

Method References:

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

SAMPLE SUMMARY

Client: Engeo, Inc. Job Number: 720-11927-1

Lab Sample ID	Client Sample ID	Client Matrix	Date/Time Sampled	Date/Time Received
720-11927-1	S @ 10'	Solid	11/27/2007 0925	11/27/2007 1655
720-11927-2	S @ 20'	Solid	11/27/2007 0930	11/27/2007 1655
720-11927-3	S @ 30'	Solid	11/27/2007 0940	11/27/2007 1655
720-11927-4	S @ 40'	Solid	11/27/2007 1010	11/27/2007 1655
720-11927-5	S @ 50'	Solid	11/27/2007 1300	11/27/2007 1655
720-11927-6	S @ 60'	Solid	11/27/2007 1300	11/27/2007 1655
720-11927-7	S @ 70'	Solid	11/27/2007 1440	11/27/2007 1655
720-11927-8	S @ 80'	Solid	11/27/2007 1515	11/27/2007 1655
720-11927-9	S @ 90'	Solid	11/27/2007 1620	11/27/2007 1655

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 10'

 Lab Sample ID:
 720-11927-1
 Date Sampled:
 11/27/2007 0925

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807006.DDilution:1.0Initial Weight/Volume:5.07 g

Date Analyzed: 11/28/2007 1526 Final Weight/Volume: 5.07 g

Date Prepared: 11/28/2007 1330

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
Methyl tert-butyl ether	ND		4.9
Acetone	ND		49
Benzene	ND		4.9
Dichlorobromomethane	ND		4.9
Bromobenzene	ND		4.9
Chlorobromomethane	ND		20
Bromoform	ND		4.9
Bromomethane	ND		9.9
2-Butanone (MEK)	ND		49
n-Butylbenzene	ND		4.9
sec-Butylbenzene	ND		4.9
tert-Butylbenzene	ND		4.9
Carbon disulfide	ND		4.9
Carbon tetrachloride	ND		4.9
Chlorobenzene	ND		4.9
Chloroethane	ND		9.9
Chloroform	ND		4.9
Chloromethane	ND		9.9
2-Chlorotoluene	ND		4.9
4-Chlorotoluene	ND		4.9
Chlorodibromomethane	ND		4.9
1,2-Dichlorobenzene	ND		4.9
1,3-Dichlorobenzene	ND		4.9
1,4-Dichlorobenzene	ND		4.9
1,3-Dichloropropane	ND		4.9
1,1-Dichloropropene	ND		4.9
1,2-Dibromo-3-Chloropropane	ND		49
Ethylene Dibromide	ND		4.9
Dibromomethane	ND		9.9
Dichlorodifluoromethane	ND		9.9
1,1-Dichloroethane	ND		4.9
1,2-Dichloroethane	ND		4.9
1,1-Dichloroethene	ND		4.9
cis-1,2-Dichloroethene	ND		4.9
trans-1,2-Dichloroethene	ND		4.9
1,2-Dichloropropane	ND		4.9
cis-1,3-Dichloropropene	ND		4.9
trans-1,3-Dichloropropene	ND		4.9
Ethylbenzene	ND		4.9
Hexachlorobutadiene	ND		4.9
2-Hexanone	ND		49
Isopropylbenzene	ND		4.9
4-Isopropyltoluene	ND		4.9
Methylene Chloride	ND		9.9

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 10'

 Lab Sample ID:
 720-11927-1
 Date Sampled:
 11/27/2007 0925

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807006.DDilution:1.0Initial Weight/Volume:5.07 g

Date Analyzed: 11/28/2007 1526 Final Weight/Volume: 10 mL

Date Prepared: 11/28/2007 1330

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		49
Naphthalene	ND		9.9
N-Propylbenzene	ND		4.9
Styrene	ND		4.9
1,1,1,2-Tetrachloroethane	ND		4.9
1,1,2,2-Tetrachloroethane	ND		4.9
Tetrachloroethene	79		4.9
Toluene	ND		4.9
1,2,3-Trichlorobenzene	ND		4.9
1,2,4-Trichlorobenzene	ND		4.9
1,1,1-Trichloroethane	ND		4.9
1,1,2-Trichloroethane	ND		4.9
Trichloroethene	ND		4.9
Trichlorofluoromethane	ND		4.9
1,2,3-Trichloropropane	ND		4.9
1,1,2-Trichloro-1,2,2-trifluoroethan	ie ND		4.9
1,2,4-Trimethylbenzene	ND		4.9
1,3,5-Trimethylbenzene	ND		4.9
Vinyl acetate	ND		49
Vinyl chloride	ND		4.9
Xylenes, Total	ND		9.9
2,2-Dichloropropane	ND		4.9
Surrogate	%Rec	<i>F</i>	Acceptance Limits
4-Bromofluorobenzene	87		50 - 138
1,2-Dichloroethane-d4 (Surr)	98		66 - 127
Toluene-d8 (Surr)	89		51 - 129

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 10'

 Lab Sample ID:
 720-11927-1
 Date Sampled:
 11/27/2007 0925

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 720-29124 Instrument ID: Varian 3900A

Preparation: 5030B Prep Batch: 720-29065 Lab File ID: c:\saturnws\data\200711\11

Dilution: 1.0 Initial Weight/Volume: 5.48 g

Date Analyzed: 11/29/2007 1057 Final Weight/Volume: 10 mL Date Prepared: 11/29/2007 0939

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL
Gasoline Range Organics (GRO)-C5-C12 ND 0.23

 Surrogate
 %Rec
 Acceptance Limits

 1,2-Dichloroethane-d4 (Surr)
 93
 60 - 140

 Toluene-d8 (Surr)
 99
 70 - 130

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 20'

 Lab Sample ID:
 720-11927-2
 Date Sampled:
 11/27/2007 0930

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807009.DDilution:1.0Initial Weight/Volume:5.04 g

Date Analyzed: 11/28/2007 1641 Final Weight/Volume: 5.04 g

Date Prepared: 11/28/2007 1330

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
Methyl tert-butyl ether	ND		5.0
Acetone	ND		50
Benzene	ND		5.0
Dichlorobromomethane	ND		5.0
Bromobenzene	ND		5.0
Chlorobromomethane	ND		20
Bromoform	ND		5.0
Bromomethane	ND		9.9
2-Butanone (MEK)	ND		50
n-Butylbenzene	ND		5.0
sec-Butylbenzene	ND		5.0
tert-Butylbenzene	ND		5.0
Carbon disulfide	ND		5.0
Carbon tetrachloride	ND		5.0
Chlorobenzene	ND		5.0
Chloroethane	ND		9.9
Chloroform	ND		5.0
Chloromethane	ND		9.9
2-Chlorotoluene	ND		5.0
4-Chlorotoluene	ND		5.0
Chlorodibromomethane	ND		5.0
1,2-Dichlorobenzene	ND		5.0
1,3-Dichlorobenzene	ND		5.0
1,4-Dichlorobenzene	ND		5.0
1,3-Dichloropropane	ND		5.0
1,1-Dichloropropene	ND		5.0
1,2-Dibromo-3-Chloropropane	ND		50
Ethylene Dibromide	ND		5.0
Dibromomethane	ND		9.9
Dichlorodifluoromethane	ND		9.9
1,1-Dichloroethane	ND		5.0
1,2-Dichloroethane	ND		5.0
1,1-Dichloroethene	ND		5.0
cis-1,2-Dichloroethene	ND		5.0
trans-1,2-Dichloroethene	ND		5.0
1,2-Dichloropropane	ND		5.0
cis-1,3-Dichloropropene	ND		5.0
trans-1,3-Dichloropropene	ND		5.0
Ethylbenzene	ND		5.0
Hexachlorobutadiene	ND		5.0
2-Hexanone	ND		50
Isopropylbenzene	ND		5.0
4-Isopropyltoluene	ND		5.0
Methylene Chloride	ND		9.9
Montylette Official	IAD		5.5

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 20'

 Lab Sample ID:
 720-11927-2
 Date Sampled:
 11/27/2007 0930

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807009.DDilution:1.0Initial Weight/Volume:5.04 g

Date Analyzed: 11/28/2007 1641 Final Weight/Volume: 10 mL

Date Prepared: 11/28/2007 1330

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		50
Naphthalene	ND		9.9
N-Propylbenzene	ND		5.0
Styrene	ND		5.0
1,1,1,2-Tetrachloroethane	ND		5.0
1,1,2,2-Tetrachloroethane	ND		5.0
Tetrachloroethene	17		5.0
Toluene	ND		5.0
1,2,3-Trichlorobenzene	ND		5.0
1,2,4-Trichlorobenzene	ND		5.0
1,1,1-Trichloroethane	ND		5.0
1,1,2-Trichloroethane	ND		5.0
Trichloroethene	ND		5.0
Trichlorofluoromethane	ND		5.0
1,2,3-Trichloropropane	ND		5.0
1,1,2-Trichloro-1,2,2-trifluoroethar	ne ND		5.0
1,2,4-Trimethylbenzene	ND		5.0
1,3,5-Trimethylbenzene	ND		5.0
Vinyl acetate	ND		50
Vinyl chloride	ND		5.0
Xylenes, Total	ND		9.9
2,2-Dichloropropane	ND		5.0
Surrogate	%Rec	Acceptanc	e Limits
4-Bromofluorobenzene	94	50 - 138	
1,2-Dichloroethane-d4 (Surr)	100	66 - 127	
Toluene-d8 (Surr)	95	51 - 129	

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 20'

 Lab Sample ID:
 720-11927-2
 Date Sampled:
 11/27/2007 0930

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 720-29124 Instrument ID: Varian 3900A

Preparation: 5030B Prep Batch: 720-29065 Lab File ID: c:\saturnws\data\200711\11

Dilution: 1.0 Initial Weight/Volume: 5.30 g

Date Analyzed: 11/29/2007 1227 Final Weight/Volume: 10 mL Date Prepared: 11/29/2007 0939

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL

Gasoline Range Organics (GRO)-C5-C12 ND 0.24

 Surrogate
 %Rec
 Acceptance Limits

 1,2-Dichloroethane-d4 (Surr)
 100
 60 - 140

 Toluene-d8 (Surr)
 95
 70 - 130

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 30'

 Lab Sample ID:
 720-11927-3
 Date Sampled:
 11/27/2007 0940

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807010.DDilution:1.0Initial Weight/Volume:5.11 g

Date Analyzed: 11/28/2007 1706 Final Weight/Volume: 10 mL

Date Prepared: 11/28/2007 1330

Analyte	, , , , , , , , , , , , , , , , , , , ,	Qualifier RL
Methyl tert-butyl ether	ND	4.9
Acetone	ND	49
Benzene	ND	4.9
Dichlorobromomethane	ND	4.9
Bromobenzene	ND	4.9
Chlorobromomethane	ND	20
Bromoform	ND	4.9
Bromomethane	ND	9.8
2-Butanone (MEK)	ND	49
n-Butylbenzene	ND	4.9
sec-Butylbenzene	ND	4.9
tert-Butylbenzene	ND	4.9
Carbon disulfide	ND	4.9
Carbon tetrachloride	ND	4.9
Chlorobenzene	ND	4.9
Chloroethane	ND	9.8
Chloroform	ND	4.9
Chloromethane	ND	9.8
2-Chlorotoluene	ND	4.9
4-Chlorotoluene	ND	4.9
Chlorodibromomethane	ND	4.9
1,2-Dichlorobenzene	ND	4.9
1,3-Dichlorobenzene	ND	4.9
1,4-Dichlorobenzene	ND	4.9
1,3-Dichloropropane	ND	4.9
1,1-Dichloropropene	ND	4.9
1,2-Dibromo-3-Chloropropane	ND	49
Ethylene Dibromide	ND	4.9
Dibromomethane	ND	9.8
Dichlorodifluoromethane	ND	9.8
1,1-Dichloroethane	ND	4.9
1,2-Dichloroethane	ND	4.9
1,1-Dichloroethene	ND	4.9
cis-1,2-Dichloroethene	ND	4.9
trans-1,2-Dichloroethene	ND	4.9
1,2-Dichloropropane	ND	4.9
cis-1,3-Dichloropropene	ND	4.9
trans-1,3-Dichloropropene	ND	4.9
Ethylbenzene	ND	4.9
Hexachlorobutadiene	ND ND	4.9
2-Hexanone	ND	49
Isopropylbenzene	ND	4.9
4-Isopropyltoluene	ND ND	4.9
Methylene Chloride	ND	9.8
wearyiene ornonae	IND	3.0

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 30'

 Lab Sample ID:
 720-11927-3
 Date Sampled:
 11/27/2007 0940

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807010.DDilution:1.0Initial Weight/Volume:5.11 g

Date Analyzed: 11/28/2007 1706 Final Weight/Volume: 10 mL

Date Prepared: 11/28/2007 1330

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		49
Naphthalene	ND		9.8
N-Propylbenzene	ND		4.9
Styrene	ND		4.9
1,1,1,2-Tetrachloroethane	ND		4.9
1,1,2,2-Tetrachloroethane	ND		4.9
Tetrachloroethene	ND		4.9
Toluene	ND		4.9
1,2,3-Trichlorobenzene	ND		4.9
1,2,4-Trichlorobenzene	ND		4.9
1,1,1-Trichloroethane	ND		4.9
1,1,2-Trichloroethane	ND		4.9
Trichloroethene	ND		4.9
Trichlorofluoromethane	ND		4.9
1,2,3-Trichloropropane	ND		4.9
1,1,2-Trichloro-1,2,2-trifluoroethar	ne ND		4.9
1,2,4-Trimethylbenzene	ND		4.9
1,3,5-Trimethylbenzene	ND		4.9
Vinyl acetate	ND		49
Vinyl chloride	ND		4.9
Xylenes, Total	ND		9.8
2,2-Dichloropropane	ND		4.9
Surrogate	%Rec		Acceptance Limits
4-Bromofluorobenzene	94		50 - 138
1,2-Dichloroethane-d4 (Surr)	95		66 - 127
Toluene-d8 (Surr)	94		51 - 129

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 30'

 Lab Sample ID:
 720-11927-3
 Date Sampled:
 11/27/2007 0940

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 720-29124 Instrument ID: Varian 3900A

Preparation: 5030B Prep Batch: 720-29065 Lab File ID: c:\saturnws\data\200711\11

Dilution: 1.0 Initial Weight/Volume: 5.16 g

Date Analyzed: 11/29/2007 1249 Final Weight/Volume: 10 mL Date Prepared: 11/29/2007 0939

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL
Gasoline Range Organics (GRO)-C5-C12 ND 0.24

 Surrogate
 %Rec
 Acceptance Limits

 1,2-Dichloroethane-d4 (Surr)
 96
 60 - 140

 Toluene-d8 (Surr)
 96
 70 - 130

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 40'

 Lab Sample ID:
 720-11927-4
 Date Sampled:
 11/27/2007 1010

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807011.DDilution:1.0Initial Weight/Volume:5.03 g

Date Analyzed: 11/28/2007 1731 Final Weight/Volume: 5.03 g

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
Methyl tert-butyl ether	ND		5.0
Acetone	ND		50
Benzene	ND		5.0
Dichlorobromomethane	ND		5.0
Bromobenzene	ND		5.0
Chlorobromomethane	ND		20
Bromoform	ND		5.0
Bromomethane	ND		9.9
2-Butanone (MEK)	ND		50
n-Butylbenzene	ND		5.0
sec-Butylbenzene	ND		5.0
tert-Butylbenzene	ND		5.0
Carbon disulfide	ND		5.0
Carbon tetrachloride	ND		5.0
Chlorobenzene	ND		5.0
Chloroethane	ND		9.9
Chloroform	ND		5.0
Chloromethane	ND		9.9
2-Chlorotoluene	ND		5.0
4-Chlorotoluene	ND		5.0
Chlorodibromomethane	ND		5.0
1,2-Dichlorobenzene	ND		5.0
1,3-Dichlorobenzene	ND		5.0
1,4-Dichlorobenzene	ND		5.0
1,3-Dichloropropane	ND		5.0
1,1-Dichloropropene	ND		5.0
1,2-Dibromo-3-Chloropropane	ND		50
Ethylene Dibromide	ND		5.0
Dibromomethane	ND		9.9
Dichlorodifluoromethane	ND		9.9
1,1-Dichloroethane	ND		5.0
1,2-Dichloroethane	ND		5.0
1,1-Dichloroethene	ND		5.0
cis-1,2-Dichloroethene	ND		5.0
trans-1,2-Dichloroethene	ND		5.0
1,2-Dichloropropane	ND		5.0
cis-1,3-Dichloropropene	ND		5.0
trans-1,3-Dichloropropene	ND		5.0
Ethylbenzene	ND		5.0
Hexachlorobutadiene	ND		5.0
2-Hexanone	ND		50
Isopropylbenzene	ND		5.0
4-Isopropyltoluene	ND		5.0
Methylene Chloride	ND		9.9

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 40'

 Lab Sample ID:
 720-11927-4
 Date Sampled:
 11/27/2007 1010

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807011.DDilution:1.0Initial Weight/Volume:5.03 g

Date Analyzed: 11/28/2007 1731 Final Weight/Volume: 10 mL

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		50
Naphthalene	ND		9.9
N-Propylbenzene	ND		5.0
Styrene	ND		5.0
1,1,1,2-Tetrachloroethane	ND		5.0
1,1,2,2-Tetrachloroethane	ND		5.0
Tetrachloroethene	ND		5.0
Toluene	ND		5.0
1,2,3-Trichlorobenzene	ND		5.0
1,2,4-Trichlorobenzene	ND		5.0
1,1,1-Trichloroethane	ND		5.0
1,1,2-Trichloroethane	ND		5.0
Trichloroethene	ND		5.0
Trichlorofluoromethane	ND		5.0
1,2,3-Trichloropropane	ND		5.0
1,1,2-Trichloro-1,2,2-trifluoroethar	ie ND		5.0
1,2,4-Trimethylbenzene	ND		5.0
1,3,5-Trimethylbenzene	ND		5.0
Vinyl acetate	ND		50
Vinyl chloride	ND		5.0
Xylenes, Total	ND		9.9
2,2-Dichloropropane	ND		5.0
Surrogate	%Rec	Accep	tance Limits
4-Bromofluorobenzene	93	50 -	138
1,2-Dichloroethane-d4 (Surr)	99	66 -	127
Toluene-d8 (Surr)	94	51 -	129

0.25

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 40'

Gasoline Range Organics (GRO)-C5-C12

Lab Sample ID: 720-11927-4 Date Sampled: 11/27/2007 1010 Client Matrix: Solid Date Received: 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS

8260B Analysis Batch: 720-29124 Varian 3900A Method: Instrument ID:

Preparation: 5030B Prep Batch: 720-29065 Lab File ID: c:\saturnws\data\200711\11

Dilution: 1.0 Initial Weight/Volume: 5.08 g

Date Analyzed: 11/29/2007 1312 Final Weight/Volume: 10 mL Date Prepared: 11/29/2007 0939

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RLND

%Rec Surrogate Acceptance Limits

1,2-Dichloroethane-d4 (Surr) 97 60 - 140 Toluene-d8 (Surr) 97 70 - 130

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 50'

 Lab Sample ID:
 720-11927-5
 Date Sampled:
 11/27/2007 1300

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807012.DDilution:1.0Initial Weight/Volume:5.12 g

Dilution: 1.0 Initial Weight/Volume: 5.12 g
Date Analyzed: 11/28/2007 1756 Final Weight/Volume: 10 mL

Methyl tert-butyl ether Acetone Benzene Dichlorobromomethane Bromobenzene	ND ND ND ND	4.9 49
Benzene Dichlorobromomethane	ND	
Dichlorobromomethane		4.0
	ND	4.9
Bromohenzene	ND	4.9
DIGITIODCIZCTIC	ND	4.9
Chlorobromomethane	ND	20
Bromoform	ND	4.9
Bromomethane	ND	9.8
2-Butanone (MEK)	ND	49
n-Butylbenzene	ND	4.9
sec-Butylbenzene	ND	4.9
tert-Butylbenzene	ND	4.9
Carbon disulfide	ND	4.9
Carbon tetrachloride	ND	4.9
Chlorobenzene	ND	4.9
Chloroethane	ND	9.8
Chloroform	ND	4.9
Chloromethane	ND	9.8
2-Chlorotoluene	ND	4.9
4-Chlorotoluene	ND	4.9
Chlorodibromomethane	ND	4.9
1,2-Dichlorobenzene	ND	4.9
1,3-Dichlorobenzene	ND	4.9
1,4-Dichlorobenzene	ND	4.9
1,3-Dichloropropane	ND	4.9
1,1-Dichloropropene	ND	4.9
1,2-Dibromo-3-Chloropropane	ND	49
Ethylene Dibromide	ND	4.9
Dibromomethane	ND	9.8
Dichlorodifluoromethane	ND	9.8
1,1-Dichloroethane	ND	4.9
1,2-Dichloroethane	ND	4.9
1,1-Dichloroethene	ND	4.9
cis-1,2-Dichloroethene	ND	4.9
trans-1,2-Dichloroethene	ND	4.9
1,2-Dichloropropane	ND	4.9
cis-1,3-Dichloropropene	ND	4.9
trans-1,3-Dichloropropene	ND	4.9
Ethylbenzene	ND	4.9
Hexachlorobutadiene	ND	4.9
2-Hexanone	ND	49
Isopropylbenzene	ND	4.9
4-Isopropyltoluene	ND	4.9
Methylene Chloride	ND	9.8

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 50'

 Lab Sample ID:
 720-11927-5
 Date Sampled:
 11/27/2007 1300

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method: 8260B Analysis Batch: 720-29067 Instrument ID: Agilent 75MSD Preparation: 5030B Prep Batch: 720-29041 Lab File ID: 112807012.D Dilution: 1.0 Initial Weight/Volume: 5.12 g Date Analyzed: 11/28/2007 1756 Final Weight/Volume: 10 mL

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		49
Naphthalene	ND		9.8
N-Propylbenzene	ND		4.9
Styrene	ND		4.9
1,1,1,2-Tetrachloroethane	ND		4.9
1,1,2,2-Tetrachloroethane	ND		4.9
Tetrachloroethene	14		4.9
Toluene	ND		4.9
1,2,3-Trichlorobenzene	ND		4.9
1,2,4-Trichlorobenzene	ND		4.9
1,1,1-Trichloroethane	ND		4.9
1,1,2-Trichloroethane	ND		4.9
Trichloroethene	ND		4.9
Trichlorofluoromethane	ND		4.9
1,2,3-Trichloropropane	ND		4.9
1,1,2-Trichloro-1,2,2-trifluoroethar	ne ND		4.9
1,2,4-Trimethylbenzene	ND		4.9
1,3,5-Trimethylbenzene	ND		4.9
Vinyl acetate	ND		49
Vinyl chloride	ND		4.9
Xylenes, Total	ND		9.8
2,2-Dichloropropane	ND		4.9
Surrogate	%Rec		Acceptance Limits
4-Bromofluorobenzene	95		50 - 138
1,2-Dichloroethane-d4 (Surr)	99		66 - 127
Toluene-d8 (Surr)	96		51 - 129

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 50'

 Lab Sample ID:
 720-11927-5
 Date Sampled:
 11/27/2007 1300

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 720-29124 Instrument ID: Varian 3900A

Preparation: 5030B Prep Batch: 720-29065 Lab File ID: c:\saturnws\data\200711\11

Dilution: 1.0 Initial Weight/Volume: 5.41 g

Date Analyzed: 11/29/2007 1334 Final Weight/Volume: 10 mL Date Prepared: 11/29/2007 0939

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL

Gasoline Range Organics (GRO)-C5-C12 ND 0.23

 Surrogate
 %Rec
 Acceptance Limits

 1,2-Dichloroethane-d4 (Surr)
 101
 60 - 140

 Toluene-d8 (Surr)
 96
 70 - 130

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 60'

 Lab Sample ID:
 720-11927-6
 Date Sampled:
 11/27/2007 1300

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807013.DDilution:1.0Initial Weight/Volume:5.00 g

Date Analyzed: 11/28/2007 1821 Final Weight/Volume: 10 mL

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier RL
Methyl tert-butyl ether	ND	5.0
Acetone	ND	50
Benzene	ND	5.0
Dichlorobromomethane	ND	5.0
Bromobenzene	ND	5.0
Chlorobromomethane	ND	20
Bromoform	ND	5.0
Bromomethane	ND	10
2-Butanone (MEK)	ND	50
n-Butylbenzene	ND	5.0
sec-Butylbenzene	ND	5.0
tert-Butylbenzene	ND	5.0
Carbon disulfide	ND	5.0
Carbon tetrachloride	ND	5.0
Chlorobenzene	ND	5.0
Chloroethane	ND	10
Chloroform	ND	5.0
Chloromethane	ND	10
2-Chlorotoluene	ND	5.0
4-Chlorotoluene	ND	5.0
Chlorodibromomethane	ND	5.0
1,2-Dichlorobenzene	ND	5.0
1,3-Dichlorobenzene	ND	5.0
1,4-Dichlorobenzene	ND	5.0
1,3-Dichloropropane	ND	5.0
1,1-Dichloropropene	ND	5.0
1,2-Dibromo-3-Chloropropane	ND	50
Ethylene Dibromide	ND	5.0
Dibromomethane	ND	10
Dichlorodifluoromethane	ND	10
1,1-Dichloroethane	ND	5.0
1,2-Dichloroethane	ND	5.0
1,1-Dichloroethene	ND	5.0
cis-1,2-Dichloroethene	ND	5.0
trans-1,2-Dichloroethene	ND	5.0
1,2-Dichloropropane	ND	5.0
cis-1,3-Dichloropropene	ND	5.0
trans-1,3-Dichloropropene	ND	5.0
Ethylbenzene	ND	5.0
Hexachlorobutadiene	ND	5.0
2-Hexanone	ND	50
Isopropylbenzene	ND	5.0
4-Isopropyltoluene	ND	5.0
Methylene Chloride	ND	10

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 60'

 Lab Sample ID:
 720-11927-6
 Date Sampled:
 11/27/2007 1300

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807013.DDilution:1.0Initial Weight/Volume:5.00 g

Date Analyzed: 11/28/2007 1821 Final Weight/Volume: 10 mL

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier R	L
4-Methyl-2-pentanone (MIBK)	ND	5	0
Naphthalene	ND	1	0
N-Propylbenzene	ND	5	.0
Styrene	ND	5	.0
1,1,1,2-Tetrachloroethane	ND	5	.0
1,1,2,2-Tetrachloroethane	ND	5	.0
Tetrachloroethene	ND	5	.0
Toluene	ND	5	.0
1,2,3-Trichlorobenzene	ND	5	.0
1,2,4-Trichlorobenzene	ND	5	.0
1,1,1-Trichloroethane	ND	5	.0
1,1,2-Trichloroethane	ND	5	.0
Trichloroethene	ND	5	.0
Trichlorofluoromethane	ND	5	.0
1,2,3-Trichloropropane	ND	5	.0
1,1,2-Trichloro-1,2,2-trifluoroethan	e ND	5	.0
1,2,4-Trimethylbenzene	ND	5	.0
1,3,5-Trimethylbenzene	ND	5	.0
Vinyl acetate	ND	5	0
Vinyl chloride	ND	5	.0
Xylenes, Total	ND	1	0
2,2-Dichloropropane	ND	5	.0
Surrogate	%Rec	Acceptance Lir	nits
4-Bromofluorobenzene	97	50 - 138	
1,2-Dichloroethane-d4 (Surr)	98	66 - 127	
Toluene-d8 (Surr)	98	51 - 129	

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 60'

 Lab Sample ID:
 720-11927-6
 Date Sampled:
 11/27/2007 1300

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 720-29124 Instrument ID: Varian 3900A

Preparation: 5030B Prep Batch: 720-29065 Lab File ID: c:\saturnws\data\200711\11

Dilution: 1.0 Initial Weight/Volume: 5.48 g

Date Analyzed: 11/29/2007 1356 Final Weight/Volume: 10 mL

Date Prepared: 11/29/2007 0939

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL

Gasoline Range Organics (GRO)-C5-C12 ND 0.23

 Surrogate
 %Rec
 Acceptance Limits

 1,2-Dichloroethane-d4 (Surr)
 98
 60 - 140

 Toluene-d8 (Surr)
 96
 70 - 130

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 70'

 Lab Sample ID:
 720-11927-7
 Date Sampled:
 11/27/2007 1440

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807014.DDilution:1.0Initial Weight/Volume:5.14

Dilution: 1.0 Initial Weight/Volume: 5.14 g
Date Analyzed: 11/28/2007 1846 Final Weight/Volume: 10 mL

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
Methyl tert-butyl ether	ND		4.9
Acetone	ND		49
Benzene	ND		4.9
Dichlorobromomethane	ND		4.9
Bromobenzene	ND		4.9
Chlorobromomethane	ND		19
Bromoform	ND		4.9
Bromomethane	ND		9.7
2-Butanone (MEK)	ND		49
n-Butylbenzene	ND		4.9
sec-Butylbenzene	ND		4.9
tert-Butylbenzene	ND		4.9
Carbon disulfide	ND		4.9
Carbon tetrachloride	ND		4.9
Chlorobenzene	ND		4.9
Chloroethane	ND		9.7
Chloroform	ND		4.9
Chloromethane	ND		9.7
2-Chlorotoluene	ND		4.9
4-Chlorotoluene	ND		4.9
Chlorodibromomethane	ND		4.9
1,2-Dichlorobenzene	ND		4.9
1,3-Dichlorobenzene	ND		4.9
1,4-Dichlorobenzene	ND		4.9
1,3-Dichloropropane	ND		4.9
1,1-Dichloropropene	ND		4.9
1,2-Dibromo-3-Chloropropane	ND		49
Ethylene Dibromide	ND		4.9
Dibromomethane	ND		9.7
Dichlorodifluoromethane	ND		9.7
1,1-Dichloroethane	ND		4.9
1,2-Dichloroethane	ND		4.9
1,1-Dichloroethene	ND		4.9
cis-1,2-Dichloroethene	ND		4.9
trans-1,2-Dichloroethene	ND		4.9
1,2-Dichloropropane	ND		4.9
cis-1,3-Dichloropropene	ND		4.9
trans-1,3-Dichloropropene	ND		4.9
Ethylbenzene	ND		4.9
Hexachlorobutadiene	ND		4.9
2-Hexanone	ND		49
Isopropylbenzene	ND		4.9
4-Isopropyltoluene	ND		4.9
Methylene Chloride	ND		9.7

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 70'

 Lab Sample ID:
 720-11927-7
 Date Sampled:
 11/27/2007
 1440

 Client Matrix:
 Solid
 Date Received:
 11/27/2007
 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807014.DDilution:1.0Initial Weight/Volume:5.14 g

Date Analyzed: 11/28/2007 1846 Final Weight/Volume: 10 mL

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		49
Naphthalene	ND		9.7
N-Propylbenzene	ND		4.9
Styrene	ND		4.9
1,1,1,2-Tetrachloroethane	ND		4.9
1,1,2,2-Tetrachloroethane	ND		4.9
Tetrachloroethene	ND		4.9
Toluene	ND		4.9
1,2,3-Trichlorobenzene	ND		4.9
1,2,4-Trichlorobenzene	ND		4.9
1,1,1-Trichloroethane	ND		4.9
1,1,2-Trichloroethane	ND		4.9
Trichloroethene	ND		4.9
Trichlorofluoromethane	ND		4.9
1,2,3-Trichloropropane	ND		4.9
1,1,2-Trichloro-1,2,2-trifluoroethar	ne ND		4.9
1,2,4-Trimethylbenzene	ND		4.9
1,3,5-Trimethylbenzene	ND		4.9
Vinyl acetate	ND		49
Vinyl chloride	ND		4.9
Xylenes, Total	ND		9.7
2,2-Dichloropropane	ND		4.9
Surrogate	%Rec	Acc	ceptance Limits
4-Bromofluorobenzene	96	50) - 138
1,2-Dichloroethane-d4 (Surr)	100		6 - 127
Toluene-d8 (Surr)	96	5.	1 - 129

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 70'

 Lab Sample ID:
 720-11927-7
 Date Sampled:
 11/27/2007 1440

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 720-29124 Instrument ID: Varian 3900A

Preparation: 5030B Prep Batch: 720-29065 Lab File ID: c:\saturnws\data\200711\11

Dilution: 1.0 Initial Weight/Volume: 5.12 g

Date Analyzed: 11/29/2007 1419 Final Weight/Volume: 10 mL Date Prepared: 11/29/2007 0939

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL
Gasoline Range Organics (GRO)-C5-C12 ND 0.24

 Surrogate
 %Rec
 Acceptance Limits

 1,2-Dichloroethane-d4 (Surr)
 91
 60 - 140

 Toluene-d8 (Surr)
 95
 70 - 130

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 80'

 Lab Sample ID:
 720-11927-8
 Date Sampled:
 11/27/2007 1515

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807015.DDilution:1.0Initial Weight/Volume:5.08 g

Date Analyzed: 11/28/2007 1911 Final Weight/Volume: 10 mL

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
Methyl tert-butyl ether	ND		4.9
Acetone	ND		49
Benzene	ND		4.9
Dichlorobromomethane	ND		4.9
Bromobenzene	ND		4.9
Chlorobromomethane	ND		20
Bromoform	ND		4.9
Bromomethane	ND		9.8
2-Butanone (MEK)	ND		49
n-Butylbenzene	ND		4.9
sec-Butylbenzene	ND		4.9
tert-Butylbenzene	ND		4.9
Carbon disulfide	ND		4.9
Carbon tetrachloride	ND		4.9
Chlorobenzene	ND		4.9
Chloroethane	ND		9.8
Chloroform	ND		4.9
Chloromethane	ND		9.8
2-Chlorotoluene	ND		4.9
4-Chlorotoluene	ND		4.9
Chlorodibromomethane	ND		4.9
1,2-Dichlorobenzene	ND		4.9
1,3-Dichlorobenzene	ND		4.9
1,4-Dichlorobenzene	ND		4.9
1,3-Dichloropropane	ND		4.9
1,1-Dichloropropene	ND		4.9
1,2-Dibromo-3-Chloropropane	ND		49
Ethylene Dibromide	ND		4.9
Dibromomethane	ND		9.8
Dichlorodifluoromethane	ND		9.8
1,1-Dichloroethane	ND		4.9
1,2-Dichloroethane	ND		4.9
1,1-Dichloroethene	ND		4.9
cis-1,2-Dichloroethene	ND		4.9
trans-1,2-Dichloroethene	ND		4.9
1,2-Dichloropropane	ND		4.9
cis-1,3-Dichloropropene	ND		4.9
trans-1,3-Dichloropropene	ND		4.9
Ethylbenzene	ND		4.9
Hexachlorobutadiene	ND		4.9
2-Hexanone	ND		49
Isopropylbenzene	ND		4.9
4-Isopropyltoluene	ND		4.9
Methylene Chloride	ND		9.8
Wodrylone Ornoride	IND		0.0

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 80'

 Lab Sample ID:
 720-11927-8
 Date Sampled:
 11/27/2007 1515

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807015.DDilution:1.0Initial Weight/Volume:5.08 g

Date Analyzed: 11/28/2007 1911 Final Weight/Volume: 10 mL

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		49
Naphthalene	ND		9.8
N-Propylbenzene	ND		4.9
Styrene	ND		4.9
1,1,1,2-Tetrachloroethane	ND		4.9
1,1,2,2-Tetrachloroethane	ND		4.9
Tetrachloroethene	ND		4.9
Toluene	ND		4.9
1,2,3-Trichlorobenzene	ND		4.9
1,2,4-Trichlorobenzene	ND		4.9
1,1,1-Trichloroethane	ND		4.9
1,1,2-Trichloroethane	ND		4.9
Trichloroethene	ND		4.9
Trichlorofluoromethane	ND		4.9
1,2,3-Trichloropropane	ND		4.9
1,1,2-Trichloro-1,2,2-trifluoroethar	e ND		4.9
1,2,4-Trimethylbenzene	ND		4.9
1,3,5-Trimethylbenzene	ND		4.9
Vinyl acetate	ND		49
Vinyl chloride	ND		4.9
Xylenes, Total	ND		9.8
2,2-Dichloropropane	ND		4.9
Surrogate	%Rec	Ac	ceptance Limits
4-Bromofluorobenzene	95	5	0 - 138
1,2-Dichloroethane-d4 (Surr)	99	6	6 - 127
Toluene-d8 (Surr)	97	5	1 - 129

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 80'

 Lab Sample ID:
 720-11927-8
 Date Sampled:
 11/27/2007 1515

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 720-29124 Instrument ID: Varian 3900A

Preparation: 5030B Prep Batch: 720-29065 Lab File ID: c:\saturnws\data\200711\11

Dilution: 1.0 Initial Weight/Volume: 5.35 g

Date Analyzed: 11/29/2007 1441 Final Weight/Volume: 10 mL

Date Prepared: 11/29/2007 0939

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL

Gasoline Range Organics (GRO)-C5-C12 ND 0.23

 Surrogate
 %Rec
 Acceptance Limits

 1,2-Dichloroethane-d4 (Surr)
 96
 60 - 140

 Toluene-d8 (Surr)
 99
 70 - 130

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 90'

 Lab Sample ID:
 720-11927-9
 Date Sampled:
 11/27/2007 1620

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807016.DDilution:1.0Initial Weight/Volume:5.26 g

Date Analyzed: 11/28/2007 1936 Final Weight/Volume: 5.26 g

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
Methyl tert-butyl ether	ND		4.8
Acetone	ND		48
Benzene	ND		4.8
Dichlorobromomethane	ND		4.8
Bromobenzene	ND		4.8
Chlorobromomethane	ND		19
Bromoform	ND		4.8
Bromomethane	ND		9.5
2-Butanone (MEK)	ND		48
n-Butylbenzene	ND		4.8
sec-Butylbenzene	ND		4.8
tert-Butylbenzene	ND		4.8
Carbon disulfide	ND		4.8
Carbon tetrachloride	ND		4.8
Chlorobenzene	ND		4.8
Chloroethane	ND		9.5
Chloroform	ND		4.8
Chloromethane	ND		9.5
2-Chlorotoluene	ND		4.8
4-Chlorotoluene	ND		4.8
Chlorodibromomethane	ND		4.8
1,2-Dichlorobenzene	ND		4.8
1,3-Dichlorobenzene	ND		4.8
1,4-Dichlorobenzene	ND		4.8
1,3-Dichloropropane	ND		4.8
1,1-Dichloropropene	ND		4.8
1,2-Dibromo-3-Chloropropane	ND		48
Ethylene Dibromide	ND		4.8
Dibromomethane	ND		9.5
Dichlorodifluoromethane	ND		9.5
1,1-Dichloroethane	ND		4.8
1,2-Dichloroethane	ND		4.8
1,1-Dichloroethene	ND		4.8
cis-1,2-Dichloroethene	ND		4.8
trans-1,2-Dichloroethene	ND		4.8
1,2-Dichloropropane	ND		4.8
cis-1,3-Dichloropropene	ND		4.8
trans-1,3-Dichloropropene	ND		4.8
Ethylbenzene	ND		4.8
Hexachlorobutadiene	ND		4.8
2-Hexanone	ND		48
Isopropylbenzene	ND		4.8
4-Isopropyltoluene	ND		4.8
Methylene Chloride	ND		9.5

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 90'

 Lab Sample ID:
 720-11927-9
 Date Sampled:
 11/27/2007 1620

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29067Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29041Lab File ID:112807016.DDilution:1.0Initial Weight/Volume:5.26 g

Date Analyzed: 11/28/2007 1936 Final Weight/Volume: 10 mL

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		48
Naphthalene	ND		9.5
N-Propylbenzene	ND		4.8
Styrene	ND		4.8
1,1,1,2-Tetrachloroethane	ND		4.8
1,1,2,2-Tetrachloroethane	ND		4.8
Tetrachloroethene	ND		4.8
Toluene	ND		4.8
1,2,3-Trichlorobenzene	ND		4.8
1,2,4-Trichlorobenzene	ND		4.8
1,1,1-Trichloroethane	ND		4.8
1,1,2-Trichloroethane	ND		4.8
Trichloroethene	ND		4.8
Trichlorofluoromethane	ND		4.8
1,2,3-Trichloropropane	ND		4.8
1,1,2-Trichloro-1,2,2-trifluoroethar	ne ND		4.8
1,2,4-Trimethylbenzene	ND		4.8
1,3,5-Trimethylbenzene	ND		4.8
Vinyl acetate	ND		48
Vinyl chloride	ND		4.8
Xylenes, Total	ND		9.5
2,2-Dichloropropane	ND		4.8
Surrogate	%Rec		Acceptance Limits
4-Bromofluorobenzene	95		50 - 138
1,2-Dichloroethane-d4 (Surr)	98		66 - 127
Toluene-d8 (Surr)	95		51 - 129

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 90'

 Lab Sample ID:
 720-11927-9
 Date Sampled:
 11/27/2007 1620

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 720-29124 Instrument ID: Varian 3900A

Preparation: 5030B Prep Batch: 720-29065 Lab File ID: c:\saturnws\data\200711\11

Dilution: 1.0 Initial Weight/Volume: 5.22 g

Date Analyzed: 11/29/2007 1503 Final Weight/Volume: 10 mL Date Prepared: 11/29/2007 0939

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL
Gasoline Range Organics (GRO)-C5-C12 ND 0.24

 Surrogate
 %Rec
 Acceptance Limits

 1,2-Dichloroethane-d4 (Surr)
 98
 60 - 140

 Toluene-d8 (Surr)
 97
 70 - 130

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 10'

 Lab Sample ID:
 720-11927-1
 Date Sampled:
 11/27/2007 0925

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8015B Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)-Silica Gel Cleanup

Method: 8015B Analysis Batch: 720-29107 Instrument ID: Varian DRO4

Preparation: 3550B Prep Batch: 720-29037 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 30.13 g
Date Analyzed: 11/29/2007 1237 Final Weight/Volume: 5 mL

Date Prepared: 11/28/2007 1304 Injection Volume:

Column ID: PRIMARY

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL
Diesel Range Organics [C10-C28] ND 1.0
Motor Oil Range Organics [C24-C36] ND 50

 Surrogate
 %Rec
 Acceptance Limits

 Capric Acid (Surr)
 0
 0 - 5

 p-Terphenyl
 89
 41 - 105

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 20'

 Lab Sample ID:
 720-11927-2
 Date Sampled:
 11/27/2007 0930

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8015B Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)-Silica Gel Cleanup

Method: 8015B Analysis Batch: 720-29107 Instrument ID: Varian DRO4

Preparation: 3550B Prep Batch: 720-29037 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 30.21 g

Date Analyzed: 11/29/2007 1303 Final Weight/Volume: 5 mL

Date Prepared: 11/28/2007 1304 Injection Volume:

Column ID: PRIMARY

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL

Diesel Range Organics [C10-C28] 17 0.99

Motor Oil Range Organics [C24-C36] ND 50

 Surrogate
 %Rec
 Acceptance Limits

 Capric Acid (Surr)
 0
 0 - 5

 p-Terphenyl
 102
 41 - 105

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 30'

 Lab Sample ID:
 720-11927-3
 Date Sampled:
 11/27/2007 0940

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8015B Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)-Silica Gel Cleanup

Method: 8015B Analysis Batch: 720-29107 Instrument ID: Varian DRO4

Preparation: 3550B Prep Batch: 720-29037 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 30.28 g

Date Analyzed: 11/29/2007 1330 Final Weight/Volume: 5 mL

Date Prepared: 11/28/2007 1304 Injection Volume:

Column ID: PRIMARY

Analyte	DryWt Corrected: N	Result (mg/Kg)	Qualifier	RL
Diesel Range Organics [C10-C28]	ND		0.99
Motor Oil Range Organics [C24-C	36]	ND		50

Surrogate	%Rec	Acceptance Limits
Capric Acid (Surr)	0	0 - 5
p-Terphenyl	93	41 - 105

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 40'

 Lab Sample ID:
 720-11927-4
 Date Sampled:
 11/27/2007 1010

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8015B Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)-Silica Gel Cleanup

Method: 8015B Analysis Batch: 720-29107 Instrument ID: Varian DRO4

Preparation: 3550B Prep Batch: 720-29037 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 30.41 g
Date Analyzed: 11/29/2007 1356 Final Weight/Volume: 5 mL

Date Prepared: 11/28/2007 1304 Injection Volume:

Column ID: PRIMARY

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL
Diesel Range Organics [C10-C28] ND 0.99
Motor Oil Range Organics [C24-C36] ND 49

 Surrogate
 %Rec
 Acceptance Limits

 Capric Acid (Surr)
 0
 0 - 5

 p-Terphenyl
 88
 41 - 105

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 50'

 Lab Sample ID:
 720-11927-5
 Date Sampled:
 11/27/2007
 1300

 Client Matrix:
 Solid
 Date Received:
 11/27/2007
 1655

8015B Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)-Silica Gel Cleanup

Method: 8015B Analysis Batch: 720-29107 Instrument ID: Varian DRO4

Preparation: 3550B Prep Batch: 720-29037 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 30.38 g

Date Analyzed: 11/29/2007 1144 Final Weight/Volume: 5 mL

Date Prepared: 11/28/2007 1304 Injection Volume:

Column ID: PRIMARY

Analyte	DryWt Corrected: N Result (mg/Kg)	Qualifier	RL
Diesel Range Organics [C10-C28] 1.1		0.99
Motor Oil Range Organics [C24-C	:36] ND		49

Surrogate	%Rec	Acceptance Limits
Capric Acid (Surr)	0	0 - 5
p-Terphenyl	89	41 - 105

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 60'

 Lab Sample ID:
 720-11927-6
 Date Sampled:
 11/27/2007 1300

 Client Matrix:
 Solid
 Date Received:
 11/27/2007 1655

8015B Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)-Silica Gel Cleanup

Method: 8015B Analysis Batch: 720-29107 Instrument ID: Varian DRO4

Preparation: 3550B Prep Batch: 720-29037 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 30.25 g
Date Analyzed: 11/29/2007 1541 Final Weight/Volume: 5 mL

Date Prepared: 11/28/2007 1304 Injection Volume:

Column ID: PRIMARY

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL
Diesel Range Organics [C10-C28] ND 0.99
Motor Oil Range Organics [C24-C36] ND 50

 Surrogate
 %Rec
 Acceptance Limits

 Capric Acid (Surr)
 0
 0 - 5

 p-Terphenyl
 92
 41 - 105

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 70'

 Lab Sample ID:
 720-11927-7
 Date Sampled:
 11/27/2007
 1440

 Client Matrix:
 Solid
 Date Received:
 11/27/2007
 1655

8015B Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)-Silica Gel Cleanup

Method: 8015B Analysis Batch: 720-29107 Instrument ID: Varian DRO4

Preparation: 3550B Prep Batch: 720-29037 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 30.46 g

Date Analyzed: 11/29/2007 1607 Final Weight/Volume: 5 mL

Date Prepared: 11/28/2007 1304 Injection Volume:

Column ID: PRIMARY

Analyte	DryWt Corrected: N Result (mg/Kg) Qualifier	RL
Diesel Range Organics [C10-C28] ND		0.98
Motor Oil Range Organics [C24-C	36] ND		49

Surrogate	%Rec	Acceptance Limits
Capric Acid (Surr)	0	0 - 5
p-Terphenyl	90	41 - 105

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 80'

 Lab Sample ID:
 720-11927-8
 Date Sampled:
 11/27/2007
 1515

 Client Matrix:
 Solid
 Date Received:
 11/27/2007
 1655

8015B Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)-Silica Gel Cleanup

Method: 8015B Analysis Batch: 720-29107 Instrument ID: Varian DRO4

Preparation: 3550B Prep Batch: 720-29037 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 30.05 g

Date Analyzed: 11/29/2007 1634 Final Weight/Volume: 5 mL
Date Prepared: 11/28/2007 1304 Injection Volume:

Column ID: PRIMARY

Odumin ID. Transaction

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL
Diesel Range Organics [C10-C28] ND 1.0
Motor Oil Range Organics [C24-C36] ND 50

Surrogate %Rec Acceptance Limits

 Capric Acid (Surr)
 0
 0 - 5

 p-Terphenyl
 85
 41 - 105

Client: Engeo, Inc. Job Number: 720-11927-1

Client Sample ID: S @ 90'

Lab Sample ID: 720-11927-9 Date Sampled: 11/27/2007 1620 Client Matrix: Solid Date Received: 11/27/2007 1655

8015B Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)-Silica Gel Cleanup

8015B Analysis Batch: 720-29107 Method: Instrument ID: Varian DRO4

Preparation: 3550B Prep Batch: 720-29037 Lab File ID: N/A

30.23 g Dilution: 1.0 Initial Weight/Volume:

Date Analyzed: 11/29/2007 1700 Final Weight/Volume: 5 mL

Date Prepared: 11/28/2007 1304 Injection Volume: Column ID:

PRIMARY

DryWt Corrected: N Result (mg/Kg) Qualifier Analyte RLDiesel Range Organics [C10-C28] ND 0.99 Motor Oil Range Organics [C24-C36] ND 50

%Rec Surrogate Acceptance Limits Capric Acid (Surr) 0 0 - 5 p-Terphenyl 92 41 - 105

DATA REPORTING QUALIFIERS

Lab Section Qualifier Description

Client: Engeo, Inc. Job Number: 720-11927-1

QC Association Summary

Lab Sample ID	Client Sample ID	Report Basis	Client Matrix	Method	Prep Batch
GC/MS VOA					
Prep Batch: 720-29041					
LCS 720-29041/1-A	Lab Control Spike	T	Solid	5030B	
LCSD 720-29041/2-A	Lab Control Spike Duplicate	Т	Solid	5030B	
MB 720-29041/3-A	Method Blank	T	Solid	5030B	
720-11927-1	S @ 10'	T	Solid	5030B	
720-11927-1MS	Matrix Spike	T	Solid	5030B	
720-11927-1MSD	Matrix Spike Duplicate	Т	Solid	5030B	
720-11927-2	S @ 20'	T	Solid	5030B	
720-11927-3	S @ 30'	T	Solid	5030B	
720-11927-4	S @ 40'	T	Solid	5030B	
720-11927-5	S @ 50'	T	Solid	5030B	
720-11927-6	S @ 60'	Т	Solid	5030B	
720-11927-7	S @ 70'	T	Solid	5030B	
720-11927-8	S @ 80'	Т	Solid	5030B	
720-11927-9	S @ 90'	Т	Solid	5030B	
Prep Batch: 720-29065					
LCS 720-29065/2-A	Lab Control Spike	T	Solid	5030B	
LCSD 720-29065/3-A	Lab Control Spike Duplicate	Т	Solid	5030B	
MB 720-29065/1-A	Method Blank	T	Solid	5030B	
720-11927-1	S @ 10'	Т	Solid	5030B	
720-11927-1MS	Matrix Spike	T	Solid	5030B	
720-11927-1MSD	Matrix Spike Duplicate	Т	Solid	5030B	
720-11927-2	S @ 20'	Т	Solid	5030B	
720-11927-3	S @ 30'	Т	Solid	5030B	
720-11927-4	S @ 40'	Т	Solid	5030B	
720-11927-5	S @ 50'	Т	Solid	5030B	
720-11927-6	S @ 60'	Т	Solid	5030B	
720-11927-7	S @ 70'	Т	Solid	5030B	
720-11927-8	S @ 80'	Т	Solid	5030B	
720-11927-9	S @ 90'	Т	Solid	5030B	

Client: Engeo, Inc. Job Number: 720-11927-1

QC Association Summary

		Report			
Lab Sample ID	Client Sample ID	Basis	Client Matrix	Method	Prep Batch
GC/MS VOA					
Analysis Batch:720-2	9067				
LCS 720-29041/1-A	Lab Control Spike	T	Solid	8260B	720-29041
LCSD 720-29041/2-A	Lab Control Spike Duplicate	T	Solid	8260B	720-29041
MB 720-29041/3-A	Method Blank	T	Solid	8260B	720-29041
720-11927-1	S @ 10'	T	Solid	8260B	720-29041
720-11927-1MS	Matrix Spike	T	Solid	8260B	720-29041
720-11927-1MSD	Matrix Spike Duplicate	Т	Solid	8260B	720-29041
720-11927-2	S @ 20'	T	Solid	8260B	720-29041
720-11927-3	S @ 30'	Т	Solid	8260B	720-29041
720-11927-4	S @ 40'	Т	Solid	8260B	720-29041
720-11927-5	S @ 50'	Т	Solid	8260B	720-29041
720-11927-6	S @ 60'	T	Solid	8260B	720-29041
720-11927-7	S @ 70'	T	Solid	8260B	720-29041
720-11927-8	S @ 80'	Т	Solid	8260B	720-29041
720-11927-9	S @ 90'	Т	Solid	8260B	720-29041
Analysis Batch:720-2	9124				
LCS 720-29065/2-A	Lab Control Spike	Т	Solid	8260B	720-29065
LCSD 720-29065/3-A	Lab Control Spike Duplicate	Т	Solid	8260B	720-29065
MB 720-29065/1-A	Method Blank	Т	Solid	8260B	720-29065
720-11927-1	S @ 10'	Т	Solid	8260B	720-29065
720-11927-1MS	Matrix Spike	Ť	Solid	8260B	720-29065
720-11927-1MSD	Matrix Spike Duplicate	Ť	Solid	8260B	720-29065
720-11927-2	S @ 20'	Ť	Solid	8260B	720-29065
720-11927-3	S @ 30'	Ť	Solid	8260B	720-29065
720-11927-4	S @ 40'	Ť	Solid	8260B	720-29065
720-11927-5	S @ 50'	Ť	Solid	8260B	720-29065
720-11927-6	S @ 60'	Ť	Solid	8260B	720-29065
720-11927-7	S @ 70'	Ť	Solid	8260B	720-29065
720-11927-8	S @ 80'	Ť	Solid	8260B	720-29065
720-11927-9	S @ 90'	Ť	Solid	8260B	720-29065

Report Basis T = Total

Client: Engeo, Inc. Job Number: 720-11927-1

QC Association Summary

		Report			
Lab Sample ID	Client Sample ID	Basis	Client Matrix	Method	Prep Batch
GC Semi VOA					
Prep Batch: 720-2903					
LCS 720-29037/2-A	Lab Control Spike	Α	Solid	3550B	
LCSD 720-29037/3-A	Lab Control Spike Duplicate	Α	Solid	3550B	
MB 720-29037/1-A	Method Blank	Α	Solid	3550B	
720-11927-1	S @ 10'	Α	Solid	3550B	
720-11927-1MS	Matrix Spike	Α	Solid	3550B	
720-11927-1MSD	Matrix Spike Duplicate	Α	Solid	3550B	
720-11927-2	S @ 20'	Α	Solid	3550B	
720-11927-3	S @ 30'	Α	Solid	3550B	
720-11927-4	S @ 40'	Α	Solid	3550B	
720-11927-5	S @ 50'	Α	Solid	3550B	
720-11927-6	S @ 60'	Α	Solid	3550B	
720-11927-7	S @ 70'	Α	Solid	3550B	
720-11927-8	S @ 80'	Α	Solid	3550B	
720-11927-9	S @ 90'	Α	Solid	3550B	
Analysis Batch:720-29	9107				
LCS 720-29037/2-A	Lab Control Spike	Α	Solid	8015B	720-29037
LCSD 720-29037/3-A	Lab Control Spike Duplicate	A	Solid	8015B	720-29037
MB 720-29037/1-A	Method Blank	A	Solid	8015B	720-29037
720-11927-1	S @ 10'	A	Solid	8015B	720-29037
720-11927-1MS	Matrix Spike	A	Solid	8015B	720-29037
720-11927-1MSD	Matrix Spike Duplicate	A	Solid	8015B	720-29037
720-11927-2	S @ 20'	A	Solid	8015B	720-29037
720-11927-3	S @ 30'	A	Solid	8015B	720-29037
720-11927-4	S @ 40'	A	Solid	8015B	720-29037
720-11927-5	S @ 50'	A	Solid	8015B	720-29037
720-11927-6	S @ 60'	A	Solid	8015B	720-29037
720-11927-7	S @ 70'	A	Solid	8015B	720-29037
720-11927-8	S @ 80'	A	Solid	8015B	720-29037
720-11927-9	S @ 90'	A	Solid	8015B	720-29037

Report Basis

A = Silica Gel Cleanup

Job Number: 720-11927-1 Client: Engeo, Inc.

Method Blank - Batch: 720-29041 Method: 8260B Preparation: 5030B

Lab Sample ID: MB 720-29041/3-A Analysis Batch: 720-29067 Instrument ID: Agilent 75MSD Client Matrix: Solid Prep Batch: 720-29041 Lab File ID: 112807005.D Dilution: 1.0 Units: ug/Kg Initial Weight/Volume: 5 g

Final Weight/Volume: 10 mL Date Analyzed: 11/28/2007 1447 Date Prepared: 11/28/2007 1330

Analyte	Result	Qual	RL
Methyl tert-butyl ether	ND		5.0
Acetone	ND		50
Benzene	ND		5.0
Dichlorobromomethane	ND		5.0
Bromobenzene	ND		5.0
Chlorobromomethane	ND		20
Bromoform	ND		5.0
Bromomethane	ND		10
2-Butanone (MEK)	ND		50
n-Butylbenzene	ND		5.0
sec-Butylbenzene	ND		5.0
tert-Butylbenzene	ND		5.0
Carbon disulfide	ND		5.0
Carbon tetrachloride	ND		5.0
Chlorobenzene	ND		5.0
Chloroethane	ND		10
Chloroform	ND		5.0
Chloromethane	ND		10
2-Chlorotoluene	ND		5.0
4-Chlorotoluene	ND		5.0
Chlorodibromomethane	ND		5.0
1,2-Dichlorobenzene	ND		5.0
1,3-Dichlorobenzene	ND		5.0
1,4-Dichlorobenzene	ND		5.0
1,3-Dichloropropane	ND		5.0
1,1-Dichloropropene	ND		5.0
1,2-Dibromo-3-Chloropropane	ND		50
Ethylene Dibromide	ND		5.0
Dibromomethane	ND		10
Dichlorodifluoromethane	ND		10
1,1-Dichloroethane	ND		5.0
1,2-Dichloroethane	ND		5.0
1,1-Dichloroethene	ND		5.0
cis-1,2-Dichloroethene	ND		5.0
trans-1,2-Dichloroethene	ND		5.0
1,2-Dichloropropane	ND		5.0
cis-1,3-Dichloropropene	ND		5.0
trans-1,3-Dichloropropene	ND		5.0
Ethylbenzene	ND		5.0
Hexachlorobutadiene	ND		5.0
2-Hexanone	ND		50

Calculations are performed before rounding to avoid round-off errors in calculated results.

Client: Engeo, Inc. Job Number: 720-11927-1

Method Blank - Batch: 720-29041 Method: 8260B Preparation: 5030B

Lab Sample ID: MB 720-29041/3-A Analysis Batch: 720-29067 Instrument ID: Agilent 75MSD

Client Matrix: Solid Prep Batch: 720-29041 Lab File ID: 112807005.D Dilution: 1.0 Units: ug/Kg Initial Weight/Volume: 5 g

Date Analyzed: 11/28/2007 1447 Final Weight/Volume: 10 mL Date Prepared: 11/28/2007 1330

Analyte	Result	Qual	RL
Isopropylbenzene	ND		5.0
4-Isopropyltoluene	ND		5.0
Methylene Chloride	ND		10
4-Methyl-2-pentanone (MIBK)	ND		50
Naphthalene	ND		10
N-Propylbenzene	ND		5.0
Styrene	ND		5.0
1,1,1,2-Tetrachloroethane	ND		5.0
1,1,2,2-Tetrachloroethane	ND		5.0
Tetrachloroethene	ND		5.0
Toluene	ND		5.0
1,2,3-Trichlorobenzene	ND		5.0
1,2,4-Trichlorobenzene	ND		5.0
1,1,1-Trichloroethane	ND		5.0
1,1,2-Trichloroethane	ND		5.0
Trichloroethene	ND		5.0
Trichlorofluoromethane	ND		5.0
1,2,3-Trichloropropane	ND		5.0
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		5.0
1,2,4-Trimethylbenzene	ND		5.0
1,3,5-Trimethylbenzene	ND		5.0
Vinyl acetate	ND		50
Vinyl chloride	ND		5.0
Xylenes, Total	ND		10
2,2-Dichloropropane	ND		5.0
Surrogate	% Rec	Acceptance Limits	
4-Bromofluorobenzene	92	50 - 138	
1,2-Dichloroethane-d4 (Surr)	93	66 - 127	
Toluene-d8 (Surr)	90	51 - 129	

Calculations are performed before rounding to avoid round-off errors in calculated results.

Client: Engeo, Inc. Job Number: 720-11927-1

Lab Control Spike/ Method: 8260B
Lab Control Spike Duplicate Recovery Report - Batch: 720-29041 Preparation: 5030B

LCS Lab Sample ID: LCS 720-29041/1-A

Client Matrix: Solid Dilution: 1.0

Date Analyzed: 11/28/2007 1357 Date Prepared: 11/28/2007 1330 Analysis Batch: 720-29067 Prep Batch: 720-29041

Units: ug/Kg

Instrument ID: Agilent 75MSD Lab File ID: 112807003.D

Initial Weight/Volume: 5 g Final Weight/Volume: 10 mL

LCSD Lab Sample ID: LCSD 720-29041/2-A

Client Matrix: Solid Dilution: 1.0

Date Analyzed: 11/28/2007 1422 Date Prepared: 11/28/2007 1330 Analysis Batch: 720-29067 Prep Batch: 720-29041

Units: ug/Kg

Instrument ID: Agilent 75MSD Lab File ID: 112807004.D

Initial Weight/Volume: 5 g Final Weight/Volume: 10 mL

	9	Rec. €					
Analyte	LCS	LCSD	Limit	RPD	RPD Limit	LCS Qual	LCSD Qual
Benzene	91	91	80 - 120	1	20		
Chlorobenzene	90	91	86 - 115	1	20		
1,1-Dichloroethene	95	94	81 - 140	0	20		
Toluene	90	92	81 - 120	2	20		
Trichloroethene	92	92	82 - 118	0	20		
Surrogate	L	.CS % Rec	LCSD %	Rec	Acce	otance Limits	:
4-Bromofluorobenzene	9	9	98		5	0 - 138	
1,2-Dichloroethane-d4 (Surr)	9	8	99		6	66 - 127	
Toluene-d8 (Surr)	9	6	96		5	1 - 129	

112807007.D

Job Number: 720-11927-1 Client: Engeo, Inc.

Matrix Spike/ Method: 8260B Matrix Spike Duplicate Recovery Report - Batch: 720-29041 Preparation: 5030B

MS Lab Sample ID: 720-11927-1 Analysis Batch: 720-29067 Instrument ID: Agilent 75MSD Client Matrix: Solid Prep Batch: 720-29041 Lab File ID:

Initial Weight/Volume: 5.00 g Dilution: 1.0 Final Weight/Volume: 10 mL Date Analyzed: 11/28/2007 1551 Date Prepared: 11/28/2007 1330

MSD Lab Sample ID: 720-11927-1 Analysis Batch: 720-29067 Instrument ID: Agilent 75MSD Client Matrix: Solid Prep Batch: 720-29041 112807008.D Lab File ID:

Dilution: 1.0 Initial Weight/Volume: 5.07 g Date Analyzed: 11/28/2007 1616 Final Weight/Volume: 10 mL Date Prepared: 11/28/2007 1330

% Rec. MS MSD **RPD RPD Limit** MS Qual MSD Qual Analyte Limit 63 - 126 Benzene 91 90 2 20 Chlorobenzene 88 88 57 - 124 1 20 1,1-Dichloroethene 93 94 66 - 149 1 20 54 - 131 Toluene 90 89 3 20 53 - 130 2 20 Trichloroethene 91 91 MS % Rec MSD % Rec Surrogate Acceptance Limits 91 4-Bromofluorobenzene 88 50 - 138 66 - 127 98 95 1,2-Dichloroethane-d4 (Surr) 51 - 129 Toluene-d8 (Surr) 89 94

Job Number: 720-11927-1 Client: Engeo, Inc.

Method Blank - Batch: 720-29065 Method: 8260B Preparation: 5030B

Lab Sample ID: MB 720-29065/1-A Analysis Batch: 720-29124 Instrument ID: Varian 3900A

Prep Batch: 720-29065 Lab File ID: c:\saturnws\data\200711\11 Client Matrix: Solid

Units: mg/Kg Initial Weight/Volume: 5.00 g Dilution: 1.0

Date Analyzed: 11/29/2007 1035 Final Weight/Volume: 10 mL Date Prepared: 11/29/2007 0939

Analyte	Result	Qual	RL
Benzene	ND		0.0050
Toluene	ND		0.0050
Gasoline Range Organics (GRO)-C5-C12	ND		0.25
Surrogate	% Rec	Acceptance	Limits
1,2-Dichloroethane-d4 (Surr)	103	60 - 140	
Toluene-d8 (Surr)	101	70 - 130	1

Lab Control Spike/ Method: 8260B Lab Control Spike Duplicate Recovery Report - Batch: 720-29065 Preparation: 5030B

LCS Lab Sample ID: LCS 720-29065/2-A Analysis Batch: 720-29124 Instrument ID: Varian 3900A

Client Matrix: Solid Prep Batch: 720-29065 Lab File ID: c:\saturnws\data\200711\11

Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 5.00 g Date Analyzed: 11/29/2007 0844 Final Weight/Volume: 10 mL

Date Prepared: 11/29/2007 0939

LCSD Lab Sample ID: LCSD 720-29065/3-A Analysis Batch: 720-29124 Instrument ID: Varian 3900A

Client Matrix: Prep Batch: 720-29065 Solid Lab File ID: c:\saturnws\data\200711\112

Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 5.00 g

Date Analyzed: 11/29/2007 0950 Final Weight/Volume: 10 mL Date Prepared: 11/29/2007 0939

% Rec. RPD Limit LCS Qual LCSD Qual Analyte LCS **LCSD** Limit **RPD** Benzene 90 103 70 - 123 13 20 Toluene 94 100 81 - 128 20 6 51 - 97 Gasoline Range Organics (GRO)-C5-C12 74 7 20 69 LCS % Rec LCSD % Rec Surrogate Acceptance Limits 1.2-Dichloroethane-d4 (Surr) 102 107 60 - 140 Toluene-d8 (Surr) 99 93 70 - 130

Client: Engeo, Inc. Job Number: 720-11927-1

Matrix Spike/ Method: 8260B
Matrix Spike Duplicate Recovery Report - Batch: 720-29065 Preparation: 5030B

MS Lab Sample ID: 720-11927-1 Analysis Batch: 720-29124 Instrument ID: Varian 3900A

Client Matrix: Solid Prep Batch: 720-29065 Lab File ID: c:\saturnws\data\200711\

Dilution: 1.0 Initial Weight/Volume: 5.32 g
Date Analyzed: 11/29/2007 1120 Final Weight/Volume: 10 mL

Date Prepared: 11/29/2007 0939

MSD Lab Sample ID: 720-11927-1 Analysis Batch: 720-29124 Instrument ID: Varian 3900A

Client Matrix: Solid Prep Batch: 720-29065 Lab File ID: c:\saturnws\data\200711\11

Dilution: 1.0 Initial Weight/Volume: 5.14 g

Date Analyzed: 11/29/2007 1142 Final Weight/Volume: 10 mL
Date Prepared: 11/29/2007 0939

	9	<u>6 Rec.</u>				
Analyte	MS	MSD	Limit	RPD	RPD Limit	MS Qual MSD Qual
Benzene	94	90	70 - 123	0	20	
Toluene	95	92	81 - 128	0	20	
Gasoline Range Organics (GRO)-C5-C12	68	63	51 - 97	4	20	
Surrogate		MS % Rec	MSD 9	% Rec	Acce	ptance Limits
1,2-Dichloroethane-d4 (Surr)		93	91		60	0 - 140
Toluene-d8 (Surr)		93	99		70	0 - 130

41 - 105

Client: Engeo, Inc. Job Number: 720-11927-1

Method Blank - Batch: 720-29037 Method: 8015B
Preparation: 3550B
Silica Gel Cleanup

Lab Sample ID: MB 720-29037/1-A Analysis Batch: 720-29107 Instrument ID: Varian DRO4

Client Matrix: Solid Prep Batch: 720-29037 Lab File ID: N/A

Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 30.31 g
Date Analyzed: 11/29/2007 1118 Final Weight/Volume: 5 mL

Date Prepared: 11/28/2007 1304 Injection Volume:

Column ID: PRIMARY

Result Qual RLAnalyte Diesel Range Organics [C10-C28] ND 0.99 Motor Oil Range Organics [C24-C36] ND 49 Surrogate % Rec Acceptance Limits Capric Acid (Surr) 0 0 - 5

p-Terphenyl 98 41 - 105

Lab Control Spike/
Lab Control Spike Duplicate Recovery Report - Batch: 720-29037

Method: 8015B

Preparation: 3550B

Silica Gel Cleanup

LCS Lab Sample ID: LCS 720-29037/2-A Analysis Batch: 720-29107 Instrument ID: Varian DRO4

Client Matrix: Solid Prep Batch: 720-29037 Lab File ID: N/A
Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 30.14

Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 30.14 g
Date Analyzed: 11/29/2007 1025 Final Weight/Volume: 5 mL
Date Prepared: 11/28/2007 1304 Injection Volume:
Column ID: PRIMARY

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LCSD Lab Sample ID: LCSD 720-29037/3-A Analysis Batch: 720-29107 Instrument ID: Varian DRO4 Client Matrix: Solid Prep Batch: 720-29037 Lab File ID: N/A

Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 30.38 g

Date Analyzed: 11/29/2007 1051 Final Weight/Volume: 5 mL
Date Prepared: 11/28/2007 1304 Injection Volume:

Column ID: PRIMARY

86

% Rec. Analyte LCS **LCSD** Limit **RPD** RPD Limit LCS Qual LCSD Qual Diesel Range Organics [C10-C28] 74 71 50 - 130 30 4 LCS % Rec Surrogate LCSD % Rec Acceptance Limits

87

Calculations are performed before rounding to avoid round-off errors in calculated results.

p-Terphenyl

41 - 105

Job Number: 720-11927-1 Client: Engeo, Inc.

Matrix Spike/ Method: 8015B Matrix Spike Duplicate Recovery Report - Batch: 720-29037 Preparation: 3550B Silica Gel Cleanup

MS Lab Sample ID: 720-11927-1 Analysis Batch: 720-29107 Instrument ID: Varian DRO4

Client Matrix: Prep Batch: 720-29037 Solid Lab File ID: N/A

Dilution: 1.0

Initial Weight/Volume: 30.33 g Date Analyzed: 11/29/2007 1726 Final Weight/Volume: 5 mL

Date Prepared: 11/28/2007 1304 Injection Volume:

Column ID: **PRIMARY**

MSD Lab Sample ID: 720-11927-1 Instrument ID: Varian DRO4 Analysis Batch: 720-29107 Client Matrix: Solid Prep Batch: 720-29037 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 30.44 g

Date Analyzed: 11/29/2007 1753 Final Weight/Volume: 5 mL

Date Prepared: 11/28/2007 1304 Injection Volume: Column ID: **PRIMARY**

96

% Rec. MS MSD Limit **RPD** MS Qual MSD Qual Analyte **RPD Limit** Diesel Range Organics [C10-C28] 50 - 130 62 62 1 30 Surrogate MS % Rec MSD % Rec Acceptance Limits

85

Calculations are performed before rounding to avoid round-off errors in calculated results.

p-Terphenyl

Selection Date Da

Login Sample Receipt Check List

Client: Engeo, Inc.

Job Number: 720-11927-1

List Source: TestAmerica San Francisco

Login Number: 11927

Creator: Bullock, Tracy

List Number: 1

Question	T / F/ NA	Comment
Radioactivity either was not measured or, if measured, is at or below background	N/A	
The cooler's custody seal, if present, is intact.	N/A	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
There are no discrepancies between the sample IDs on the containers and the COC.	True	
Samples are received within Holding Time.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
VOA sample vials do not have headspace or bubble is <6mm (1/4") in diameter.	True	
If necessary, staff have been informed of any short hold time or quick TAT needs	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	



ANALYTICAL REPORT

Job Number: 720-11961-1

Job Description: 224 Rickenbacker Circle

For:

Engeo, Inc. 2010 Crow Canyon Place Suite 250 San Ramon, CA 94583

Attention: Ms. Kelly Krohn



Melissa Brewer
Project Manager I
melissa.brewer@testamericainc.com
12/07/2007

Job Narrative 720-J11961-1

Comments

No additional comments.

Receipt

DB-1 (GW@70') 1 of 2 ambers did not have the depth on bottles. Through the process of elimination labeled as 70'. DB-1 (GW@95') 1 of 2 of the ambers: depth was 90" time matched with COC.

All other samples were received in good condition within temperature requirements.

GC/MS VOA

Method 8260B: Surrogate 1,2-DCA-d4 was outside control limits for 11931-1-MSD for batch 29173. The associated laboratory control standard (LCS) met acceptance criteria.

No other analytical or quality issues were noted.

GC Semi VOA

No analytical or quality issues were noted.

Organic Prep

No analytical or quality issues were noted.

EXECUTIVE SUMMARY - Detections

Client: Engeo, Inc. Job Number: 720-11961-1

Lab Sample ID	Client Sample ID		Reporting			
Analyte		Result / Qualifier	Limit	Units	Method	
720-11961-2	DB-1 (GW@70')					
Benzene		2.7	0.50	ug/L	8260B	
Gasoline Range Or	rganics (GRO)-C5-C12	98	50	ug/L	8260B	
Ethylbenzene		1.2	0.50	ug/L	8260B	
Toluene		2.3	0.50	ug/L	8260B	
Xylenes, Total		1.0	1.0	ug/L	8260B	
Silica Gel Cleanu	v					
Diesel Range Orga	nics [C10-C28]	190	50	ug/L	8015B	

METHOD SUMMARY

Client: Engeo, Inc.

Job Number: 720-11961-1

Description	Lab Location	Method	Preparation Method
Matrix: Water			
Volatile Organic Compounds by GC/MS	TAL SF	SW846 8260B	
Volatile Organic Compounds by GC/MS (Low Level)	TAL SF	SW846 8260B	
Purge-and-Trap for Aqueous Samples/Unpreserved	TAL SF		SW846 5030B
Purge-and-Trap for Aqueous Samples/Unpreserved	TAL SF		SW846 5030B
Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)	TAL SF	SW846 8015B	
Separatory Funnel Liquid-Liquid Extraction	TAL SF		SW846 3510C SGC

Lab References:

TAL SF = TestAmerica San Francisco

Method References:

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

SAMPLE SUMMARY

Client: Engeo, Inc. Job Number: 720-11961-1

Lab Sample ID	Client Sample ID	Client Matrix	Date/Time Sampled	Date/Time Received
720-11961-1	DB-1 (GW@35')	Water	11/28/2007 1000	11/28/2007 1810
720-11961-2	DB-1 (GW@70')	Water	11/28/2007 1340	11/28/2007 1810
720-11961-3	DB-1 (GW@95')	Water	11/28/2007 1520	11/28/2007 1810

Client: Engeo, Inc. Job Number: 720-11961-1

Client Sample ID: DB-1 (GW@35')

 Lab Sample ID:
 720-11961-1
 Date Sampled:
 11/28/2007 1000

 Client Matrix:
 Water
 Date Received:
 11/28/2007 1810

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method: 8260B Analysis Batch: 720-29171 Instrument ID: Varian 3900G

Preparation: 5030B Lab File ID: c:\saturnws\data\200712\12

Dilution: 1.0 Initial Weight/Volume: 40 mL Date Analyzed: 12/03/2007 1351 Final Weight/Volume: 40 mL

Date Prepared: 12/03/2007 1351

Analyte	Result (ug/L)	Qualifier	RL
Methyl tert-butyl ether	ND		5.0
Acetone	ND		50
Benzene	ND		0.50
Dichlorobromomethane	ND		0.50
Bromobenzene	ND		1.0
Chlorobromomethane	ND		1.0
Bromoform	ND		1.0
Bromomethane	ND		1.0
2-Butanone (MEK)	ND		50
n-Butylbenzene	ND		1.0
sec-Butylbenzene	ND		1.0
tert-Butylbenzene	ND		1.0
Carbon disulfide	ND		5.0
Carbon tetrachloride	ND		0.50
Chlorobenzene	ND		0.50
Chloroethane	ND		1.0
Chloroform	ND		1.0
Chloromethane	ND		1.0
2-Chlorotoluene	ND		0.50
4-Chlorotoluene	ND		0.50
Chlorodibromomethane	ND		0.50
1,2-Dichlorobenzene	ND		0.50
1,3-Dichlorobenzene	ND		0.50
1,4-Dichlorobenzene	ND		0.50
1,3-Dichloropropane	ND		1.0
1,1-Dichloropropene	ND		0.50
1,2-Dibromo-3-Chloropropane	ND		1.0
Ethylene Dibromide	ND		0.50
Dibromomethane	ND		0.50
Dichlorodifluoromethane	ND		0.50
1,1-Dichloroethane	ND		0.50
1,2-Dichloroethane	ND		0.50
1,1-Dichloroethene	ND		0.50
cis-1,2-Dichloroethene	ND		0.50
trans-1,2-Dichloroethene	ND		0.50
1,2-Dichloropropane	ND		0.50
cis-1,3-Dichloropropene	ND		0.50
trans-1,3-Dichloropropene	ND		0.50
Ethylbenzene	ND		0.50
Hexachlorobutadiene	ND		1.0
2-Hexanone	ND		50
Isopropylbenzene	ND		0.50
4-Isopropyltoluene	ND		1.0
Methylene Chloride	ND		5.0

Client: Engeo, Inc. Job Number: 720-11961-1

Client Sample ID: DB-1 (GW@35')

 Lab Sample ID:
 720-11961-1
 Date Sampled:
 11/28/2007 1000

 Client Matrix:
 Water
 Date Received:
 11/28/2007 1810

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method: 8260B Analysis Batch: 720-29171 Instrument ID: Varian 3900G

Preparation: 5030B Lab File ID: c:\saturnws\data\200712\12

Dilution: 1.0 Initial Weight/Volume: 40 mL Date Analyzed: 12/03/2007 1351 Final Weight/Volume: 40 mL

Date Prepared: 12/03/2007 1351

Analyte	Result (ug/L)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		50
Naphthalene	ND		1.0
N-Propylbenzene	ND		1.0
Styrene	ND		0.50
1,1,1,2-Tetrachloroethane	ND		0.50
1,1,2,2-Tetrachloroethane	ND		0.50
Tetrachloroethene	ND		0.50
Toluene	ND		0.50
1,2,3-Trichlorobenzene	ND		1.0
1,2,4-Trichlorobenzene	ND		1.0
1,1,1-Trichloroethane	ND		0.50
1,1,2-Trichloroethane	ND		0.50
Trichloroethene	ND		0.50
Trichlorofluoromethane	ND		1.0
1,2,3-Trichloropropane	ND		0.50
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		0.50
1,2,4-Trimethylbenzene	ND		0.50
1,3,5-Trimethylbenzene	ND		0.50
Vinyl acetate	ND		50
Vinyl chloride	ND		0.50
Xylenes, Total	ND		1.0
2,2-Dichloropropane	ND		0.50
Surrogate	%Rec	Acceptance	Limits
4-Bromofluorobenzene	115	83 - 127	
1,2-Dichloroethane-d4 (Surr)	103	86 - 129	
Toluene-d8 (Surr)	107	82 - 126	

Client: Engeo, Inc. Job Number: 720-11961-1

Client Sample ID: DB-1 (GW@35')

 Lab Sample ID:
 720-11961-1
 Date Sampled:
 11/28/2007 1000

 Client Matrix:
 Water
 Date Received:
 11/28/2007 1810

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 720-29173 Instrument ID: Varian 3900A

Preparation: 5030B Lab File ID: c:\saturnws\data\200711\11

Dilution: 1.0 Initial Weight/Volume: 10 mL

Date Analyzed: 11/30/2007 1817 Final Weight/Volume: 10 mL

Date Prepared: 11/30/2007 1817

Analyte Result (ug/L) Qualifier RL
Gasoline Range Organics (GRO)-C5-C12 ND 50

 Surrogate
 %Rec
 Acceptance Limits

 Toluene-d8 (Surr)
 97
 77 - 121

 1,2-Dichloroethane-d4 (Surr)
 101
 73 - 130

Client: Engeo, Inc. Job Number: 720-11961-1

Client Sample ID: DB-1 (GW@70')

 Lab Sample ID:
 720-11961-2
 Date Sampled:
 11/28/2007
 1340

 Client Matrix:
 Water
 Date Received:
 11/28/2007
 1810

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method: 8260B Analysis Batch: 720-29171 Instrument ID: Varian 3900G

Preparation: 5030B Lab File ID: c:\saturnws\data\200712\12

Dilution: 1.0 Initial Weight/Volume: 40 mL Date Analyzed: 12/03/2007 1605 Final Weight/Volume: 40 mL

Date Prepared: 12/03/2007 1605

Analyte	Result (ug/L)	Qualifier	RL
Methyl tert-butyl ether	ND		5.0
Acetone	ND		50
Benzene	2.7		0.50
Dichlorobromomethane	ND		0.50
Bromobenzene	ND		1.0
Chlorobromomethane	ND		1.0
Bromoform	ND		1.0
Bromomethane	ND		1.0
2-Butanone (MEK)	ND		50
n-Butylbenzene	ND		1.0
sec-Butylbenzene	ND		1.0
tert-Butylbenzene	ND		1.0
Carbon disulfide	ND		5.0
Carbon tetrachloride	ND		0.50
Chlorobenzene	ND		0.50
Chloroethane	ND		1.0
Chloroform	ND		1.0
Chloromethane	ND		1.0
2-Chlorotoluene	ND		0.50
4-Chlorotoluene	ND		0.50
Chlorodibromomethane	ND		0.50
1,2-Dichlorobenzene	ND		0.50
1,3-Dichlorobenzene	ND		0.50
1,4-Dichlorobenzene	ND		0.50
1,3-Dichloropropane	ND		1.0
1,1-Dichloropropene	ND		0.50
1,2-Dibromo-3-Chloropropane	ND		1.0
Ethylene Dibromide	ND		0.50
Dibromomethane	ND		0.50
Dichlorodifluoromethane	ND		0.50
1,1-Dichloroethane	ND		0.50
1,2-Dichloroethane	ND		0.50
1,1-Dichloroethene	ND		0.50
cis-1,2-Dichloroethene	ND		0.50
trans-1,2-Dichloroethene	ND		0.50
1,2-Dichloropropane	ND		0.50
cis-1,3-Dichloropropene	ND		0.50
trans-1,3-Dichloropropene	ND		0.50
Ethylbenzene	1.2		0.50
Hexachlorobutadiene	ND		1.0
2-Hexanone	ND		50
Isopropylbenzene	ND		0.50
4-Isopropyltoluene	ND		1.0
Methylene Chloride	ND		5.0

Client: Engeo, Inc. Job Number: 720-11961-1

Client Sample ID: DB-1 (GW@70')

 Lab Sample ID:
 720-11961-2
 Date Sampled:
 11/28/2007
 1340

 Client Matrix:
 Water
 Date Received:
 11/28/2007
 1810

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method: 8260B Analysis Batch: 720-29171 Instrument ID: Varian 3900G

Preparation: 5030B Lab File ID: c:\saturnws\data\200712\12

Dilution: 1.0 Initial Weight/Volume: 40 mL Date Analyzed: 12/03/2007 1605 Final Weight/Volume: 40 mL

Date Prepared: 12/03/2007 1605

Analyte	Result (ug/L)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		50
Naphthalene	ND		1.0
N-Propylbenzene	ND		1.0
Styrene	ND		0.50
1,1,1,2-Tetrachloroethane	ND		0.50
1,1,2,2-Tetrachloroethane	ND		0.50
Tetrachloroethene	ND		0.50
Toluene	2.3		0.50
1,2,3-Trichlorobenzene	ND		1.0
1,2,4-Trichlorobenzene	ND		1.0
1,1,1-Trichloroethane	ND		0.50
1,1,2-Trichloroethane	ND		0.50
Trichloroethene	ND		0.50
Trichlorofluoromethane	ND		1.0
1,2,3-Trichloropropane	ND		0.50
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		0.50
1,2,4-Trimethylbenzene	ND		0.50
1,3,5-Trimethylbenzene	ND		0.50
Vinyl acetate	ND		50
Vinyl chloride	ND		0.50
Xylenes, Total	1.0		1.0
2,2-Dichloropropane	ND		0.50
Surrogate	%Rec	Acceptance	Limits
4-Bromofluorobenzene	111	83 - 127	
1,2-Dichloroethane-d4 (Surr)	106	86 - 129	
Toluene-d8 (Surr)	103	82 - 126	

Client: Engeo, Inc. Job Number: 720-11961-1

Client Sample ID: DB-1 (GW@70')

 Lab Sample ID:
 720-11961-2
 Date Sampled:
 11/28/2007 1340

 Client Matrix:
 Water
 Date Received:
 11/28/2007 1810

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 720-29173 Instrument ID: Varian 3900A

Preparation: 5030B Lab File ID: c:\saturnws\data\200711\11

Dilution: 1.0 Initial Weight/Volume: 10 mL

Date Analyzed: 11/30/2007 1840 Final Weight/Volume: 10 mL

Date Prepared: 11/30/2007 1840

Analyte Result (ug/L) Qualifier RL
Gasoline Range Organics (GRO)-C5-C12 98 50

 Surrogate
 %Rec
 Acceptance Limits

 Toluene-d8 (Surr)
 99
 77 - 121

 1,2-Dichloroethane-d4 (Surr)
 101
 73 - 130

Client: Engeo, Inc. Job Number: 720-11961-1

Client Sample ID: DB-1 (GW@95')

 Lab Sample ID:
 720-11961-3
 Date Sampled:
 11/28/2007 1520

 Client Matrix:
 Water
 Date Received:
 11/28/2007 1810

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method: 8260B Analysis Batch: 720-29171 Instrument ID: Varian 3900G

Preparation: 5030B Lab File ID: c:\saturnws\data\200712\12

Dilution: 1.0 Initial Weight/Volume: 40 mL Date Analyzed: 12/03/2007 1638 Final Weight/Volume: 40 mL

Date Prepared: 12/03/2007 1638

Analyte	Result (ug/L)	Qualifier	RL
Methyl tert-butyl ether	ND		5.0
Acetone	ND		50
Benzene	ND		0.50
Dichlorobromomethane	ND		0.50
Bromobenzene	ND		1.0
Chlorobromomethane	ND		1.0
Bromoform	ND		1.0
Bromomethane	ND		1.0
2-Butanone (MEK)	ND		50
n-Butylbenzene	ND		1.0
sec-Butylbenzene	ND		1.0
tert-Butylbenzene	ND		1.0
Carbon disulfide	ND		5.0
Carbon tetrachloride	ND		0.50
Chlorobenzene	ND		0.50
Chloroethane	ND		1.0
Chloroform	ND		1.0
Chloromethane	ND		1.0
2-Chlorotoluene	ND		0.50
4-Chlorotoluene	ND		0.50
Chlorodibromomethane	ND		0.50
1,2-Dichlorobenzene	ND		0.50
1,3-Dichlorobenzene	ND		0.50
1,4-Dichlorobenzene	ND		0.50
1,3-Dichloropropane	ND		1.0
1,1-Dichloropropene	ND		0.50
1,2-Dibromo-3-Chloropropane	ND		1.0
Ethylene Dibromide	ND		0.50
Dibromomethane	ND		0.50
Dichlorodifluoromethane	ND		0.50
1,1-Dichloroethane	ND		0.50
1,2-Dichloroethane	ND		0.50
1,1-Dichloroethene	ND		0.50
cis-1,2-Dichloroethene	ND		0.50
trans-1,2-Dichloroethene	ND		0.50
1,2-Dichloropropane	ND		0.50
cis-1,3-Dichloropropene	ND		0.50
trans-1,3-Dichloropropene	ND		0.50
Ethylbenzene	ND		0.50
Hexachlorobutadiene	ND		1.0
2-Hexanone	ND		50
Isopropylbenzene	ND		0.50
4-Isopropyltoluene	ND		1.0
Methylene Chloride	ND		5.0

Client: Engeo, Inc. Job Number: 720-11961-1

Client Sample ID: DB-1 (GW@95')

 Lab Sample ID:
 720-11961-3
 Date Sampled:
 11/28/2007 1520

 Client Matrix:
 Water
 Date Received:
 11/28/2007 1810

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method: 8260B Analysis Batch: 720-29171 Instrument ID: Varian 3900G

Preparation: 5030B Lab File ID: c:\saturnws\data\200712\12

Dilution: 1.0 Initial Weight/Volume: 40 mL Date Analyzed: 12/03/2007 1638 Final Weight/Volume: 40 mL

Date Prepared: 12/03/2007 1638

Analyte	Result (ug/L)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		50
Naphthalene	ND		1.0
N-Propylbenzene	ND		1.0
Styrene	ND		0.50
1,1,1,2-Tetrachloroethane	ND		0.50
1,1,2,2-Tetrachloroethane	ND		0.50
Tetrachloroethene	ND		0.50
Toluene	ND		0.50
1,2,3-Trichlorobenzene	ND		1.0
1,2,4-Trichlorobenzene	ND		1.0
1,1,1-Trichloroethane	ND		0.50
1,1,2-Trichloroethane	ND		0.50
Trichloroethene	ND		0.50
Trichlorofluoromethane	ND		1.0
1,2,3-Trichloropropane	ND		0.50
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		0.50
1,2,4-Trimethylbenzene	ND		0.50
1,3,5-Trimethylbenzene	ND		0.50
Vinyl acetate	ND		50
Vinyl chloride	ND		0.50
Xylenes, Total	ND		1.0
2,2-Dichloropropane	ND		0.50
Surrogate	%Rec	Acceptance	e Limits
4-Bromofluorobenzene	111	83 - 127	
1,2-Dichloroethane-d4 (Surr)	106	86 - 129	
Toluene-d8 (Surr)	101	82 - 126	

Client: Engeo, Inc. Job Number: 720-11961-1

Client Sample ID: DB-1 (GW@95')

 Lab Sample ID:
 720-11961-3
 Date Sampled:
 11/28/2007 1520

 Client Matrix:
 Water
 Date Received:
 11/28/2007 1810

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 720-29163 Instrument ID: Varian 3900E

Preparation: 5030B Lab File ID: c:\varianws\data\200711\11

Dilution: 1.0 Initial Weight/Volume: 10 mL

Date Analyzed: 12/01/2007 0338 Final Weight/Volume: 10 mL Date Prepared: 12/01/2007 0338

Analyte Result (ug/L) Qualifier RL
Gasoline Range Organics (GRO)-C5-C12 ND 50

 Surrogate
 %Rec
 Acceptance Limits

 Toluene-d8 (Surr)
 91
 77 - 121

 1,2-Dichloroethane-d4 (Surr)
 93
 73 - 130

Client: Engeo, Inc. Job Number: 720-11961-1

Client Sample ID: DB-1 (GW@35')

 Lab Sample ID:
 720-11961-1
 Date Sampled:
 11/28/2007 1000

 Client Matrix:
 Water
 Date Received:
 11/28/2007 1810

8015B Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)-Silica Gel Cleanup

Method: 8015B Analysis Batch: 720-29381 Instrument ID: Varian DRO4

Preparation: 3510C SGC Prep Batch: 720-29325 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 250 mL

Date Analyzed: 12/07/2007 1038 Final Weight/Volume: 1 mL

Date Prepared: 12/06/2007 1438 Injection Volume:

Column ID: PRIMARY

Analyte	Result (ug/L)	Qualifier	RL
Diesel Range Organics [C10-C28]	ND		50
Motor Oil Range Organics [C24-C36]	ND		500

 Surrogate
 %Rec
 Acceptance Limits

 Capric Acid (Surr)
 0
 0 - 5

 p-Terphenyl
 66
 50 - 130

Client: Engeo, Inc. Job Number: 720-11961-1

Client Sample ID: DB-1 (GW@70')

Lab Sample ID: 720-11961-2 Date Sampled: 11/28/2007 1340 Client Matrix: Water Date Received: 11/28/2007 1810

8015B Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)-Silica Gel Cleanup

8015B Analysis Batch: 720-29381 Varian DRO4 Method: Instrument ID:

Preparation: 3510C SGC Prep Batch: 720-29325 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume:

250 mL Date Analyzed: 12/07/2007 1105 Final Weight/Volume: 1 mL

Date Prepared: 12/06/2007 1438 Injection Volume:

Column ID: **PRIMARY**

Analyte	Result (ug/L)	Qualifier	RL
Diesel Range Organics [C10-C28]	190		50
Motor Oil Range Organics [C24-C36]	ND		500

Surrogate	%Rec	Acceptance Limits
Capric Acid (Surr)	0	0 - 5
p-Terphenyl	67	50 - 130

Client: Engeo, Inc. Job Number: 720-11961-1

Client Sample ID: DB-1 (GW@95')

 Lab Sample ID:
 720-11961-3
 Date Sampled:
 11/28/2007 1520

 Client Matrix:
 Water
 Date Received:
 11/28/2007 1810

8015B Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)-Silica Gel Cleanup

Method: 8015B Analysis Batch: 720-29381 Instrument ID: Varian DRO4

Preparation: 3510C SGC Prep Batch: 720-29325 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 250 mL

Date Analyzed: 12/07/2007 1131 Final Weight/Volume: 1 mL

Date Prepared: 12/06/2007 1438 Injection Volume:

Column ID: PRIMARY

Analyte	Result (ug/L)	Qualifier	RL
Diesel Range Organics [C10-C28]	ND		50
Motor Oil Range Organics [C24-C36]	ND		500

Surrogate	%Rec	Acceptance Limits
Capric Acid (Surr)	0	0 - 5
p-Terphenyl	60	50 - 130

DATA REPORTING QUALIFIERS

Client: Engeo, Inc. Job Number: 720-11961-1

Lab Section	Qualifier	Description
GC/MS VOA		
	X	Surrogate exceeds the control limits

Client: Engeo, Inc. Job Number: 720-11961-1

QC Association Summary

		Report			
Lab Sample ID	Client Sample ID	Basis	Client Matrix	Method	Prep Batch
GC/MS VOA					
Analysis Batch:720-2	9163				
LCS 720-29163/15	Lab Control Spike	T	Water	8260B	
LCSD 720-29163/14	Lab Control Spike Duplicate	T	Water	8260B	
MB 720-29163/9	Method Blank	T	Water	8260B	
720-11961-3	DB-1 (GW@95')	T	Water	8260B	
720-11962-B-8 MS	Matrix Spike	T	Water	8260B	
720-11962-B-8 MSD	Matrix Spike Duplicate	T	Water	8260B	
Analysis Batch:720-2	9171				
LCS 720-29171/2	Lab Control Spike	T	Water	8260B	
LCSD 720-29171/1	Lab Control Spike Duplicate	Т	Water	8260B	
MB 720-29171/3	Method Blank	Т	Water	8260B	
720-11935-C-7 MS	Matrix Spike	T	Water	8260B	
720-11935-D-7 MSD	Matrix Spike Duplicate	Т	Water	8260B	
720-11961-1	DB-1 (GW@35')	T	Water	8260B	
720-11961-2	DB-1 (GW@70')	Т	Water	8260B	
720-11961-3	DB-1 (GW@95')	T	Water	8260B	
Analysis Batch:720-2	9173				
LCS 720-29173/3	Lab Control Spike	T	Water	8260B	
MB 720-29173/4	Method Blank	T	Water	8260B	
720-11931-A-1 MS	Matrix Spike	T	Water	8260B	
720-11931-A-1 MSD	Matrix Spike Duplicate	T	Water	8260B	
720-11961-1	DB-1 (GW@35')	T	Water	8260B	
720-11961-2	DB-1 (GW@70')	T	Water	8260B	

Report Basis

T = Total

Client: Engeo, Inc. Job Number: 720-11961-1

QC Association Summary

		Report			
Lab Sample ID	Client Sample ID	Basis	Client Matrix	Method	Prep Batch
GC Semi VOA					
Prep Batch: 720-29325					
LCS 720-29325/2-A	Lab Control Spike	Α	Water	3510C SGC	
LCSD 720-29325/3-A	Lab Control Spike Duplicate	Α	Water	3510C SGC	
MB 720-29325/1-A	Method Blank	Α	Water	3510C SGC	
720-11961-1	DB-1 (GW@35')	Α	Water	3510C SGC	
720-11961-2	DB-1 (GW@70')	Α	Water	3510C SGC	
720-11961-3	DB-1 (GW@95')	Α	Water	3510C SGC	
720-12002-E-2-B MS	Matrix Spike	Α	Water	3510C SGC	
720-12002-E-2-C MSD	Matrix Spike Duplicate	Α	Water	3510C SGC	
Analysis Batch:720-293	881				
LCS 720-29325/2-A	Lab Control Spike	Α	Water	8015B	720-29325
LCSD 720-29325/3-A	Lab Control Spike Duplicate	Α	Water	8015B	720-29325
MB 720-29325/1-A	Method Blank	Α	Water	8015B	720-29325
720-11961-1	DB-1 (GW@35')	Α	Water	8015B	720-29325
720-11961-2	DB-1 (GW@70')	Α	Water	8015B	720-29325
720-11961-3	DB-1 (GW@95')	Α	Water	8015B	720-29325
Analysis Batch:720-293	382				
720-12002-E-2-B MS	Matrix Spike	Α	Water	8015B	720-29325
720-12002-E-2-C MSD	Matrix Spike Duplicate	Α	Water	8015B	720-29325

Report Basis
A = Silica Gel Cleanup

Client: Engeo, Inc. Job Number: 720-11961-1

Method Blank - Batch: 720-29163 Method: 8260B Preparation: 5030B

Lab Sample ID: MB 720-29163/9 Analysis Batch: 720-29163 Instrument ID: Varian 3900E

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\varianws\data\200711\11

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 10 mL Date Analyzed: 11/30/2007 2141 Final Weight/Volume: 10 mL Date Prepared: 11/30/2007 2141

Analyte	Result	Qual	RL
Benzene	ND		0.50
Toluene	ND		0.50
Gasoline Range Organics (GRO)-C5-C12	ND		50
Surrogate	% Rec	Acceptance	Limits
Toluene-d8 (Surr)	97	77 - 121	
1,2-Dichloroethane-d4 (Surr)	90	73 - 130)

Lab Control Spike/ Method: 8260B
Lab Control Spike Duplicate Recovery Report - Batch: 720-29163 Preparation: 5030B

LCS Lab Sample ID: LCS 720-29163/15 Analysis Batch: 720-29163 Instrument ID: Varian 3900E

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\varianws\\data\200711\11

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 10 mL

Date Analyzed: 11/30/2007 2131 Final Weight/Volume: 10 mL

Date Analyzed: 11/30/2007 2131 Final Weight/Volume: 10 mL Date Prepared: 11/30/2007 2131

LCSD Lab Sample ID: LCSD 720-29163/14 Analysis Batch: 720-29163 Instrument ID: Varian 3900E

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\varianws\data\200711\113
Dilution: 1.0 Units: ug/L Initial Weight/Volume: 10 mL

Date Analyzed: 11/30/2007 2155 Final Weight/Volume: 10 mL

Date Prepared: 11/30/2007 2155

% Rec. RPD RPD Limit LCS Qual LCSD Qual Analyte LCS **LCSD** Limit Benzene 90 97 69 - 129 8 20 Toluene 95 100 70 - 130 5 20 LCS % Rec LCSD % Rec Surrogate Acceptance Limits 99 77 - 121 Toluene-d8 (Surr) 96 1,2-Dichloroethane-d4 (Surr) 88 73 - 130 90

Client: Engeo, Inc. Job Number: 720-11961-1

Matrix Spike/ Method: 8260B
Matrix Spike Duplicate Recovery Report - Batch: 720-29163 Preparation: 5030B

MS Lab Sample ID: 720-11962-B-8 MS Analysis Batch: 720-29163 Instrument ID: Varian 3900E

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\varianws\data\200711\'

Dilution: 5.0 Initial Weight/Volume: 10 mL

Date Analyzed: 12/01/2007 0329 Final Weight/Volume: 10 mL

Date Prepared: 12/01/2007 0329 Final Weight/Volume: 10 mL

MSD Lab Sample ID: 720-11962-B-8 MSD Analysis Batch: 720-29163 Instrument ID: Varian 3900E

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\varianws\data\200711\11

Dilution: 5.0 Initial Weight/Volume: 10 mL

Date Analyzed: 12/01/2007 0352 Final Weight/Volume: 10 mL Date Prepared: 12/01/2007 0352

% Rec. MS MSD Limit RPD **RPD** Limit MS Qual MSD Qual Analyte Benzene 83 91 69 - 129 9 20 Toluene 91 93 70 - 130 2 20 Surrogate MS % Rec MSD % Rec Acceptance Limits Toluene-d8 (Surr) 92 90 77 - 121 1,2-Dichloroethane-d4 (Surr) 80 83 73 - 130

Client: Engeo, Inc. Job Number: 720-11961-1

Method Blank - Batch: 720-29171 Method: 8260B Preparation: 5030B

Lab Sample ID: MB 720-29171/3 Analysis Batch: 720-29171 Instrument ID: Varian 3900G

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\saturnws\data\200712\12

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 40 mL Date Analyzed: 12/03/2007 1104 Final Weight/Volume: 40 mL

Date Prepared: 12/03/2007 1104

Analyte	Result	Qual	RL
Methyl tert-butyl ether	ND		5.0
Acetone	ND		50
Benzene	ND		0.50
Dichlorobromomethane	ND		0.50
Bromobenzene	ND		1.0
Chlorobromomethane	ND		1.0
Bromoform	ND		1.0
Bromomethane	ND		1.0
2-Butanone (MEK)	ND		50
n-Butylbenzene	ND		1.0
sec-Butylbenzene	ND		1.0
tert-Butylbenzene	ND		1.0
Carbon disulfide	ND		5.0
Carbon tetrachloride	ND		0.50
Chlorobenzene	ND		0.50
Chloroethane	ND		1.0
Chloroform	ND		1.0
Chloromethane	ND		1.0
2-Chlorotoluene	ND		0.50
4-Chlorotoluene	ND		0.50
Chlorodibromomethane	ND		0.50
1,2-Dichlorobenzene	ND		0.50
1,3-Dichlorobenzene	ND		0.50
1,4-Dichlorobenzene	ND		0.50
1,3-Dichloropropane	ND		1.0
1,1-Dichloropropene	ND		0.50
1,2-Dibromo-3-Chloropropane	ND		1.0
Ethylene Dibromide	ND		0.50
Dibromomethane	ND		0.50
Dichlorodifluoromethane	ND		0.50
1,1-Dichloroethane	ND		0.50
1,2-Dichloroethane	ND		0.50
1,1-Dichloroethene	ND		0.50
cis-1,2-Dichloroethene	ND		0.50
trans-1,2-Dichloroethene	ND		0.50
1,2-Dichloropropane	ND		0.50
cis-1,3-Dichloropropene	ND		0.50
trans-1,3-Dichloropropene	ND		0.50
Ethylbenzene	ND		0.50
Hexachlorobutadiene	ND		1.0
2-Hexanone	ND		50

Client: Engeo, Inc. Job Number: 720-11961-1

Method Blank - Batch: 720-29171 Method: 8260B Preparation: 5030B

Lab Sample ID: MB 720-29171/3 Analysis Batch: 720-29171 Instrument ID: Varian 3900G

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\saturnws\data\200712\12

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 40 mL Date Analyzed: 12/03/2007 1104 Final Weight/Volume: 40 mL

Date Prepared: 12/03/2007 1104

4-Isopropyltoluene Methylene Chloride A-Methyl-2-pentanone (MIBK) Naphthalene N-Propylbenzene Styrene 1,1,1,2-Tetrachloroethane	ID I		0.50 1.0 5.0 50 1.0 1.0 0.50 0.50 0.50
Methylene Chloride 1-Methyl-2-pentanone (MIBK) Naphthalene N-Propylbenzene Styrene 1,1,1,2-Tetrachloroethane	ID ID ID ID ID ID ID ID		5.0 50 1.0 1.0 0.50 0.50 0.50 0.50
4-Methyl-2-pentanone (MIBK) Naphthalene N-Propylbenzene Styrene N,1,1,2-Tetrachloroethane	ID ID ID ID ID ID ID ID		50 1.0 1.0 0.50 0.50 0.50 0.50
Naphthalene N-Propylbenzene NStyrene N1,1,1,2-Tetrachloroethane NST	ID ID ID ID ID ID ID		1.0 1.0 0.50 0.50 0.50 0.50
N-Propylbenzene Styrene N,1,1,2-Tetrachloroethane	ID ID ID ID ID ID		1.0 0.50 0.50 0.50 0.50
Styrene N 1,1,1,2-Tetrachloroethane N	ID ID ID ID ID		0.50 0.50 0.50 0.50
1,1,1,2-Tetrachloroethane	ID ID ID ID		0.50 0.50 0.50
1,1,1,2-Tetrachloroethane	ID ID ID		0.50 0.50
1.1.2.2-Tetrachloroethane	ID ID		0.50
	ID		
Fetrachloroethene N			
Foluene N	ID		0.50
1,2,3-Trichlorobenzene	10		1.0
1,2,4-Trichlorobenzene	ID		1.0
1,1,1-Trichloroethane	ID		0.50
1,1,2-Trichloroethane	ID		0.50
Frichloroethene N	ID		0.50
Frichlorofluoromethane N	ID		1.0
1,2,3-Trichloropropane N	ID		0.50
1,1,2-Trichloro-1,2,2-trifluoroethane	ID		0.50
1,2,4-Trimethylbenzene N	ID		0.50
1,3,5-Trimethylbenzene	ID		0.50
/inyl acetate N	ID		50
/inyl chloride N	ID		0.50
Kylenes, Total N	ID		1.0
2,2-Dichloropropane	ID		0.50
Surrogate	% Rec	Acceptance Limits	
4-Bromofluorobenzene	107	83 - 127	
1,2-Dichloroethane-d4 (Surr)	99	86 - 129	
Foluene-d8 (Surr)	101	82 - 126	

Client: Engeo, Inc. Job Number: 720-11961-1

Lab Control Spike/ Method: 8260B
Lab Control Spike Duplicate Recovery Report - Batch: 720-29171 Preparation: 5030B

LCS Lab Sample ID: LCS 720-29171/2 Analysis Batch: 720-29171 Instrument ID: Varian 3900G

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\saturnws\data\200712\12

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 40 mL

Date Analyzed: 12/03/2007 0958 Final Weight/Volume: 40 mL Date Prepared: 12/03/2007 0958

LCSD Lab Sample ID: LCSD 720-29171/1 Analysis Batch: 720-29171 Instrument ID: Varian 3900G

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\saturnws\data\200712\12\

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 40 mL Date Analyzed: 12/03/2007 1031 Final Weight/Volume: 40 mL

Date Prepared:

12/03/2007 1031

% Rec. LCS **LCSD RPD** RPD Limit LCS Qual LCSD Qual Analyte Limit Benzene 85 82 69 - 129 4 20 99 96 61 - 121 3 20 Chlorobenzene 1,1-Dichloroethene 87 90 65 - 125 3 20 Toluene 90 87 70 - 130 20 4 Trichloroethene 74 - 134 20 85 81 5 LCS % Rec LCSD % Rec Acceptance Limits

 Trichloroethene
 85
 81
 74 - 134
 5
 20

 Surrogate
 LCS % Rec
 LCSD % Rec
 Acceptance Limits

 4-Bromofluorobenzene
 99
 103
 83 - 127

 1,2-Dichloroethane-d4 (Surr)
 91
 98
 86 - 129

 Toluene-d8 (Surr)
 92
 96
 82 - 126

Client: Engeo, Inc. Job Number: 720-11961-1

Matrix Spike/ Method: 8260B
Matrix Spike Duplicate Recovery Report - Batch: 720-29171 Preparation: 5030B

MS Lab Sample ID: 720-11935-C-7 MS Analysis Batch: 720-29171 Instrument ID: Varian 3900G

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\saturnws\data\200712\

Dilution: 1.0 Initial Weight/Volume: 40 mL Date Analyzed: 12/03/2007 1425 Final Weight/Volume: 40 mL

Date Prepared: 12/03/2007 1425

MSD Lab Sample ID: 720-11935-D-7 MSD Analysis Batch: 720-29171 Instrument ID: Varian 3900G

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\saturnws\data\200712\12

Dilution: 1.0 Initial Weight/Volume: 40 mL

Date Analyzed: 12/03/2007 1458 Final Weight/Volume: 40 mL Date Prepared: 12/03/2007 1458

	<u>%</u>	Rec.				
Analyte	MS	MSD	Limit	RPD	RPD Limit	MS Qual MSD Qual
Benzene	92	91	69 - 129	2	20	
Chlorobenzene	110	110	61 - 121	0	20	
1,1-Dichloroethene	97	97	65 - 125	0	20	
Toluene	99	97	70 - 130	2	20	
Trichloroethene	100	92	74 - 134	9	20	
Surrogate		MS % Rec	MSD % Rec		Acceptance Limits	
4-Bromofluorobenzene		109	103		83	3 - 127
1,2-Dichloroethane-d4 (Surr)		104	105		86	6 - 129
Toluene-d8 (Surr)		107	103		82	2 - 126

Client: Engeo, Inc. Job Number: 720-11961-1

Method Blank - Batch: 720-29173 Method: 8260B Preparation: 5030B

Lab Sample ID: MB 720-29173/4 Analysis Batch: 720-29173 Instrument ID: Varian 3900A

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\saturnws\data\200711\11

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 10 mL

Date Analyzed: 11/30/2007 1100 Final Weight/Volume: 10 mL Date Prepared: 11/30/2007 1100

Analyte	Result	Qual	RL
Benzene	ND		0.50
Toluene	ND		0.50
Gasoline Range Organics (GRO)-C5-C12	ND		50
Surrogate	% Rec	Acceptano	ce Limits
Toluene-d8 (Surr)	98	77 - 1	21
1,2-Dichloroethane-d4 (Surr)	94	73 - 1	30

Lab Control Spike - Batch: 720-29173 Method: 8260B Preparation: 5030B

Lab Sample ID: LCS 720-29173/3 Analysis Batch: 720-29173 Instrument ID: Varian 3900A

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\saturnws\data\200711\11

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 10 mL Date Analyzed: 11/30/2007 0937 Final Weight/Volume: 10 mL

Date Analyzed: 11/30/2007 0937 Final Weight/Volume: 10 mL

Date Prepared: 11/30/2007 0937

Analyte	Spike Amount	Result	% Rec.	Limit	Qual
Benzene	5.40	5.19	96	69 - 129	
Toluene	40.3	38.7	96	70 - 130	
Surrogate	% Rec		Acceptance Limits		
Toluene-d8 (Surr)	99			77 - 121	
1,2-Dichloroethane-d4 (Surr)	109		73 - 130		

Client: Engeo, Inc. Job Number: 720-11961-1

Matrix Spike/ Method: 8260B
Matrix Spike Duplicate Recovery Report - Batch: 720-29173 Preparation: 5030B

MS Lab Sample ID: 720-11931-A-1 MS Analysis Batch: 720-29173 Instrument ID: Varian 3900A

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\saturnws\data\200711\

Dilution: 1.0 Initial Weight/Volume: 10 mL

Date Analyzed: 11/30/2007 1242 Final Weight/Volume: 10 ml

Date Analyzed: 11/30/2007 1242 Final Weight/Volume: 10 mL Date Prepared: 11/30/2007 1242

MSD Lab Sample ID: 720-11931-A-1 MSD Analysis Batch: 720-29173 Instrument ID: Varian 3900A

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\saturnws\data\200711\11

Dilution: 1.0 Initial Weight/Volume: 10 mL

Date Analyzed: 11/30/2007 1305 Final Weight/Volume: 10 mL Date Prepared: 11/30/2007 1305

% Rec. MS MSD Limit RPD **RPD** Limit MS Qual MSD Qual Analyte Benzene 94 95 69 - 129 20 1 Toluene 102 101 70 - 130 0 20 Surrogate MS % Rec MSD % Rec Acceptance Limits Toluene-d8 (Surr) 100 102 77 - 121 1,2-Dichloroethane-d4 (Surr) 100 59 Χ 73 - 130

Client: Engeo, Inc. Job Number: 720-11961-1

Method Blank - Batch: 720-29325 Method: 8015B

Preparation: 3510C SGC Silica Gel Cleanup

Lab Sample ID: MB 720-29325/1-A

Client Matrix: Water Dilution: 1.0

Date Analyzed: 12/07/2007 1012 Date Prepared: 12/06/2007 1438 Analysis Batch: 720-29381 Prep Batch: 720-29325

Units: ug/L

Instrument ID: Varian DRO4

Lab File ID: N/A

Initial Weight/Volume: 250 mL Final Weight/Volume: 1 mL

Injection Volume:

Column ID: PRIMARY

Analyte	Result	Qual	RL
Diesel Range Organics [C10-C28]	ND		50
Motor Oil Range Organics [C24-C36]	ND		500
Surrogate	% Rec	Acceptance Limits	
Capric Acid (Surr)	0	0 - 5	
p-Terphenyl	72	50 - 130	

Lab Control Spike/ Method: 8015B

Lab Control Spike Duplicate Recovery Report - Batch: 720-29325 Preparation: 3510C SGC

Silica Gel Cleanup

LCS Lab Sample ID: LCS 720-29325/2-A Analysis Batch: 720-29381 Instrument ID: Varian DRO4

Client Matrix: Water

Dilution: 1.0 Date Analyzed: 12/07

Date Analyzed: 12/07/2007 0920 Date Prepared: 12/06/2007 1438 Analysis Batch: 720-29381 Prep Batch: 720-29325

Units: ug/L

720-29325 Lab File ID: N/A

Initial Weight/Volume: 250 mL Final Weight/Volume: 1 mL

Injection Volume:

Column ID: PRIMARY

LCSD Lab Sample ID: LCSD 720-29325/3-A

Client Matrix: Water Dilution: 1.0

Date Analyzed: 12/07/2007 0946 Date Prepared: 12/06/2007 1438 Analysis Batch: 720-29381

Prep Batch: 720-29325

Units: ug/L

Instrument ID: Varian DRO4

Lab File ID: N/A

Initial Weight/Volume: 250 mL Final Weight/Volume: 1 mL

Injection Volume:

Column ID: PRIMARY

% Rec. Analyte LCS **LCSD** Limit **RPD** RPD Limit LCS Qual LCSD Qual Diesel Range Organics [C10-C28] 58 62 50 - 130 7 30 LCS % Rec Surrogate LCSD % Rec Acceptance Limits 71 73 50 - 130 p-Terphenyl

PRIMARY

50 - 130

Job Number: 720-11961-1 Client: Engeo, Inc.

Matrix Spike/ Method: 8015B

Matrix Spike Duplicate Recovery Report - Batch: 720-29325 Preparation: 3510C SGC

Silica Gel Cleanup

Instrument ID: HP DRO5 MS Lab Sample ID: 720-12002-E-2-B MS Analysis Batch: 720-29382

Client Matrix: Prep Batch: 720-29325 Water Lab File ID: N/A Dilution: 1.0

Initial Weight/Volume: 250 mL Date Analyzed: 12/07/2007 1208 Final Weight/Volume: 1 mL

Date Prepared: 12/06/2007 1438 Injection Volume: Column ID:

MSD Lab Sample ID: 720-12002-E-2-C MSD Instrument ID: HP DRO5 Analysis Batch: 720-29382

Client Matrix: Water Prep Batch: 720-29325 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 250 mL Date Analyzed: 12/07/2007 1235 Final Weight/Volume: 1 mL

Date Prepared: 12/06/2007 1438 Injection Volume:

Column ID: **PRIMARY**

66

p-Terphenyl

% Rec. MS MSD Limit **RPD** MS Qual MSD Qual Analyte **RPD Limit** Diesel Range Organics [C10-C28] 50 - 130 56 64 12 30 Surrogate MS % Rec MSD % Rec Acceptance Limits

70

Login Sample Receipt Check List

Client: Engeo, Inc.

Job Number: 720-11961-1

Login Number: 11961 List Source: TestAmerica San Francisco

Creator: Lewis, Trenton L List Number: 1

Question	T / F/ NA	Comment
Radioactivity either was not measured or, if measured, is at or below background	N/A	
The cooler's custody seal, if present, is intact.	N/A	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
There are no discrepancies between the sample IDs on the containers and the COC.	False	See Narrative
Samples are received within Holding Time.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
VOA sample vials do not have headspace or bubble is <6mm (1/4") in diameter.	True	
If necessary, staff have been informed of any short hold time or quick TAT needs	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	



ANALYTICAL REPORT

Job Number: 720-12284-1

Job Description: Rickenbacker

For:
Engeo, Inc.
580 N Wilma Avenue
Suite A
Ripon, CA 95366-9502

Attention: Mr. Richard Gandolfo

Akanef Sal

Designee for
Melissa Brewer
Project Manager I
melissa.brewer@testamericainc.com
01/04/2008

Job Narrative 720-J12284-1

Comments

No additional comments.

Receipt

All samples were received in good condition within temperature requirements.

GC/MS VOA

Method(s) 8260B: The matrix spike duplicate (MS/MSD) recoveries for batch 30082 were outside control limits. The associated laboratory control standard (LCS) met acceptance criteria.

Method(s) 8260B: Surrogate and internal standard recoveries for the following sample was outside control limits: B-8-2.5 (720-12227-1). Evidence of matrix interference is present; confirmed by MS and MSD.

Method(s) 8260B: The matrix spike and matrix spike duplicate (MS/MSD) recoveries for batch 29888 were outside control limits. The associated laboratory control standard and laboratory control standard duplicate (LCS/LCSD) met acceptance criteria.

No other analytical or quality issues were noted.

GC VOA

No analytical or quality issues were noted.

GC Semi VOA

Method(s) 8015B: The continuing calibration verification (CCV) for PTP recovered above the upper control limit. The samples associated with this CCV were non-detects for the affected analytes; therefore, the data have been reported.

No other analytical or quality issues were noted.

Organic Prep

No analytical or quality issues were noted.

EXECUTIVE SUMMARY - Detections

Client: Engeo, Inc. Job Number: 720-12284-1

Lab Sample ID Analyte	Client Sample ID	Result / Qualifier	Reporting Limit	Units	Method	
720-12284-1	MWB1 @ 5-1/2'					
Tetrachloroethene		81	4.7	ug/Kg	8260B	
720-12284-2	MWB1 @ 10-1/2'					
Tetrachloroethene		68	4.9	ug/Kg	8260B	

METHOD SUMMARY

Client: Engeo, Inc. Job Number: 720-12284-1

Description	Lab Location	Method	Preparation Method
Matrix: Solid			
Volatile Organic Compounds by GC/MS	TAL SF	SW846 8260B	
Volatile Organic Compounds by GC/MS (Low Level)	TAL SF	SW846 8260B	
Purge and Trap for Solids	TAL SF		SW846 5030B
Purge and Trap for Solids	TAL SF		SW846 5030B
Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)	TAL SF	SW846 8015B	
Ultrasonic Extraction	TAL SF		SW846 3550B

Lab References:

TAL SF = TestAmerica San Francisco

Method References:

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

METHOD / ANALYST SUMMARY

Client: Engeo, Inc. Job Number: 720-12284-1

Method	Analyst	Analyst ID
SW846 8260B SW846 8260B	Ali, Badri Le, Lien	BA LL
SW846 8015B	Hayashi, Derek	DH

SAMPLE SUMMARY

Client: Engeo, Inc. Job Number: 720-12284-1

Lab Sample ID	Client Sample ID	Client Matrix	Date/Time Sampled	Date/Time Received
720-12284-1	MWB1 @ 5-1/2'	Solid	12/18/2007 0900	12/18/2007 1532
720-12284-2	MWB1 @ 10-1/2'	Solid	12/18/2007 0910	12/18/2007 1532
720-12284-3	MWB2 @ 25-1/2'	Solid	12/18/2007 1330	12/18/2007 1532

Client: Engeo, Inc. Job Number: 720-12284-1

Client Sample ID: MWB1 @ 5-1/2'

 Lab Sample ID:
 720-12284-1
 Date Sampled:
 12/18/2007 0900

 Client Matrix:
 Solid
 Date Received:
 12/18/2007 1532

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29888Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29834Lab File ID:121907012.DDilution:1.0Initial Weight/Volume:5.27 g

Date Analyzed: 12/19/2007 1541 Final Weight/Volume: 5.27 g

Date Prepared: 12/19/2007 1200

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
Methyl tert-butyl ether	ND		4.7
Acetone	ND		47
Benzene	ND		4.7
Dichlorobromomethane	ND		4.7
Bromobenzene	ND		4.7
Chlorobromomethane	ND		19
Bromoform	ND		4.7
Bromomethane	ND		9.5
2-Butanone (MEK)	ND		47
n-Butylbenzene	ND		4.7
sec-Butylbenzene	ND		4.7
tert-Butylbenzene	ND		4.7
Carbon disulfide	ND		4.7
Carbon tetrachloride	ND		4.7
Chlorobenzene	ND		4.7
Chloroethane	ND		9.5
Chloroform	ND		4.7
Chloromethane	ND		9.5
2-Chlorotoluene	ND		4.7
4-Chlorotoluene	ND		4.7
Chlorodibromomethane	ND		4.7
1,2-Dichlorobenzene	ND		4.7
1,3-Dichlorobenzene	ND		4.7
1,4-Dichlorobenzene	ND		4.7
1,3-Dichloropropane	ND		4.7
1,1-Dichloropropene	ND		4.7
1,2-Dibromo-3-Chloropropane	ND		47
Ethylene Dibromide	ND		4.7
Dibromomethane	ND		9.5
Dichlorodifluoromethane	ND		9.5
1,1-Dichloroethane	ND		4.7
1,2-Dichloroethane	ND		4.7
1,1-Dichloroethene	ND		4.7
cis-1,2-Dichloroethene	ND		4.7
trans-1,2-Dichloroethene	ND		4.7
1,2-Dichloropropane	ND		4.7
cis-1,3-Dichloropropene	ND		4.7
trans-1,3-Dichloropropene	ND		4.7
Ethylbenzene	ND		4.7
Hexachlorobutadiene	ND		4.7
2-Hexanone	ND		47
Isopropylbenzene	ND		4.7
4-Isopropyltoluene	ND		4.7
Methylene Chloride	ND		9.5

Client: Engeo, Inc. Job Number: 720-12284-1

Client Sample ID: MWB1 @ 5-1/2'

 Lab Sample ID:
 720-12284-1
 Date Sampled:
 12/18/2007 0900

 Client Matrix:
 Solid
 Date Received:
 12/18/2007 1532

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29888Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29834Lab File ID:121907012.DDilution:1.0Initial Weight/Volume:5.27 g

Dilution: 1.0 Initial Weight/Volume: 5.27 g

Date Analyzed: 12/19/2007 1541 Final Weight/Volume: 10 mL

Date Prepared: 12/19/2007 1200

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		47
Naphthalene	ND		9.5
N-Propylbenzene	ND		4.7
Styrene	ND		4.7
1,1,1,2-Tetrachloroethane	ND		4.7
1,1,2,2-Tetrachloroethane	ND		4.7
Tetrachloroethene	81		4.7
Toluene	ND		4.7
1,2,3-Trichlorobenzene	ND		4.7
1,2,4-Trichlorobenzene	ND		4.7
1,1,1-Trichloroethane	ND		4.7
1,1,2-Trichloroethane	ND		4.7
Trichloroethene	ND		4.7
Trichlorofluoromethane	ND		4.7
1,2,3-Trichloropropane	ND		4.7
1,1,2-Trichloro-1,2,2-trifluoroethan	ne ND		4.7
1,2,4-Trimethylbenzene	ND		4.7
1,3,5-Trimethylbenzene	ND		4.7
Vinyl acetate	ND		47
Vinyl chloride	ND		4.7
Xylenes, Total	ND		9.5
2,2-Dichloropropane	ND		4.7
Surrogate	%Rec	Accep	tance Limits
4-Bromofluorobenzene	85	50 - 1	138
1,2-Dichloroethane-d4 (Surr)	94	66 - 1	127
Toluene-d8 (Surr)	89	51 - 1	129

Client: Engeo, Inc. Job Number: 720-12284-1

Client Sample ID: MWB1 @ 5-1/2'

 Lab Sample ID:
 720-12284-1
 Date Sampled:
 12/18/2007 0900

 Client Matrix:
 Solid
 Date Received:
 12/18/2007 1532

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 720-30082 Instrument ID: Varian 3900E

Preparation: 5030B Prep Batch: 720-30083 Lab File ID: c:\varianws\data\200712\12

Dilution: 1.0 Initial Weight/Volume: 5.50 g

Date Analyzed: 12/26/2007 1306 Final Weight/Volume: 10 mL

Date Prepared: 12/26/2007 0835

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL
Gasoline Range Organics (GRO)-C5-C12 ND 0.23

 Surrogate
 %Rec
 Acceptance Limits

 1,2-Dichloroethane-d4 (Surr)
 106
 60 - 140

 Toluene-d8 (Surr)
 97
 70 - 130

Client: Engeo, Inc. Job Number: 720-12284-1

Client Sample ID: MWB1 @ 10-1/2'

 Lab Sample ID:
 720-12284-2
 Date Sampled:
 12/18/2007 0910

 Client Matrix:
 Solid
 Date Received:
 12/18/2007 1532

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29888Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29834Lab File ID:121907013.DDilution:1.0Initial Weight/Volume:5.13 g

Dilution: 1.0 Initial Weight/Volume: 5.13 g
Date Analyzed: 12/19/2007 1606 Final Weight/Volume: 10 mL

Date Prepared: 12/19/2007 1200

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier RL
Methyl tert-butyl ether	ND	4.9
Acetone	ND	49
Benzene	ND	4.9
Dichlorobromomethane	ND	4.9
Bromobenzene	ND	4.9
Chlorobromomethane	ND	19
Bromoform	ND	4.9
Bromomethane	ND	9.7
2-Butanone (MEK)	ND	49
n-Butylbenzene	ND	4.9
sec-Butylbenzene	ND	4.9
tert-Butylbenzene	ND	4.9
Carbon disulfide	ND	4.9
Carbon tetrachloride	ND	4.9
Chlorobenzene	ND	4.9
Chloroethane	ND	9.7
Chloroform	ND	4.9
Chloromethane	ND	9.7
2-Chlorotoluene	ND	4.9
4-Chlorotoluene	ND	4.9
Chlorodibromomethane	ND	4.9
1,2-Dichlorobenzene	ND	4.9
1,3-Dichlorobenzene	ND	4.9
1,4-Dichlorobenzene	ND	4.9
1,3-Dichloropropane	ND	4.9
1,1-Dichloropropene	ND	4.9
1,2-Dibromo-3-Chloropropane	ND	49
Ethylene Dibromide	ND	4.9
Dibromomethane	ND	9.7
Dichlorodifluoromethane	ND	9.7
1,1-Dichloroethane	ND	4.9
1,2-Dichloroethane	ND	4.9
1,1-Dichloroethene	ND	4.9
cis-1,2-Dichloroethene	ND	4.9
trans-1,2-Dichloroethene	ND	4.9
1,2-Dichloropropane	ND	4.9
cis-1,3-Dichloropropene	ND	4.9
trans-1,3-Dichloropropene	ND	4.9
Ethylbenzene	ND	4.9
Hexachlorobutadiene	ND	4.9
2-Hexanone	ND	49
Isopropylbenzene	ND	4.9
4-Isopropyltoluene	ND	4.9
Methylene Chloride	ND	9.7
Modificial Official	IND	J.1

Client: Engeo, Inc. Job Number: 720-12284-1

Client Sample ID: MWB1 @ 10-1/2'

 Lab Sample ID:
 720-12284-2
 Date Sampled:
 12/18/2007 0910

 Client Matrix:
 Solid
 Date Received:
 12/18/2007 1532

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29888Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29834Lab File ID:121907013.DDilution:1.0Initial Weight/Volume:5.13 g

Date Analyzed: 12/19/2007 1606 Final Weight/Volume: 10 mL

Date Prepared: 12/19/2007 1200

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		49
Naphthalene	ND		9.7
N-Propylbenzene	ND		4.9
Styrene	ND		4.9
1,1,1,2-Tetrachloroethane	ND		4.9
1,1,2,2-Tetrachloroethane	ND		4.9
Tetrachloroethene	68		4.9
Toluene	ND		4.9
1,2,3-Trichlorobenzene	ND		4.9
1,2,4-Trichlorobenzene	ND		4.9
1,1,1-Trichloroethane	ND		4.9
1,1,2-Trichloroethane	ND		4.9
Trichloroethene	ND		4.9
Trichlorofluoromethane	ND		4.9
1,2,3-Trichloropropane	ND		4.9
1,1,2-Trichloro-1,2,2-trifluoroethan	ne ND		4.9
1,2,4-Trimethylbenzene	ND		4.9
1,3,5-Trimethylbenzene	ND		4.9
Vinyl acetate	ND		49
Vinyl chloride	ND		4.9
Xylenes, Total	ND		9.7
2,2-Dichloropropane	ND		4.9
Surrogate	%Rec	Λο	ceptance Limits
4-Bromofluorobenzene	85) - 138
			0 - 130 6 - 127
1,2-Dichloroethane-d4 (Surr)	95		
Toluene-d8 (Surr)	88	5	1 - 129

Client: Engeo, Inc. Job Number: 720-12284-1

Client Sample ID: MWB1 @ 10-1/2'

 Lab Sample ID:
 720-12284-2
 Date Sampled:
 12/18/2007 0910

 Client Matrix:
 Solid
 Date Received:
 12/18/2007 1532

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 720-30082 Instrument ID: Varian 3900E

Preparation: 5030B Prep Batch: 720-30083 Lab File ID: c:\varianws\data\200712\12

Dilution: 1.0 Initial Weight/Volume: 5.34 g

Date Analyzed: 12/26/2007 1330 Final Weight/Volume: 10 mL

Date Prepared: 12/26/2007 0835

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL
Gasoline Range Organics (GRO)-C5-C12 ND 0.23

 Surrogate
 %Rec
 Acceptance Limits

 1,2-Dichloroethane-d4 (Surr)
 101
 60 - 140

 Toluene-d8 (Surr)
 94
 70 - 130

Client: Engeo, Inc. Job Number: 720-12284-1

Client Sample ID: MWB2 @ 25-1/2'

 Lab Sample ID:
 720-12284-3
 Date Sampled:
 12/18/2007 1330

 Client Matrix:
 Solid
 Date Received:
 12/18/2007 1532

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29888Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29834Lab File ID:121907014.DDilution:1.0Initial Weight/Volume:5.00

Dilution: 1.0 Initial Weight/Volume: 5.00 g
Date Analyzed: 12/19/2007 1631 Final Weight/Volume: 10 mL

Date Prepared: 12/19/2007 1200

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier RL
Methyl tert-butyl ether	ND	5.0
Acetone	ND	50
Benzene	ND	5.0
Dichlorobromomethane	ND	5.0
Bromobenzene	ND	5.0
Chlorobromomethane	ND	20
Bromoform	ND	5.0
Bromomethane	ND	10
2-Butanone (MEK)	ND	50
n-Butylbenzene	ND	5.0
sec-Butylbenzene	ND	5.0
tert-Butylbenzene	ND	5.0
Carbon disulfide	ND	5.0
Carbon tetrachloride	ND	5.0
Chlorobenzene	ND	5.0
Chloroethane	ND	10
Chloroform	ND	5.0
Chloromethane	ND	10
2-Chlorotoluene	ND	5.0
4-Chlorotoluene	ND	5.0
Chlorodibromomethane	ND	5.0
1,2-Dichlorobenzene	ND	5.0
1,3-Dichlorobenzene	ND	5.0
1,4-Dichlorobenzene	ND	5.0
1,3-Dichloropropane	ND	5.0
1,1-Dichloropropene	ND	5.0
1,2-Dibromo-3-Chloropropane	ND	50
Ethylene Dibromide	ND	5.0
Dibromomethane	ND	10
Dichlorodifluoromethane	ND	10
1,1-Dichloroethane	ND	5.0
1,2-Dichloroethane	ND	5.0
1,1-Dichloroethene	ND	5.0
cis-1,2-Dichloroethene	ND	5.0
trans-1,2-Dichloroethene	ND	5.0
1,2-Dichloropropane	ND	5.0
cis-1,3-Dichloropropene	ND	5.0
trans-1,3-Dichloropropene	ND	5.0
Ethylbenzene	ND	5.0
Hexachlorobutadiene	ND	5.0
2-Hexanone	ND	50
Isopropylbenzene	ND	5.0
4-Isopropyltoluene	ND	5.0
Methylene Chloride	ND	10

Client: Engeo, Inc. Job Number: 720-12284-1

Client Sample ID: MWB2 @ 25-1/2'

 Lab Sample ID:
 720-12284-3
 Date Sampled:
 12/18/2007
 1330

 Client Matrix:
 Solid
 Date Received:
 12/18/2007
 1532

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29888Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29834Lab File ID:121907014.DDilution:1.0Initial Weight/Volume:5.00 g

Date Analyzed: 12/19/2007 1631 Final Weight/Volume: 10 mL

Date Prepared: 12/19/2007 1200

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		50
Naphthalene	ND		10
N-Propylbenzene	ND		5.0
Styrene	ND		5.0
1,1,1,2-Tetrachloroethane	ND		5.0
1,1,2,2-Tetrachloroethane	ND		5.0
Tetrachloroethene	ND		5.0
Toluene	ND		5.0
1,2,3-Trichlorobenzene	ND		5.0
1,2,4-Trichlorobenzene	ND		5.0
1,1,1-Trichloroethane	ND		5.0
1,1,2-Trichloroethane	ND		5.0
Trichloroethene	ND		5.0
Trichlorofluoromethane	ND		5.0
1,2,3-Trichloropropane	ND		5.0
1,1,2-Trichloro-1,2,2-trifluoroethar	ne ND		5.0
1,2,4-Trimethylbenzene	ND		5.0
1,3,5-Trimethylbenzene	ND		5.0
Vinyl acetate	ND		50
Vinyl chloride	ND		5.0
Xylenes, Total	ND		10
2,2-Dichloropropane	ND		5.0
Surrogate	%Rec	Acceptance	e Limits
4-Bromofluorobenzene	101	50 - 138	
1,2-Dichloroethane-d4 (Surr)	97	66 - 127	
Toluene-d8 (Surr)	101	51 - 129	

Client: Engeo, Inc. Job Number: 720-12284-1

Client Sample ID: MWB2 @ 25-1/2'

 Lab Sample ID:
 720-12284-3
 Date Sampled:
 12/18/2007 1330

 Client Matrix:
 Solid
 Date Received:
 12/18/2007 1532

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 720-30082 Instrument ID: Varian 3900E

Preparation: 5030B Prep Batch: 720-30083 Lab File ID: c:\varianws\data\200712\12

Dilution: 1.0 Initial Weight/Volume: 5.21 g

Date Analyzed: 12/26/2007 1353 Final Weight/Volume: 10 mL Date Prepared: 12/26/2007 0835

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL
Gasoline Range Organics (GRO)-C5-C12 ND 0.24

 Surrogate
 %Rec
 Acceptance Limits

 1,2-Dichloroethane-d4 (Surr)
 96
 60 - 140

 Toluene-d8 (Surr)
 95
 70 - 130

Client: Engeo, Inc. Job Number: 720-12284-1

Client Sample ID: MWB1 @ 5-1/2'

 Lab Sample ID:
 720-12284-1
 Date Sampled:
 12/18/2007 0900

 Client Matrix:
 Solid
 Date Received:
 12/18/2007 1532

8015B Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)-Silica Gel Cleanup

Method: 8015B Analysis Batch: 720-30339 Instrument ID: HP DRO5

Preparation: 3550B Prep Batch: 720-30218 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 30.15 g
Date Analyzed: 01/02/2008 1256 Final Weight/Volume: 5 mL

Date Prepared: 12/28/2007 1556 Injection Volume:

Column ID: PRIMARY

Analyte	DryWt Corrected: N Result (mg/Kg)	Qualifier	RL
Diesel Range Organics [C10-C28] ND		1.0
Motor Oil Range Organics [C24-0	[36] ND		50

Surrogate	%Rec	Acceptance Limits
Capric Acid (Surr)	5	0 - 5
p-Terphenyl	72	41 - 105

Client: Engeo, Inc. Job Number: 720-12284-1

Client Sample ID: MWB1 @ 10-1/2'

 Lab Sample ID:
 720-12284-2
 Date Sampled:
 12/18/2007 0910

 Client Matrix:
 Solid
 Date Received:
 12/18/2007 1532

8015B Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)-Silica Gel Cleanup

Method: 8015B Analysis Batch: 720-30339 Instrument ID: HP DRO5

Preparation: 3550B Prep Batch: 720-30218 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 30.07 g
Date Analyzed: 01/02/2008 1323 Final Weight/Volume: 5 mL

Date Prepared: 12/28/2007 1556 Injection Volume:

Column ID: PRIMARY

Analyte	DryWt Corrected: N Result (mg/Kg)	Qualifier	RL
Diesel Range Organics [C10-C28] ND		1.0
Motor Oil Range Organics [C24-0	[36] ND		50

Surrogate	%Rec	Acceptance Limits
Capric Acid (Surr)	3	0 - 5
p-Terphenyl	71	41 - 105

Client: Engeo, Inc. Job Number: 720-12284-1

Client Sample ID: MWB2 @ 25-1/2'

 Lab Sample ID:
 720-12284-3
 Date Sampled:
 12/18/2007 1330

 Client Matrix:
 Solid
 Date Received:
 12/18/2007 1532

8015B Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)-Silica Gel Cleanup

Method: 8015B Analysis Batch: 720-30339 Instrument ID: HP DRO5

Preparation: 3550B Prep Batch: 720-30218 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 30.03 g
Date Analyzed: 01/02/2008 2034 Final Weight/Volume: 5 mL

Date Prepared: 12/28/2007 1556 Injection Volume:

Column ID: PRIMARY

Analyte	DryWt Corrected: N Result (mg/Kg)	Qualifier	RL
Diesel Range Organics [C10-C	28] ND		1.0
Motor Oil Range Organics [C24	I-C361 ND		50

Surrogate	%Rec	Acceptance Limits
Capric Acid (Surr)	0	0 - 5
p-Terphenyl	87	41 - 105

DATA REPORTING QUALIFIERS

Client: Engeo, Inc. Job Number: 720-12284-1

Lab Section	Qualifier	Description
00/140 \/ (0.4		
GC/MS VOA		
	F	MS or MSD exceeds the control limits
	X	Surrogate exceeds the control limits

Client: Engeo, Inc. Job Number: 720-12284-1

QC Association Summary

		Report			
Lab Sample ID	Client Sample ID	Basis	Client Matrix	Method	Prep Batch
GC/MS VOA					
Prep Batch: 720-29834					
LCS 720-29834/1-A	Lab Control Spike	T	Solid	5030B	
LCSD 720-29834/2-A	Lab Control Spike Duplicate	T	Solid	5030B	
MB 720-29834/3-A	Method Blank	T	Solid	5030B	
720-12227-A-1-F MS	Matrix Spike	T	Solid	5030B	
720-12227-A-1-G MSD	Matrix Spike Duplicate	T	Solid	5030B	
720-12284-1	MWB1 @ 5-1/2'	T	Solid	5030B	
720-12284-2	MWB1 @ 10-1/2'	T	Solid	5030B	
720-12284-3	MWB2 @ 25-1/2'	T	Solid	5030B	
Analysis Batch:720-29888	3				
LCS 720-29834/1-A	Lab Control Spike	T	Solid	8260B	720-29834
LCSD 720-29834/2-A	Lab Control Spike Duplicate	T	Solid	8260B	720-29834
MB 720-29834/3-A	Method Blank	T	Solid	8260B	720-29834
720-12227-A-1-F MS	Matrix Spike	Т	Solid	8260B	720-29834
720-12227-A-1-G MSD	Matrix Spike Duplicate	T	Solid	8260B	720-29834
720-12284-1	MWB1 @ 5-1/2'	Т	Solid	8260B	720-29834
720-12284-2	MWB1 @ 10-1/2'	Т	Solid	8260B	720-29834
720-12284-3	MWB2 @ 25-1/2'	Т	Solid	8260B	720-29834
Analysis Batch:720-30082	2				
LCS 720-30083/1-A	Lab Control Spike	Т	Solid	8260B	720-30083
LCSD 720-30083/2-A	Lab Control Spike Duplicate	Т	Solid	8260B	720-30083
MB 720-30083/3-A	Method Blank	Т	Solid	8260B	720-30083
720-12284-1	MWB1 @ 5-1/2'	Т	Solid	8260B	720-30083
720-12284-1MS	Matrix Spike	Т	Solid	8260B	720-30083
720-12284-1MSD	Matrix Spike Duplicate	T	Solid	8260B	720-30083
720-12284-2	MWB1 @ 10-1/2'	T	Solid	8260B	720-30083
720-12284-3	MWB2 @ 25-1/2'	Т	Solid	8260B	720-30083
Prep Batch: 720-30083					
LCS 720-30083/1-A	Lab Control Spike	Т	Solid	5030B	
LCSD 720-30083/2-A	Lab Control Spike Duplicate	Т	Solid	5030B	
MB 720-30083/3-A	Method Blank	T	Solid	5030B	
720-12284-1	MWB1 @ 5-1/2'	T	Solid	5030B	
720-12284-1MS	Matrix Spike	Т	Solid	5030B	
720-12284-1MSD	Matrix Spike Duplicate	T	Solid	5030B	
720-12284-2	MWB1 @ 10-1/2'	Т	Solid	5030B	
720-12284-3	MWB2 @ 25-1/2'	Т	Solid	5030B	

Report Basis

T = Total

Client: Engeo, Inc. Job Number: 720-12284-1

QC Association Summary

		Report			
Lab Sample ID	Client Sample ID	Basis	Client Matrix	Method	Prep Batch
GC Semi VOA					
Prep Batch: 720-30218	3				
LCS 720-30218/2-A	Lab Control Spike	Α	Solid	3550B	
LCSD 720-30218/3-A	Lab Control Spike Duplicate	Α	Solid	3550B	
MB 720-30218/1-A	Method Blank	Α	Solid	3550B	
720-12284-1	MWB1 @ 5-1/2'	Α	Solid	3550B	
720-12284-2	MWB1 @ 10-1/2'	Α	Solid	3550B	
720-12284-3	MWB2 @ 25-1/2'	Α	Solid	3550B	
Analysis Batch:720-30	339				
LCS 720-30218/2-A	Lab Control Spike	Α	Solid	8015B	720-30218
LCSD 720-30218/3-A	Lab Control Spike Duplicate	Α	Solid	8015B	720-30218
MB 720-30218/1-A	Method Blank	Α	Solid	8015B	720-30218
720-12284-1	MWB1 @ 5-1/2'	Α	Solid	8015B	720-30218
720-12284-2	MWB1 @ 10-1/2'	Α	Solid	8015B	720-30218
720-12284-3	MWB2 @ 25-1/2'	Α	Solid	8015B	720-30218

Report Basis

A = Silica Gel Cleanup

Job Number: 720-12284-1 Client: Engeo, Inc.

Method Blank - Batch: 720-29834 Method: 8260B Preparation: 5030B

Lab Sample ID: MB 720-29834/3-A Analysis Batch: 720-29888 Instrument ID: Agilent 75MSD Client Matrix: Solid Prep Batch: 720-29834 Lab File ID: 121907005.D Units: ug/Kg Dilution: Initial Weight/Volume: 5 g 1.0 Final Weight/Volume: 10 mL

Date Analyzed: 12/19/2007 1247 Date Prepared: 12/19/2007 1200

Analyte	Result	Qual	RL
Methyl tert-butyl ether	ND		5.0
Acetone	ND		50
Benzene	ND		5.0
Dichlorobromomethane	ND		5.0
Bromobenzene	ND		5.0
Chlorobromomethane	ND		20
Bromoform	ND		5.0
Bromomethane	ND		10
2-Butanone (MEK)	ND		50
n-Butylbenzene	ND		5.0
sec-Butylbenzene	ND		5.0
tert-Butylbenzene	ND		5.0
Carbon disulfide	ND		5.0
Carbon tetrachloride	ND		5.0
Chlorobenzene	ND		5.0
Chloroethane	ND		10
Chloroform	ND		5.0
Chloromethane	ND		10
2-Chlorotoluene	ND		5.0
4-Chlorotoluene	ND		5.0
Chlorodibromomethane	ND		5.0
1,2-Dichlorobenzene	ND		5.0
1,3-Dichlorobenzene	ND		5.0
1,4-Dichlorobenzene	ND		5.0
1,3-Dichloropropane	ND		5.0
1,1-Dichloropropene	ND		5.0
1,2-Dibromo-3-Chloropropane	ND		50
Ethylene Dibromide	ND		5.0
Dibromomethane	ND		10
Dichlorodifluoromethane	ND		10
1,1-Dichloroethane	ND		5.0
1,2-Dichloroethane	ND		5.0
1,1-Dichloroethene	ND		5.0
cis-1,2-Dichloroethene	ND		5.0
trans-1,2-Dichloroethene	ND		5.0
1,2-Dichloropropane	ND		5.0
cis-1,3-Dichloropropene	ND		5.0
trans-1,3-Dichloropropene	ND		5.0
Ethylbenzene	ND		5.0
Hexachlorobutadiene	ND		5.0
2-Hexanone	ND		50

Calculations are performed before rounding to avoid round-off errors in calculated results.

Job Number: 720-12284-1 Client: Engeo, Inc.

Method Blank - Batch: 720-29834 Method: 8260B Preparation: 5030B

Lab Sample ID: MB 720-29834/3-A Analysis Batch: 720-29888 Instrument ID: Agilent 75MSD Client Matrix: Solid Prep Batch: 720-29834 Lab File ID: 121907005.D

Dilution: 1.0 Units: ug/Kg Initial Weight/Volume: 5 g Final Weight/Volume: 10 mL Date Analyzed: 12/19/2007 1247 Date Prepared: 12/19/2007 1200

Analyte	Result	Qual	RL
Isopropylbenzene	ND		5.0
4-Isopropyltoluene	ND		5.0
Methylene Chloride	ND		10
4-Methyl-2-pentanone (MIBK)	ND		50
Naphthalene	ND		10
N-Propylbenzene	ND		5.0
Styrene	ND		5.0
1,1,1,2-Tetrachloroethane	ND		5.0
1,1,2,2-Tetrachloroethane	ND		5.0
Tetrachloroethene	ND		5.0
Toluene	ND		5.0
1,2,3-Trichlorobenzene	ND		5.0
1,2,4-Trichlorobenzene	ND		5.0
1,1,1-Trichloroethane	ND		5.0
1,1,2-Trichloroethane	ND		5.0
Trichloroethene	ND		5.0
Trichlorofluoromethane	ND		5.0
1,2,3-Trichloropropane	ND		5.0
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		5.0
1,2,4-Trimethylbenzene	ND		5.0
1,3,5-Trimethylbenzene	ND		5.0
Vinyl acetate	ND		50
Vinyl chloride	ND		5.0
Xylenes, Total	ND		10
2,2-Dichloropropane	ND		5.0
Surrogate	% Rec	Acceptance Limits	
4-Bromofluorobenzene	99	50 - 138	
1,2-Dichloroethane-d4 (Surr)	97	66 - 127	
Toluene-d8 (Surr)	97	51 - 129	

Calculations are performed before rounding to avoid round-off errors in calculated results.

Client: Engeo, Inc. Job Number: 720-12284-1

Lab Control Spike/ Method: 8260B
Lab Control Spike Duplicate Recovery Report - Batch: 720-29834 Preparation: 5030B

LCS Lab Sample ID: LCS 720-29834/1-A

Client Matrix: Solid Dilution: 1.0

Date Analyzed: 12/19/2007 1157 Date Prepared: 12/19/2007 1200 Analysis Batch: 720-29888 Prep Batch: 720-29834

Units: ug/Kg

Instrument ID: Agilent 75MSD Lab File ID: 121907003.D

Initial Weight/Volume: 5 g Final Weight/Volume: 10 mL

LCSD Lab Sample ID: LCSD 720-29834/2-A

Client Matrix: Solid Dilution: 1.0

Date Analyzed: 12/19/2007 1222 Date Prepared: 12/19/2007 1200 Analysis Batch: 720-29888 Prep Batch: 720-29834

Units: ug/Kg

Instrument ID: Agilent 75MSD Lab File ID: 121907004.D Initial Weight/Volume: 5 g

Final Weight/Volume: 10 mL

	9	<u> 6 Rec.</u>					
Analyte	LCS	LCSD	Limit	RPD	RPD Limit	LCS Qual	LCSD Qual
Benzene	97	91	80 - 120	6	20		
Chlorobenzene	99	92	86 - 115	7	20		
1,1-Dichloroethene	108	102	81 - 140	6	20		
Toluene	98	91	81 - 120	8	20		
Trichloroethene	100	93	82 - 118	6	20		
Surrogate	L	.CS % Rec	LCSD %	Rec	Acce	otance Limits	;
4-Bromofluorobenzene	1	00	102		5	0 - 138	
1,2-Dichloroethane-d4 (Surr)	1	04	98		6	66 - 127	
Toluene-d8 (Surr)	9	8	97		5	1 - 129	

Client: Engeo, Inc. Job Number: 720-12284-1

Matrix Spike/ Method: 8260B
Matrix Spike Duplicate Recovery Report - Batch: 720-29834 Preparation: 5030B

MS Lab Sample ID: 720-12227-A-1-F MS Analysis Batch: 720-29888 Instrument ID: Agilent 75MSD

Client Matrix: Solid Prep Batch: 720-29834 Lab File ID: 121907010.D Dilution: 1.0 Initial Weight/Volume: 5.26 q

Dilution: 1.0 Initial Weight/Volume: 5.26 g

Date Analyzed: 12/19/2007 1451 Final Weight/Volume: 10 mL

Date Prepared: 12/19/2007 1200

MSD Lab Sample ID: 720-12227-A-1-G MSD Analysis Batch: 720-29888 Instrument ID: Agilent 75MSD

Client Matrix: Solid Prep Batch: 720-29834 Lab File ID: 121907011.D Dilution: 1.0 Initial Weight/Volume: 5.02 g

Dilution: 1.0 Initial Weight/Volume: 5.02 g
Date Analyzed: 12/19/2007 1516 Final Weight/Volume: 10 mL
Date Prepared: 12/19/2007 1200

% Rec. MSD **RPD** MS Qual MSD Qual Analyte MS Limit **RPD Limit** Benzene 56 63 - 126 11 20 F F 53 Chlorobenzene 22 25 57 - 124 20 20 F F 1,1-Dichloroethene 76 80 66 - 149 10 20 F F Toluene 35 38 54 - 131 12 20 53 - 130 F F Trichloroethene 46 50 14 20 MS % Rec MSD % Rec Surrogate Acceptance Limits 21 4-Bromofluorobenzene 16 Χ Χ 50 - 138 66 - 127 76 1,2-Dichloroethane-d4 (Surr) 73 51 - 129 Toluene-d8 (Surr) 38 Χ 37 Χ

Client: Engeo, Inc. Job Number: 720-12284-1

Method Blank - Batch: 720-30083 Method: 8260B Preparation: 5030B

Lab Sample ID: MB 720-30083/3-A Analysis Batch: 720-30082 Instrument ID: Varian 3900E

Client Matrix: Solid Prep Batch: 720-30083 Lab File ID: c:\varianws\data\200712\12

Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 5 g
Date Analyzed: 12/26/2007 1035 Final Weight/Volume: 10 mL
Date Prepared: 12/26/2007 0835

Qual RLAnalyte Result Benzene ND 0.0050 Toluene ND 0.0050 Gasoline Range Organics (GRO)-C5-C12 ND 0.25 Surrogate % Rec Acceptance Limits 1,2-Dichloroethane-d4 (Surr) 102 60 - 140 Toluene-d8 (Surr) 99 70 - 130

Lab Control Spike/ Method: 8260B
Lab Control Spike Duplicate Recovery Report - Batch: 720-30083 Preparation: 5030B

LCS Lab Sample ID: LCS 720-30083/1-A Analysis Batch: 720-30082 Instrument ID: Varian 3900E

Client Matrix: Solid Prep Batch: 720-30083 Lab File ID: c:\varianws\data\200712\12

Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 5 g
Date Analyzed: 12/26/2007 1058 Final Weight/Volume: 10 mL

Date Prepared: 12/26/2007 0835

LCSD Lab Sample ID: LCSD 720-30083/2-A Analysis Batch: 720-30082 Instrument ID: Varian 3900E

Client Matrix: Solid Prep Batch: 720-30083 Lab File ID: c:\varianws\data\200712\122 Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 5 g

Date Analyzed: 12/26/2007 1121 Final Weight/Volume: 10 mL
Date Prepared: 12/26/2007 0835

% Rec. Analyte LCS **LCSD** Limit **RPD** RPD Limit LCS Qual LCSD Qual Benzene 94 91 70 - 123 3 20 Toluene 101 97 81 - 128 4 20 2 Gasoline Range Organics (GRO)-C5-C12 51 - 97 20 65 66 LCS % Rec Surrogate LCSD % Rec Acceptance Limits 1.2-Dichloroethane-d4 (Surr) 99 95 60 - 140 Toluene-d8 (Surr) 99 97 70 - 130

Calculations are performed before rounding to avoid round-off errors in calculated results.

Client: Engeo, Inc. Job Number: 720-12284-1

Matrix Spike/ Method: 8260B
Matrix Spike Duplicate Recovery Report - Batch: 720-30083 Preparation: 5030B

MS Lab Sample ID: 720-12284-1 Analysis Batch: 720-30082 Instrument ID: Varian 3900E

Client Matrix: Solid Prep Batch: 720-30083 Lab File ID: c:\varianws\data\200712\

Dilution: 1.0 Initial Weight/Volume: 5.06 g

Date Analyzed: 12/26/2007 1416 Final Weight/Volume: 10 mL Date Prepared: 12/26/2007 0835

MSD Lab Sample ID: 720-12284-1 Analysis Batch: 720-30082 Instrument ID: Varian 3900E

Client Matrix: Solid Prep Batch: 720-30083 Lab File ID: c:\varianws\data\200712\12

Dilution: 1.0 Initial Weight/Volume: 5.49 g

Date Analyzed: 12/26/2007 1442 Final Weight/Volume: 10 mL
Date Prepared: 12/26/2007 0835

% Rec. MSD Limit **RPD RPD** Limit MS Qual MSD Qual Analyte MS Benzene 70 - 123 89 85 12 20 Toluene 95 90 81 - 128 14 20 F Gasoline Range Organics (GRO)-C5-C12 53 49 51 - 97 15 20 MS % Rec MSD % Rec Surrogate Acceptance Limits 1,2-Dichloroethane-d4 (Surr) 100 107 60 - 140 Toluene-d8 (Surr) 101 97 70 - 130

Client: Engeo, Inc. Job Number: 720-12284-1

Method Blank - Batch: 720-30218 Method: 8015B
Preparation: 3550B
Silica Gel Cleanup

Lab Sample ID: MB 720-30218/1-A

Analysis Batch: 720-30339

Client Matrix: Solid

Analysis Batch: 720-30339

Instrument ID: HP DRO5

Lab File ID: N/A

Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 30.14 g

Date Analyzed: 01/02/2008 1350 Final Weight/Volume: 5 mL
Date Prepared: 12/28/2007 1556 Injection Volume:

Column ID: PRIMARY

 Analyte
 Result
 Qual
 RL

 Diesel Range Organics [C10-C28]
 ND
 1.0

 Motor Oil Range Organics [C24-C36]
 ND
 50

 Surrogate
 % Rec
 Acceptance Limits

 Capric Acid (Surr)
 1
 0 - 5

 p-Terphenyl
 88
 41 - 105

Lab Control Spike/
Lab Control Spike Duplicate Recovery Report - Batch: 720-30218

Method: 8015B
Preparation: 3550B
Silica Gel Cleanup

LCS Lab Sample ID: LCS 720-30218/2-A Analysis Batch: 720-30339 Instrument ID: HP DRO5

Client Matrix: Solid Prep Batch: 720-30218 Lab File ID: N/A
Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 3

Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 30.13 g
Date Analyzed: 01/02/2008 1256 Final Weight/Volume: 5 mL
Date Prepared: 12/28/2007 1556 Injection Volume:

Column ID:

PRIMARY

LCSD Lab Sample ID: LCSD 720-30218/3-A Analysis Batch: 720-30339 Instrument ID: HP DRO5

Client Matrix: Solid Prep Batch: 720-30218 Lab File ID: N/A

Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 30.07 g
Date Analyzed: 01/02/2008 1323 Final Weight/Volume: 5 mL

Date Prepared: 12/28/2007 1556 Injection Volume:

Column ID: PRIMARY

% Rec. Analyte LCS **LCSD** Limit **RPD** RPD Limit LCS Qual LCSD Qual Diesel Range Organics [C10-C28] 64 50 - 130 3 30 66 LCS % Rec Surrogate LCSD % Rec Acceptance Limits 103 105 41 - 105 p-Terphenyl

Calculations are performed before rounding to avoid round-off errors in calculated results.

STL San Francisco Chain of Custody 1220 Quarry Lane • Pleasanton CA 94566-4756 Reference #: 168683 Phone: (925) 484-1919 • Fax: (925) 484-1096 Date 12/18/07 Page 1 of 1 Email: sflogin@stl-inc.com Report To Analysis Request ENGEO ampany: RICHARD Gandafo 808 Low Level Metals by EPA 200,8/6020 (ICP-MS): Metals: Cl Lead Cl LUFT Cl RCRA SPON. Wilnow Are Ripon Hexavalent Chromlum pH (24h hold time far H₂O) Alkalinity TDS [Purgeable Arematics BTEX EPA - D 8021 🗀 02608 D NO. hone: 209 831 ObloEmail: rgandulfo ocage CAM17 Metals (EPA 6010/7470/7471) EPA (0.00 ill To: Sampled By: W.E.T (STLC) TCLP Containers Spec Cond. TSS R. Gandulfo ti di ttn: Phone: ö Baimple ID The Pres Time - Mat. Pres tix erv. MWB1 B51/12/18 9:00 ice 5 MWB1 0 1012 9:10 MWB20258 13:30 Project Info. Sample Receipt 1) Relinguished by: 2) Relinquished by: 3) Relinquished by: # of Containers: RICKENBACKER Signature Time Signature Head Space: Time Rich Gardor Signature Time 7584.1.001.01 Printed Name Date Printed Name Date ENGED adit Card# Conforms to record: Company Company Company 1) Received by: 48h 2) Received by: 24h Other: 3) Received by: port: Routine | Level 3 | Level 4 | EDD State Tenk Fund EDF cial Instructions / Comments; Signature Time Signature ☐ Global ID Time 1/04/2006 reports 8015M from C₈-C₂₄ (Industry norm). Default for 8015B is C₁₀-C₂₈. Printed Name Date Printed Name Date Company Company

Login Sample Receipt Check List

Client: Engeo, Inc. Job Number: 720-12284-1

List Source: TestAmerica San Francisco

Login Number: 12284 Creator: Bullock, Tracy

List Number: 1

Question T / F/ NA Comment Radioactivity either was not measured or, if measured, is at or below N/A background The cooler's custody seal, if present, is intact. N/A The cooler or samples do not appear to have been compromised or True tampered with. Samples were received on ice. True True Cooler Temperature is acceptable. Cooler Temperature is recorded. True COC is present. True COC is filled out in ink and legible. True COC is filled out with all pertinent information. True There are no discrepancies between the sample IDs on the containers and True the COC. Samples are received within Holding Time. True Sample containers have legible labels. True Containers are not broken or leaking. True Sample collection date/times are provided. True Appropriate sample containers are used. True Sample bottles are completely filled. True There is sufficient vol. for all requested analyses, incl. any requested True MS/MSDs VOA sample vials do not have headspace or bubble is <6mm (1/4") in True diameter. If necessary, staff have been informed of any short hold time or quick TAT True needs Multiphasic samples are not present. True Samples do not require splitting or compositing. True



ANALYTICAL REPORT

Job Number: 720-12301-1

Job Description: Rickenbacker

For:
Engeo, Inc.
580 N Wilma Avenue
Suite A
Ripon, CA 95366-9502

Attention: Mr. Richard Gandolfo

Akanef Sal

Designee for
Melissa Brewer
Project Manager I
melissa.brewer@testamericainc.com
01/04/2008

Job Narrative 720-J12301-1

Comments

No additional comments.

Receipt

All samples were received in good condition within temperature requirements.

GC/MS VOA

No analytical or quality issues were noted.

GC VOA

No analytical or quality issues were noted.

GC Semi VOA

Method(s) 8015B: Capric Surrogate recovery for the MB was outside the upper control limit: (MB 720-30239/1-A). This MB did not contain any target analytes; therefore, re-extraction and/or re-analysis was not performed.

Method(s) 8015B: The matrix spike / matrix spike duplicate (MS/MSD) recoveries for batch 30306 were outside control limits. The associated laboratory control standard (LCS) met acceptance criteria.

No other analytical or quality issues were noted.

Organic Prep

No analytical or quality issues were noted.

EXECUTIVE SUMMARY - Detections

Client: Engeo, Inc. Job Number: 720-12301-1

Lab Sample ID Analyte	Client Sample ID	Result / Qualifier	Reporting Limit	Units	Method
720-12301-1	MW3 @ 26'				
Silica Gel Cleanup Diesel Range Orga		2.2	0.99	mg/Kg	8015B

METHOD SUMMARY

Client: Engeo, Inc.

Job Number: 720-12301-1

Description	Lab Location	Method	Preparation Method
Matrix: Solid			
Volatile Organic Compounds by GC/MS	TAL SF	SW846 8260B	
Volatile Organic Compounds by GC/MS (Low Level) Purge and Trap for Solids Purge and Trap for Solids	TAL SF TAL SF TAL SF	SW846 8260B	SW846 5030B SW846 5030B
Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)	TAL SF	SW846 8015B	
Ultrasonic Extraction	TAL SF		SW846 3550B

Lab References:

TAL SF = TestAmerica San Francisco

Method References:

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

METHOD / ANALYST SUMMARY

Client: Engeo, Inc. Job Number: 720-12301-1

Method	Analyst	Analyst ID
SW846 8260B SW846 8260B	Ali, Badri Le, Lien	BA LL
SW846 8015B	Relja, Marlene	MR

SAMPLE SUMMARY

Client: Engeo, Inc. Job Number: 720-12301-1

Lab Sample ID	Client Sample ID	Client Matrix	Date/Time Sampled	Date/Time Received
720-12301-1	MW3 @ 26'	Solid	12/19/2007 0845	12/19/2007 1134

Client: Engeo, Inc. Job Number: 720-12301-1

Client Sample ID: MW3 @ 26'

 Lab Sample ID:
 720-12301-1
 Date Sampled:
 12/19/2007 0845

 Client Matrix:
 Solid
 Date Received:
 12/19/2007 1134

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29955Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29901Lab File ID:122007012.DDilution:1.0Initial Weight/Volume:5.38

Dilution: 1.0 Initial Weight/Volume: 5.38 g
Date Analyzed: 12/20/2007 1536 Final Weight/Volume: 10 mL

Date Prepared: 12/20/2007 1100

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
Methyl tert-butyl ether	ND		4.6
Acetone	ND		46
Benzene	ND		4.6
Dichlorobromomethane	ND		4.6
Bromobenzene	ND		4.6
Chlorobromomethane	ND		19
Bromoform	ND		4.6
Bromomethane	ND		9.3
2-Butanone (MEK)	ND		46
n-Butylbenzene	ND		4.6
sec-Butylbenzene	ND		4.6
tert-Butylbenzene	ND		4.6
Carbon disulfide	ND		4.6
Carbon tetrachloride	ND		4.6
Chlorobenzene	ND		4.6
Chloroethane	ND		9.3
Chloroform	ND		4.6
Chloromethane	ND		9.3
2-Chlorotoluene	ND		4.6
4-Chlorotoluene	ND		4.6
Chlorodibromomethane	ND		4.6
1,2-Dichlorobenzene	ND		4.6
1,3-Dichlorobenzene	ND		4.6
1,4-Dichlorobenzene	ND		4.6
1,3-Dichloropropane	ND		4.6
1,1-Dichloropropene	ND		4.6
1,2-Dibromo-3-Chloropropane	ND		46
Ethylene Dibromide	ND		4.6
Dibromomethane	ND		9.3
Dichlorodifluoromethane	ND		9.3
1,1-Dichloroethane	ND		4.6
1,2-Dichloroethane	ND		4.6
1,1-Dichloroethene	ND		4.6
cis-1,2-Dichloroethene	ND		4.6
trans-1,2-Dichloroethene	ND		4.6
1,2-Dichloropropane	ND		4.6
cis-1,3-Dichloropropene	ND		4.6
trans-1,3-Dichloropropene	ND		4.6
Ethylbenzene	ND		4.6
Hexachlorobutadiene	ND		4.6
2-Hexanone	ND		46
Isopropylbenzene	ND		4.6
4-Isopropyltoluene	ND		4.6
Methylene Chloride	ND		9.3

Client: Engeo, Inc. Job Number: 720-12301-1

Client Sample ID: MW3 @ 26'

 Lab Sample ID:
 720-12301-1
 Date Sampled:
 12/19/2007 0845

 Client Matrix:
 Solid
 Date Received:
 12/19/2007 1134

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method:8260BAnalysis Batch: 720-29955Instrument ID:Agilent 75MSDPreparation:5030BPrep Batch: 720-29901Lab File ID:122007012.DDilution:1.0Initial Weight/Volume:5.38 g

Date Analyzed: 12/20/2007 1536 Final Weight/Volume: 10 mL

Date Prepared: 12/20/2007 1100

Analyte	DryWt Corrected: N Result (ug/Kg)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		46
Naphthalene	ND		9.3
N-Propylbenzene	ND		4.6
Styrene	ND		4.6
1,1,1,2-Tetrachloroethane	ND		4.6
1,1,2,2-Tetrachloroethane	ND		4.6
Tetrachloroethene	ND		4.6
Toluene	ND		4.6
1,2,3-Trichlorobenzene	ND		4.6
1,2,4-Trichlorobenzene	ND		4.6
1,1,1-Trichloroethane	ND		4.6
1,1,2-Trichloroethane	ND		4.6
Trichloroethene	ND		4.6
Trichlorofluoromethane	ND		4.6
1,2,3-Trichloropropane	ND		4.6
1,1,2-Trichloro-1,2,2-trifluoroethan	e ND		4.6
1,2,4-Trimethylbenzene	ND		4.6
1,3,5-Trimethylbenzene	ND		4.6
Vinyl acetate	ND		46
Vinyl chloride	ND		4.6
Xylenes, Total	ND		9.3
2,2-Dichloropropane	ND		4.6
Surrogate	%Rec		Acceptance Limits
4-Bromofluorobenzene	99		50 - 138
1,2-Dichloroethane-d4 (Surr)	109		66 - 127
			51 - 129
Toluene-d8 (Surr)	107		31 - 129

Client: Engeo, Inc. Job Number: 720-12301-1

Client Sample ID: MW3 @ 26'

 Lab Sample ID:
 720-12301-1
 Date Sampled:
 12/19/2007 0845

 Client Matrix:
 Solid
 Date Received:
 12/19/2007 1134

8260B Volatile Organic Compounds by GC/MS

Method: 8260B Analysis Batch: 720-30155 Instrument ID: Varian 3900A

Preparation: 5030B Prep Batch: 720-30114 Lab File ID: c:\saturnws\data\200712\12

Dilution: 1.0 Initial Weight/Volume: 5.43 g

Date Analyzed: 12/27/2007 1330 Final Weight/Volume: 10 mL

Date Prepared: 12/27/2007 0930

Analyte DryWt Corrected: N Result (mg/Kg) Qualifier RL

Gasoline Range Organics (GRO)-C5-C12 ND 0.23

 Surrogate
 %Rec
 Acceptance Limits

 1,2-Dichloroethane-d4 (Surr)
 94
 60 - 140

 Toluene-d8 (Surr)
 104
 70 - 130

Client: Engeo, Inc. Job Number: 720-12301-1

Client Sample ID: MW3 @ 26'

 Lab Sample ID:
 720-12301-1
 Date Sampled:
 12/19/2007 0845

 Client Matrix:
 Solid
 Date Received:
 12/19/2007 1134

8015B Nonhalogenated Organics using GC/FID -Modified (Diesel Range Organics)-Silica Gel Cleanup

Method: 8015B Analysis Batch: 720-30306 Instrument ID: HP DRO5

Preparation: 3550B Prep Batch: 720-30239 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 30.38 g
Date Analyzed: 12/31/2007 1853 Final Weight/Volume: 5 mL

Date Prepared: 12/31/2007 0905 Injection Volume:

Column ID: PRIMARY

Analyte	DryWt Corrected: N Result (mg/Kg)	Qualifier	RL
Diesel Range Organics [C10-C2	28] 2.2		0.99
Motor Oil Range Organics [C24	-C361 ND		49

Surrogate	%Rec	Acceptance Limits
Capric Acid (Surr)	1	0 - 5
p-Terphenyl	83	41 - 105

DATA REPORTING QUALIFIERS

Client: Engeo, Inc. Job Number: 720-12301-1

Lab Section	Qualifier	Description
GC Semi VOA		
	4	MS, MSD: The analyte present in the original sample is 4 times greater than the matrix spike concentration; therefore, control limits are not applicable.
	Χ	Surrogate exceeds the control limits

Client: Engeo, Inc. Job Number: 720-12301-1

QC Association Summary

		Report			
Lab Sample ID	Client Sample ID	Basis	Client Matrix	Method	Prep Batch
GC/MS VOA					
Prep Batch: 720-29901					
LCS 720-29901/1-A	Lab Control Spike	T	Solid	5030B	
LCSD 720-29901/2-A	Lab Control Spike Duplicate	T	Solid	5030B	
MB 720-29901/3-A	Method Blank	T	Solid	5030B	
720-12297-A-1-H MS	Matrix Spike	Т	Solid	5030B	
720-12297-A-1-I MSD	Matrix Spike Duplicate	Т	Solid	5030B	
720-12301-1	MW3 @ 26'	Т	Solid	5030B	
Analysis Batch:720-299	955				
LCS 720-29901/1-A	Lab Control Spike	Т	Solid	8260B	720-29901
LCSD 720-29901/2-A	Lab Control Spike Duplicate	Ť	Solid	8260B	720-29901
MB 720-29901/3-A	Method Blank	Т	Solid	8260B	720-29901
720-12297-A-1-H MS	Matrix Spike	Т	Solid	8260B	720-29901
720-12297-A-1-I MSD	Matrix Spike Duplicate	T	Solid	8260B	720-29901
720-12301-1	MW3 @ 26'	Т	Solid	8260B	720-29901
Prep Batch: 720-30114					
LCS 720-30114/2-A	Lab Control Spike	Т	Solid	5030B	
LCSD 720-30114/3-A	Lab Control Spike Duplicate	Т	Solid	5030B	
MB 720-30114/1-A	Method Blank	Т	Solid	5030B	
720-12301-1	MW3 @ 26'	Т	Solid	5030B	
720-12380-A-5-D MS	Matrix Spike	Т	Solid	5030B	
720-12380-A-5-E MSD	Matrix Spike Duplicate	Т	Solid	5030B	
Analysis Batch:720-30	155				
LCS 720-30114/2-A	Lab Control Spike	Т	Solid	8260B	720-30114
LCSD 720-30114/3-A	Lab Control Spike Duplicate	Т	Solid	8260B	720-30114
MB 720-30114/1-A	Method Blank	Ť	Solid	8260B	720-30114
720-12301-1	MW3 @ 26'	Т	Solid	8260B	720-30114
720-12380-A-5-D MS	Matrix Spike	Т	Solid	8260B	720-30114
720-12380-A-5-E MSD	Matrix Spike Duplicate	Ť	Solid	8260B	720-30114

Report Basis T = Total

Client: Engeo, Inc. Job Number: 720-12301-1

QC Association Summary

		Report			
Lab Sample ID	Client Sample ID	Basis	Client Matrix	Method	Prep Batch
GC Semi VOA					
Prep Batch: 720-30239					
LCS 720-30239/2-A	Lab Control Spike	Α	Solid	3550B	
LCSD 720-30239/3-A	Lab Control Spike Duplicate	Α	Solid	3550B	
MB 720-30239/1-A	Method Blank	Α	Solid	3550B	
720-12301-1	MW3 @ 26'	Α	Solid	3550B	
720-12365-H-1-C MS	Matrix Spike	Α	Solid	3550B	
720-12365-H-1-D MSD	Matrix Spike Duplicate	Α	Solid	3550B	
Analysis Batch:720-30	306				
LCS 720-30239/2-A	Lab Control Spike	Α	Solid	8015B	720-30239
LCSD 720-30239/3-A	Lab Control Spike Duplicate	Α	Solid	8015B	720-30239
MB 720-30239/1-A	Method Blank	Α	Solid	8015B	720-30239
720-12301-1	MW3 @ 26'	Α	Solid	8015B	720-30239
720-12365-H-1-C MS	Matrix Spike	Α	Solid	8015B	720-30239
720-12365-H-1-D MSD	Matrix Spike Duplicate	Α	Solid	8015B	720-30239

Report Basis

A = Silica Gel Cleanup

RL

Client: Engeo, Inc. Job Number: 720-12301-1

Method Blank - Batch: 720-29901 Method: 8260B Preparation: 5030B

Lab Sample ID: MB 720-29901/3-A Analysis Batch: 720-29955 Instrument ID: Agilent 75MSD Client Matrix: Solid Prep Batch: 720-29901 Lab File ID: 122007005.D Dilution: 1.0 Units: ug/Kg Initial Weight/Volume: 5 g Date Analyzed: 12/20/2007 1242 Final Weight/Volume: 10 mL

Result

Qual

Date Analyzed: 12/20/2007 1242 Date Prepared: 12/20/2007 1100

Analyte

Analyte	Result Qu	al RL
Methyl tert-butyl ether	ND	5.0
Acetone	ND	50
Benzene	ND	5.0
Dichlorobromomethane	ND	5.0
Bromobenzene	ND	5.0
Chlorobromomethane	ND	20
Bromoform	ND	5.0
Bromomethane	ND	10
2-Butanone (MEK)	ND	50
n-Butylbenzene	ND	5.0
sec-Butylbenzene	ND	5.0
tert-Butylbenzene	ND	5.0
Carbon disulfide	ND	5.0
Carbon tetrachloride	ND	5.0
Chlorobenzene	ND	5.0
Chloroethane	ND	10
Chloroform	ND	5.0
Chloromethane	ND	10
2-Chlorotoluene	ND	5.0
4-Chlorotoluene	ND	5.0
Chlorodibromomethane	ND	5.0
1,2-Dichlorobenzene	ND	5.0
1,3-Dichlorobenzene	ND	5.0
1,4-Dichlorobenzene	ND	5.0
1,3-Dichloropropane	ND	5.0
1,1-Dichloropropene	ND	5.0
1,2-Dibromo-3-Chloropropane	ND	50
Ethylene Dibromide	ND	5.0
Dibromomethane	ND	10
Dichlorodifluoromethane	ND	10
1,1-Dichloroethane	ND	5.0
1,2-Dichloroethane	ND	5.0
1,1-Dichloroethene	ND	5.0
cis-1,2-Dichloroethene	ND	5.0
trans-1,2-Dichloroethene	ND	5.0
1,2-Dichloropropane	ND	5.0
cis-1,3-Dichloropropene	ND	5.0
trans-1,3-Dichloropropene	ND	5.0
Ethylbenzene	ND	5.0
Hexachlorobutadiene	ND	5.0
2-Hexanone	ND	50

Client: Engeo, Inc. Job Number: 720-12301-1

Method Blank - Batch: 720-29901 Method: 8260B Preparation: 5030B

Lab Sample ID: MB 720-29901/3-A Analysis Batch: 720-29955 Instrument ID: Agilent 75MSD Client Matrix: Solid Prep Batch: 720-29901 Lab File ID: 122007005.D

Dilution: 1.0 Units: ug/Kg Initial Weight/Volume: 5 g
Date Analyzed: 12/20/2007 1242 Final Weight/Volume: 10 mL
Date Prepared: 12/20/2007 1100

Analyte	Result	Qual	RL
Isopropylbenzene	ND		5.0
4-Isopropyltoluene	ND		5.0
Methylene Chloride	ND		10
4-Methyl-2-pentanone (MIBK)	ND		50
Naphthalene	ND		10
N-Propylbenzene	ND		5.0
Styrene	ND		5.0
1,1,1,2-Tetrachloroethane	ND		5.0
1,1,2,2-Tetrachloroethane	ND		5.0
Tetrachloroethene	ND		5.0
Toluene	ND		5.0
1,2,3-Trichlorobenzene	ND		5.0
1,2,4-Trichlorobenzene	ND		5.0
1,1,1-Trichloroethane	ND		5.0
1,1,2-Trichloroethane	ND		5.0
Trichloroethene	ND		5.0
Trichlorofluoromethane	ND		5.0
1,2,3-Trichloropropane	ND		5.0
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		5.0
1,2,4-Trimethylbenzene	ND		5.0
1,3,5-Trimethylbenzene	ND		5.0
Vinyl acetate	ND		50
Vinyl chloride	ND		5.0
Xylenes, Total	ND		10
2,2-Dichloropropane	ND		5.0
Surrogate	% Rec	Acceptance Limits	
4-Bromofluorobenzene	98	50 - 138	
1,2-Dichloroethane-d4 (Surr)	106	66 - 127	
Toluene-d8 (Surr)	101	51 - 129	

Client: Engeo, Inc. Job Number: 720-12301-1

Lab Control Spike/ Method: 8260B
Lab Control Spike Duplicate Recovery Report - Batch: 720-29901 Preparation: 5030B

LCS Lab Sample ID: LCS 720-29901/1-A

Client Matrix: Solid Dilution: 1.0

Date Analyzed: 12/20/2007 1152 Date Prepared: 12/20/2007 1100 Analysis Batch: 720-29955 Prep Batch: 720-29901

Units: ug/Kg

Instrument ID: Agilent 75MSD Lab File ID: 122007003.D

Initial Weight/Volume: 5 g Final Weight/Volume: 10 mL

LCSD Lab Sample ID: LCSD 720-29901/2-A

Client Matrix: Solid Dilution: 1.0

Date Analyzed: 12/20/2007 1217 Date Prepared: 12/20/2007 1100 Analysis Batch: 720-29955 Prep Batch: 720-29901

Units: ug/Kg

Instrument ID: Agilent 75MSD Lab File ID: 122007004.D

Initial Weight/Volume: 5 g Final Weight/Volume: 10 mL

	9	<u> 6 Rec.</u>					
Analyte	LCS	LCSD	Limit	RPD	RPD Limit	LCS Qual	LCSD Qual
Benzene	94	96	80 - 120	2	20		
Chlorobenzene	94	94	86 - 115	0	20		
1,1-Dichloroethene	109	110	81 - 140	0	20		
Toluene	95	96	81 - 120	1	20		
Trichloroethene	96	96	82 - 118	0	20		
Surrogate	L	.CS % Rec	LCSD %	Rec	Acce	ptance Limits	;
4-Bromofluorobenzene	1	01	95		Ę	50 - 138	
1,2-Dichloroethane-d4 (Surr)	1	04	98		(66 - 127	
Toluene-d8 (Surr)	1	01	94		į.	51 - 129	

Client: Engeo, Inc. Job Number: 720-12301-1

Matrix Spike/ Method: 8260B
Matrix Spike Duplicate Recovery Report - Batch: 720-29901 Preparation: 5030B

MS Lab Sample ID: 720-12297-A-1-H MS Analysis Batch: 720-29955 Instrument ID: Agilent 75MSD

Client Matrix: Solid Prep Batch: 720-29901 Lab File ID: 122007013.D Dilution: 1.0 Initial Weight/Volume: 5.09 q

Date Analyzed: 12/20/2007 1602 Final Weight/Volume: 10 mL

Date Prepared: 12/20/2007 1100

MSD Lab Sample ID: 720-12297-A-1-I MSD Analysis Batch: 720-29955 Instrument ID: Agilent 75MSD

Client Matrix: Solid Prep Batch: 720-29901 Lab File ID: 122007014.D

Dilution: 1.0 Initial Weight/Volume: 5.21 g
Date Analyzed: 12/20/2007 1627 Final Weight/Volume: 10 mL
Date Prepared: 12/20/2007 1100

% Rec. MS MSD **RPD** MS Qual MSD Qual Analyte Limit **RPD Limit** Benzene 92 87 63 - 126 8 20 Chlorobenzene 83 75 57 - 124 12 20 1,1-Dichloroethene 106 100 66 - 149 20 Toluene 88 81 54 - 131 10 20 53 - 130 20 Trichloroethene 86 78 12 MS % Rec MSD % Rec Surrogate Acceptance Limits 86 70 4-Bromofluorobenzene 50 - 138 66 - 127 1,2-Dichloroethane-d4 (Surr) 93 81 51 - 129 Toluene-d8 (Surr) 79 91

Client: Engeo, Inc. Job Number: 720-12301-1

Method Blank - Batch: 720-30114 Method: 8260B Preparation: 5030B

Lab Sample ID: MB 720-30114/1-A Analysis Batch: 720-30155 Instrument ID: Varian 3900A

Client Matrix: Solid Prep Batch: 720-30114 Lab File ID: c:\saturnws\data\200712\12

Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 5.00 g

Date Analyzed: 12/27/2007 1151 Final Weight/Volume: 10 mL Date Prepared: 12/27/2007 0930

Analyte	Result	Qual	RL
Benzene	ND		0.0050
Toluene	ND		0.0050
Gasoline Range Organics (GRO)-C5-C12	ND		0.25
Surrogate	% Rec	Acceptance Lim	its
1,2-Dichloroethane-d4 (Surr)	92	60 - 140	
Toluene-d8 (Surr)	100	70 - 130	

Lab Control Spike/ Method: 8260B
Lab Control Spike Duplicate Recovery Report - Batch: 720-30114 Preparation: 5030B

LCS Lab Sample ID: LCS 720-30114/2-A Analysis Batch: 720-30155 Instrument ID: Varian 3900A

Client Matrix: Solid Prep Batch: 720-30114 Lab File ID: c:\saturnws\data\200712\12

Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 5.00 g
Date Analyzed: 12/27/2007 1214 Final Weight/Volume: 10 mL

Date Analyzed: 12/27/2007 1214 Final Weight/Volume: 10 mL Date Prepared: 12/27/2007 0930

LCSD Lab Sample ID: LCSD 720-30114/3-A Analysis Batch: 720-30155 Instrument ID: Varian 3900A

Client Matrix: Solid Prep Batch: 720-30114 Lab File ID: c:\saturnws\data\200712\122 Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 5.00 g

Date Analyzed: 12/27/2007 1236 Final Weight/Volume: 10 mL Date Prepared: 12/27/2007 0930

% Rec. RPD RPD Limit LCS Qual LCSD Qual Analyte LCS LCSD Limit Benzene 77 83 70 - 123 7 20 Toluene 84 92 81 - 128 9 20 51 - 97 Gasoline Range Organics (GRO)-C5-C12 62 20 58 6 LCS % Rec Surrogate LCSD % Rec Acceptance Limits 1.2-Dichloroethane-d4 (Surr) 96 89 60 - 140 Toluene-d8 (Surr) 99 105 70 - 130

Client: Engeo, Inc. Job Number: 720-12301-1

Matrix Spike/ Method: 8260B
Matrix Spike Duplicate Recovery Report - Batch: 720-30114 Preparation: 5030B

MS Lab Sample ID: 720-12380-A-5-D MS Analysis Batch: 720-30155 Instrument ID: Varian 3900A

Client Matrix: Solid Prep Batch: 720-30114 Lab File ID: c:\saturnws\data\200712\

Dilution: 1.0 Initial Weight/Volume: 5.52 g

Date Analyzed: 12/27/2007 1521 Final Weight/Volume: 10 mL Date Prepared: 12/27/2007 0930

MSD Lab Sample ID: 720-12380-A-5-E MSD Analysis Batch: 720-30155 Instrument ID: Varian 3900A

Client Matrix: Solid Prep Batch: 720-30114 Lab File ID: c:\saturnws\data\200712\12

Dilution: 1.0 Initial Weight/Volume: 5.46 g

Date Analyzed: 12/27/2007 1544 Final Weight/Volume: 10 mL
Date Prepared: 12/27/2007 0930

% Rec. MSD Limit **RPD RPD Limit** MS Qual MSD Qual Analyte MS Benzene 70 - 123 88 85 2 20 Toluene 94 91 81 - 128 2 20 Gasoline Range Organics (GRO)-C5-C12 58 56 51 - 97 20 MS % Rec MSD % Rec Surrogate Acceptance Limits 1,2-Dichloroethane-d4 (Surr) 93 91 60 - 140 Toluene-d8 (Surr) 107 99 70 - 130

Job Number: 720-12301-1 Client: Engeo, Inc.

Method Blank - Batch: 720-30239 Method: 8015B Preparation: 3550B

Silica Gel Cleanup

Lab Sample ID: MB 720-30239/1-A Analysis Batch: 720-30306 Instrument ID: HP DRO5 Client Matrix: Solid Prep Batch: 720-30239 Lab File ID: N/A

Date Prepared: 12/31/2007 0905

Units: mg/Kg Dilution: 1.0 Initial Weight/Volume: 30.17 g Date Analyzed: 12/31/2007 1826 Final Weight/Volume: 5 mL

Injection Volume:

Column ID: **PRIMARY**

Result Qual RLAnalyte Diesel Range Organics [C10-C28] ND 0.99 Motor Oil Range Organics [C24-C36] ND 50 Surrogate % Rec Acceptance Limits Capric Acid (Surr) 8 Χ 0 - 5 p-Terphenyl 96 41 - 105

Lab Control Spike/ Method: 8015B Lab Control Spike Duplicate Recovery Report - Batch: 720-30239 Preparation: 3550B

Silica Gel Cleanup

LCS Lab Sample ID: LCS 720-30239/2-A Analysis Batch: 720-30306 Instrument ID: HP DRO5

Prep Batch: 720-30239 Client Matrix: Solid Lab File ID: N/A Dilution: 1.0 Units: mg/Kg Initial Weight/Volume:

30.20 g Final Weight/Volume: Date Analyzed: 12/31/2007 1732 5 mL Date Prepared: 12/31/2007 0905 Injection Volume:

Column ID: **PRIMARY**

LCSD Lab Sample ID: LCSD 720-30239/3-A Analysis Batch: 720-30306 Instrument ID: HP DRO5

Prep Batch: 720-30239 Client Matrix: Solid Lab File ID: N/A Dilution: 1.0 Units: mg/Kg Initial Weight/Volume: 30.27 g

Date Analyzed: 12/31/2007 1759 Final Weight/Volume: 5 mL

Injection Volume: Date Prepared: 12/31/2007 0905 Column ID: **PRIMARY**

% Rec.

Analyte LCS **LCSD** Limit **RPD** RPD Limit LCS Qual LCSD Qual Diesel Range Organics [C10-C28] 79 77 50 - 130 3 30 Surrogate LCS % Rec LCSD % Rec Acceptance Limits 87 41 - 105 p-Terphenyl 86

PRIMARY

Job Number: 720-12301-1 Client: Engeo, Inc.

Matrix Spike/ Method: 8015B Matrix Spike Duplicate Recovery Report - Batch: 720-30239 Preparation: 3550B Silica Gel Cleanup

Instrument ID: HP DRO5 MS Lab Sample ID: 720-12365-H-1-C MS Analysis Batch: 720-30306

Prep Batch: 720-30239 Client Matrix: Solid Lab File ID: N/A

Initial Weight/Volume: 30.00 g Dilution: 2.0 Date Analyzed: 12/31/2007 2106 Final Weight/Volume: 5 mL

Date Prepared: 12/31/2007 0905 Injection Volume: Column ID:

Instrument ID: HP DRO5 MSD Lab Sample ID: 720-12365-H-1-D MSD Analysis Batch: 720-30306

Client Matrix: Solid Prep Batch: 720-30239 Lab File ID: N/A

Dilution: 2.0 Initial Weight/Volume: 30.22 g Date Analyzed: 12/31/2007 2133 Final Weight/Volume: 5 mL

Date Prepared: 12/31/2007 0905 Injection Volume:

Column ID: **PRIMARY**

% Rec. MS MSD Limit **RPD** MS Qual MSD Qual Analyte **RPD Limit** Diesel Range Organics [C10-C28] 50 - 130 -190 1 46 30 4 4 Surrogate MS % Rec MSD % Rec Acceptance Limits 53 59 41 - 105 p-Terphenyl

STL San Francisco Chain of Custody 1220 Quarry Lane • Pleasanton CA 94566-4756
Phone: (925) 484-1919 • Fax: (925) 484-1096
Email: sflogin@stl-inc.com
Analysis Reg Reference #: 10869C 1220 Quarry Lane • Pleasanton CA 94566-4756 Date 12/19/07 Page 1 Report To Analysis Request RICHARD GANDOLFO Fuel Tests EPA 8250B: □ Gas □ BTEX □ Five Oxyenates □ 0CA, EDB □ Ethand Company: ENGEO 808 Metals: C Lead C LUFT D RCRA C Other Low Level Metals by EPA 200.8/6020 (ICP-MS): Vojatile Organics GC/MS (VOCs) D NO. ddress: 580 N.W. Ima Are EPA 8081 EPA 8082 hone: 00 SO. W.E.T (STLC) TCLP 00 ill To: Sampled By: 03 and Grease (EPA 1664) R. Gandolfo Phone: Sample ID Dale Time Mat Pres tix erv. MW3026 Project Info. 1) Relinguished by: 2) Relinquished by: 3) Relinquished by: oject Name: # of Containers: Richenbacker Signature Time Signature Time Head Space: 12/19/07 7584-1.001.01 Printed Name Date Printed Name Temp: ENGEO edit Card#: Conforms to record: Company Company 1) Received by: 2) Received by: 3) Received by: 48h Other: port: SO Routine D Level 3 D Level 4 D EDD D Slate Tank Fund EDF Signature Time Signature Time icial Instructions / Comments: Printed Name Date Printed Name Date Company Company 'L SF reports 8015M from C₀-C₂₄ (industry norm). Default for 8015B is C₁₀-C₂₈. Rev 06/04

Login Sample Receipt Check List

Client: Engeo, Inc.

Job Number: 720-12301-1

Login Number: 12301 List Source: TestAmerica San Francisco

Creator: Mullen, Joan List Number: 1

Question	T / F/ NA	Comment
Radioactivity either was not measured or, if measured, is at or below background	N/A	
The cooler's custody seal, if present, is intact.	N/A	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
There are no discrepancies between the sample IDs on the containers and the COC.	True	
Samples are received within Holding Time.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
VOA sample vials do not have headspace or bubble is <6mm (1/4") in diameter.	True	
If necessary, staff have been informed of any short hold time or quick TAT needs	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	



ANALYTICAL REPORT

Job Number: 720-12803-1

Job Description: Rickenbacker

For:
Engeo, Inc.
580 N Wilma Avenue
Suite A
Ripon, CA 95366-9502

Attention: Mr. Richard Gandolfo



Melissa Brewer
Project Manager I
melissa.brewer@testamericainc.com
01/31/2008

Job Narrative 720-J12803-1

Comments

No additional comments.

Receipt All samples were received in good condition within temperature requirements.

GC/MS VOANo analytical or quality issues were noted.

EXECUTIVE SUMMARY - Detections

Client: Engeo, Inc. Job Number: 720-12803-1

Lab Sample ID Analyte	Client Sample ID	Result / Qualifier	Reporting Limit	Units	Method	
720-12803-1 Tetrachloroethene	MW-1	0.80	0.50	ug/L	8260B	
720-12803-2	MW-2		•••	39.1		
Tetrachloroethene		0.95	0.50	ug/L	8260B	

METHOD SUMMARY

Client: Engeo, Inc. Job Number: 720-12803-1

Description	Lab Location	Method	Preparation Method
Matrix: Water			
Volatile Organic Compounds by GC/MS (Low Level)	TAL SF	SW846 8260B	
Purge-and-Trap	TAL SF		SW846 5030B

Lab References:

TAL SF = TestAmerica San Francisco

Method References:

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

SAMPLE SUMMARY

Client: Engeo, Inc. Job Number: 720-12803-1

Lab Sample ID	Client Sample ID	Client Matrix	Date/Time Sampled	Date/Time Received	
720-12803-1	MW-1	Water	01/28/2008 0000	01/28/2008 1214	
720-12803-2	MW-2	Water	01/28/2008 0000	01/28/2008 1214	
720-12803-3	MW-3	Water	01/28/2008 0000	01/28/2008 1214	

Client: Engeo, Inc. Job Number: 720-12803-1

Client Sample ID: MW-1

 Lab Sample ID:
 720-12803-1
 Date Sampled:
 01/28/2008 0000

 Client Matrix:
 Water
 Date Received:
 01/28/2008 1214

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method: 8260B Analysis Batch: 720-31331 Instrument ID: Varian 3900G

Preparation: 5030B Lab File ID: c:\saturnws\data\200801\01

Dilution: 1.0 Initial Weight/Volume: 40 mL Date Analyzed: 01/30/2008 1922 Final Weight/Volume: 40 mL

Analyte	Result (ug/L)	Qualifier	RL
Methyl tert-butyl ether	ND		5.0
Acetone	ND		50
Benzene	ND		0.50
Dichlorobromomethane	ND		0.50
Bromobenzene	ND		1.0
Chlorobromomethane	ND		1.0
Bromoform	ND		1.0
Bromomethane	ND		1.0
2-Butanone (MEK)	ND		50
n-Butylbenzene	ND		1.0
sec-Butylbenzene	ND		1.0
tert-Butylbenzene	ND		1.0
Carbon disulfide	ND		5.0
Carbon tetrachloride	ND		0.50
Chlorobenzene	ND		0.50
Chloroethane	ND		1.0
Chloroform	ND		1.0
Chloromethane	ND		1.0
2-Chlorotoluene	ND		0.50
4-Chlorotoluene	ND		0.50
Chlorodibromomethane	ND		0.50
1,2-Dichlorobenzene	ND		0.50
1,3-Dichlorobenzene	ND		0.50
1,4-Dichlorobenzene	ND		0.50
1,3-Dichloropropane	ND		1.0
1,1-Dichloropropene	ND		0.50
1,2-Dibromo-3-Chloropropane	ND		1.0
Ethylene Dibromide	ND		0.50
Dibromomethane	ND		0.50
Dichlorodifluoromethane	ND		0.50
1,1-Dichloroethane	ND		0.50
1,2-Dichloroethane	ND		0.50
1,1-Dichloroethene	ND		0.50
cis-1,2-Dichloroethene	ND		0.50
trans-1,2-Dichloroethene	ND		0.50
1,2-Dichloropropane	ND		0.50
cis-1,3-Dichloropropene	ND		0.50
trans-1,3-Dichloropropene	ND		0.50
Ethylbenzene	ND		0.50
Hexachlorobutadiene	ND		1.0
2-Hexanone	ND		50
Isopropylbenzene	ND		0.50
4-Isopropyltoluene	ND		1.0
Methylene Chloride	ND		5.0

Client: Engeo, Inc. Job Number: 720-12803-1

Client Sample ID: MW-1

 Lab Sample ID:
 720-12803-1
 Date Sampled:
 01/28/2008 0000

 Client Matrix:
 Water
 Date Received:
 01/28/2008 1214

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method: 8260B Analysis Batch: 720-31331 Instrument ID: Varian 3900G

Preparation: 5030B Lab File ID: c:\saturnws\data\200801\01

Dilution: 1.0 Initial Weight/Volume: 40 mL Date Analyzed: 01/30/2008 1922 Final Weight/Volume: 40 mL

Analyte	Result (ug/L)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		50
Naphthalene	ND		1.0
N-Propylbenzene	ND		1.0
Styrene	ND		0.50
1,1,1,2-Tetrachloroethane	ND		0.50
1,1,2,2-Tetrachloroethane	ND		0.50
Tetrachloroethene	0.80		0.50
Toluene	ND		0.50
1,2,3-Trichlorobenzene	ND		1.0
1,2,4-Trichlorobenzene	ND		1.0
1,1,1-Trichloroethane	ND		0.50
1,1,2-Trichloroethane	ND		0.50
Trichloroethene	ND		0.50
Trichlorofluoromethane	ND		1.0
1,2,3-Trichloropropane	ND		0.50
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		0.50
1,2,4-Trimethylbenzene	ND		0.50
1,3,5-Trimethylbenzene	ND		0.50
Vinyl acetate	ND		50
Vinyl chloride	ND		0.50
Xylenes, Total	ND		1.0
2,2-Dichloropropane	ND		0.50
Surrogate	%Rec		Acceptance Limits
4-Bromofluorobenzene	108		71 - 139
1,2-Dichloroethane-d4 (Surr)	96		62 - 118
Toluene-d8 (Surr)	101		73 - 117

Client: Engeo, Inc. Job Number: 720-12803-1

Client Sample ID: MW-2

 Lab Sample ID:
 720-12803-2
 Date Sampled:
 01/28/2008 0000

 Client Matrix:
 Water
 Date Received:
 01/28/2008 1214

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method: 8260B Analysis Batch: 720-31331 Instrument ID: Varian 3900G

Preparation: 5030B Lab File ID: c:\saturnws\data\200801\01

Dilution: 1.0 Initial Weight/Volume: 40 mL Date Analyzed: 01/30/2008 1955 Final Weight/Volume: 40 mL

Analyte	Result (ug/L)	Qualifier	RL
Methyl tert-butyl ether	ND		5.0
Acetone	ND		50
Benzene	ND		0.50
Dichlorobromomethane	ND		0.50
Bromobenzene	ND		1.0
Chlorobromomethane	ND		1.0
Bromoform	ND		1.0
Bromomethane	ND		1.0
2-Butanone (MEK)	ND		50
n-Butylbenzene	ND		1.0
sec-Butylbenzene	ND		1.0
tert-Butylbenzene	ND		1.0
Carbon disulfide	ND		5.0
Carbon tetrachloride	ND		0.50
Chlorobenzene	ND		0.50
Chloroethane	ND		1.0
Chloroform	ND		1.0
Chloromethane	ND		1.0
2-Chlorotoluene	ND		0.50
4-Chlorotoluene	ND		0.50
Chlorodibromomethane	ND		0.50
1,2-Dichlorobenzene	ND		0.50
1,3-Dichlorobenzene	ND		0.50
1,4-Dichlorobenzene	ND		0.50
1,3-Dichloropropane	ND		1.0
1,1-Dichloropropene	ND		0.50
1,2-Dibromo-3-Chloropropane	ND		1.0
Ethylene Dibromide	ND		0.50
Dibromomethane	ND		0.50
Dichlorodifluoromethane	ND		0.50
1,1-Dichloroethane	ND		0.50
1,2-Dichloroethane	ND		0.50
1,1-Dichloroethene	ND		0.50
cis-1,2-Dichloroethene	ND		0.50
trans-1,2-Dichloroethene	ND		0.50
1,2-Dichloropropane	ND		0.50
cis-1,3-Dichloropropene	ND		0.50
trans-1,3-Dichloropropene	ND		0.50
Ethylbenzene	ND		0.50
Hexachlorobutadiene	ND		1.0
2-Hexanone	ND		50
Isopropylbenzene	ND		0.50
4-Isopropyltoluene	ND		1.0
Methylene Chloride	ND		5.0

Client: Engeo, Inc. Job Number: 720-12803-1

Client Sample ID: MW-2

 Lab Sample ID:
 720-12803-2
 Date Sampled:
 01/28/2008 0000

 Client Matrix:
 Water
 Date Received:
 01/28/2008 1214

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method: 8260B Analysis Batch: 720-31331 Instrument ID: Varian 3900G

Preparation: 5030B Lab File ID: c:\saturnws\data\200801\01

Dilution: 1.0 Initial Weight/Volume: 40 mL Date Analyzed: 01/30/2008 1955 Final Weight/Volume: 40 mL

Analyte	Result (ug/L)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		50
Naphthalene	ND		1.0
N-Propylbenzene	ND		1.0
Styrene	ND		0.50
1,1,1,2-Tetrachloroethane	ND		0.50
1,1,2,2-Tetrachloroethane	ND		0.50
Tetrachloroethene	0.95		0.50
Toluene	ND		0.50
1,2,3-Trichlorobenzene	ND		1.0
1,2,4-Trichlorobenzene	ND		1.0
1,1,1-Trichloroethane	ND		0.50
1,1,2-Trichloroethane	ND		0.50
Trichloroethene	ND		0.50
Trichlorofluoromethane	ND		1.0
1,2,3-Trichloropropane	ND		0.50
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		0.50
1,2,4-Trimethylbenzene	ND		0.50
1,3,5-Trimethylbenzene	ND		0.50
Vinyl acetate	ND		50
Vinyl chloride	ND		0.50
Xylenes, Total	ND		1.0
2,2-Dichloropropane	ND		0.50
Surrogate	%Rec		Acceptance Limits
4-Bromofluorobenzene	109		71 - 139
1,2-Dichloroethane-d4 (Surr)	94		62 - 118
Toluene-d8 (Surr)	100		73 - 117

Client: Engeo, Inc. Job Number: 720-12803-1

Client Sample ID: MW-3

 Lab Sample ID:
 720-12803-3
 Date Sampled:
 01/28/2008 0000

 Client Matrix:
 Water
 Date Received:
 01/28/2008 1214

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method: 8260B Analysis Batch: 720-31331 Instrument ID: Varian 3900G

Preparation: 5030B Lab File ID: c:\saturnws\data\200801\01

Dilution: 1.0 Initial Weight/Volume: 40 mL Date Analyzed: 01/30/2008 2029 Final Weight/Volume: 40 mL

Analyte	Result (ug/L)	Qualifier	RL
Methyl tert-butyl ether	ND		5.0
Acetone	ND		50
Benzene	ND		0.50
Dichlorobromomethane	ND		0.50
Bromobenzene	ND		1.0
Chlorobromomethane	ND		1.0
Bromoform	ND		1.0
Bromomethane	ND		1.0
2-Butanone (MEK)	ND		50
n-Butylbenzene	ND		1.0
sec-Butylbenzene	ND		1.0
tert-Butylbenzene	ND		1.0
Carbon disulfide	ND		5.0
Carbon tetrachloride	ND		0.50
Chlorobenzene	ND		0.50
Chloroethane	ND		1.0
Chloroform	ND		1.0
Chloromethane	ND		1.0
2-Chlorotoluene	ND		0.50
4-Chlorotoluene	ND		0.50
Chlorodibromomethane	ND		0.50
1,2-Dichlorobenzene	ND		0.50
1,3-Dichlorobenzene	ND		0.50
1,4-Dichlorobenzene	ND		0.50
1,3-Dichloropropane	ND		1.0
1,1-Dichloropropene	ND		0.50
1,2-Dibromo-3-Chloropropane	ND		1.0
Ethylene Dibromide	ND		0.50
Dibromomethane	ND		0.50
Dichlorodifluoromethane	ND		0.50
1,1-Dichloroethane	ND		0.50
1,2-Dichloroethane	ND		0.50
1,1-Dichloroethene	ND		0.50
cis-1,2-Dichloroethene	ND		0.50
trans-1,2-Dichloroethene	ND		0.50
1,2-Dichloropropane	ND		0.50
cis-1,3-Dichloropropene	ND		0.50
trans-1,3-Dichloropropene	ND		0.50
Ethylbenzene	ND		0.50
Hexachlorobutadiene	ND		1.0
2-Hexanone	ND		50
Isopropylbenzene	ND		0.50
4-Isopropyltoluene	ND		1.0
			1.0

Client: Engeo, Inc. Job Number: 720-12803-1

Client Sample ID: MW-3

 Lab Sample ID:
 720-12803-3
 Date Sampled:
 01/28/2008 0000

 Client Matrix:
 Water
 Date Received:
 01/28/2008 1214

8260B Volatile Organic Compounds by GC/MS (Low Level)

Method: 8260B Analysis Batch: 720-31331 Instrument ID: Varian 3900G

Preparation: 5030B Lab File ID: c:\saturnws\data\200801\01

Dilution: 1.0 Initial Weight/Volume: 40 mL Date Analyzed: 01/30/2008 2029 Final Weight/Volume: 40 mL

Analyte	Result (ug/L)	Qualifier	RL
4-Methyl-2-pentanone (MIBK)	ND		50
Naphthalene	ND		1.0
N-Propylbenzene	ND		1.0
Styrene	ND		0.50
1,1,1,2-Tetrachloroethane	ND		0.50
1,1,2,2-Tetrachloroethane	ND		0.50
Tetrachloroethene	ND		0.50
Toluene	ND		0.50
1,2,3-Trichlorobenzene	ND		1.0
1,2,4-Trichlorobenzene	ND		1.0
1,1,1-Trichloroethane	ND		0.50
1,1,2-Trichloroethane	ND		0.50
Trichloroethene	ND		0.50
Trichlorofluoromethane	ND		1.0
1,2,3-Trichloropropane	ND		0.50
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		0.50
1,2,4-Trimethylbenzene	ND		0.50
1,3,5-Trimethylbenzene	ND		0.50
Vinyl acetate	ND		50
Vinyl chloride	ND		0.50
Xylenes, Total	ND		1.0
2,2-Dichloropropane	ND		0.50
Surrogate	%Rec		Acceptance Limits
4-Bromofluorobenzene	103		71 - 139
1,2-Dichloroethane-d4 (Surr)	91		62 - 118
Toluene-d8 (Surr)	99		73 - 117

DATA REPORTING QUALIFIERS

Lab Section Qualifier Description

Client: Engeo, Inc. Job Number: 720-12803-1

QC Association Summary

		Report			
Lab Sample ID	Client Sample ID	Basis	Client Matrix	Method	Prep Batch
GC/MS VOA					
Analysis Batch:720-31	331				
LCS 720-31331/2	Lab Control Spike	T	Water	8260B	
LCSD 720-31331/1	Lab Control Spike Duplicate	Т	Water	8260B	
MB 720-31331/3	Method Blank	T	Water	8260B	
720-12761-B-1 MS	Matrix Spike	Т	Water	8260B	
720-12761-C-1 MSD	Matrix Spike Duplicate	Т	Water	8260B	
720-12803-1	MW-1	Т	Water	8260B	
720-12803-2	MW-2	Т	Water	8260B	
720-12803-3	MW-3	Т	Water	8260B	

Report Basis

T = Total

Client: Engeo, Inc. Job Number: 720-12803-1

Method Blank - Batch: 720-31331 Method: 8260B Preparation: 5030B

Lab Sample ID: MB 720-31331/3 Analysis Batch: 720-31331 Instrument ID: Varian 3900G

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\saturnws\data\200801\01

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 40 mL Date Analyzed: 01/30/2008 1059 Final Weight/Volume: 40 mL

Date Prepared: 01/30/2008 1059

Analyte	Result	Qual	RL
Methyl tert-butyl ether	ND		5.0
Acetone	ND		50
Benzene	ND		0.50
Dichlorobromomethane	ND		0.50
Bromobenzene	ND		1.0
Chlorobromomethane	ND		1.0
Bromoform	ND		1.0
Bromomethane	ND		1.0
2-Butanone (MEK)	ND		50
n-Butylbenzene	ND		1.0
sec-Butylbenzene	ND		1.0
tert-Butylbenzene	ND		1.0
Carbon disulfide	ND		5.0
Carbon tetrachloride	ND		0.50
Chlorobenzene	ND		0.50
Chloroethane	ND		1.0
Chloroform	ND		1.0
Chloromethane	ND		1.0
2-Chlorotoluene	ND		0.50
4-Chlorotoluene	ND		0.50
Chlorodibromomethane	ND		0.50
1,2-Dichlorobenzene	ND		0.50
1,3-Dichlorobenzene	ND		0.50
1,4-Dichlorobenzene	ND		0.50
1,3-Dichloropropane	ND		1.0
1,1-Dichloropropene	ND		0.50
1,2-Dibromo-3-Chloropropane	ND		1.0
Ethylene Dibromide	ND		0.50
Dibromomethane	ND		0.50
Dichlorodifluoromethane	ND		0.50
1,1-Dichloroethane	ND		0.50
1,2-Dichloroethane	ND		0.50
1,1-Dichloroethene	ND		0.50
cis-1,2-Dichloroethene	ND		0.50
trans-1,2-Dichloroethene	ND		0.50
1,2-Dichloropropane	ND		0.50
cis-1,3-Dichloropropene	ND		0.50
trans-1,3-Dichloropropene	ND		0.50
Ethylbenzene	ND		0.50
Hexachlorobutadiene	ND		1.0
2-Hexanone	ND		50

Job Number: 720-12803-1 Client: Engeo, Inc.

Method Blank - Batch: 720-31331 Method: 8260B Preparation: 5030B

Lab Sample ID: MB 720-31331/3 Analysis Batch: 720-31331 Instrument ID: Varian 3900G

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\saturnws\data\200801\01

Units: ug/L Initial Weight/Volume: 40 mL Dilution: 1.0

Date Analyzed: 01/30/2008 1059 Final Weight/Volume: 40 mL Date Prepared: 01/30/2008 1059

Analyte	Result	Qual	RL
Isopropylbenzene	ND		0.50
4-Isopropyltoluene	ND		1.0
Methylene Chloride	ND		5.0
4-Methyl-2-pentanone (MIBK)	ND		50
Naphthalene	ND		1.0
N-Propylbenzene	ND		1.0
Styrene	ND		0.50
1,1,1,2-Tetrachloroethane	ND		0.50
1,1,2,2-Tetrachloroethane	ND		0.50
Tetrachloroethene	ND		0.50
Toluene	ND		0.50
1,2,3-Trichlorobenzene	ND		1.0
1,2,4-Trichlorobenzene	ND		1.0
1,1,1-Trichloroethane	ND		0.50
1,1,2-Trichloroethane	ND		0.50
Trichloroethene	ND		0.50
Trichlorofluoromethane	ND		1.0
1,2,3-Trichloropropane	ND		0.50
1,1,2-Trichloro-1,2,2-trifluoroethane	ND		0.50
1,2,4-Trimethylbenzene	ND		0.50
1,3,5-Trimethylbenzene	ND		0.50
Vinyl acetate	ND		50
Vinyl chloride	ND		0.50
Xylenes, Total	ND		1.0
2,2-Dichloropropane	ND		0.50
Surrogate	% Rec	Acceptance Limits	
4-Bromofluorobenzene	105	71 - 139	
1,2-Dichloroethane-d4 (Surr)	93	62 - 118	
Toluene-d8 (Surr)	100	73 - 117	

Client: Engeo, Inc. Job Number: 720-12803-1

Lab Control Spike/ Method: 8260B
Lab Control Spike Duplicate Recovery Report - Batch: 720-31331 Preparation: 5030B

LCS Lab Sample ID: LCS 720-31331/2 Analysis Batch: 720-31331 Instrument ID: Varian 3900G

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\saturnws\data\200801\0'

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 40 mL

Date Analyzed: 01/30/2008 0952 Final Weight/Volume: 40 mL Date Prepared: 01/30/2008 0952

LCSD Lab Sample ID: LCSD 720-31331/1 Analysis Batch: 720-31331 Instrument ID: Varian 3900G

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\saturnws\data\200801\013

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 40 mL

Date Analyzed: 01/30/2008 1026 Final Weight/Volume: 40 mL Date Prepared: 01/30/2008 1026

		% Rec.			
Analyte	LCS	LCSD	Limit	RPD	RPD Limit LCS Qual LCSD Qual
Benzene	82	93	69 - 129	12	20
Chlorobenzene	98	106	61 - 121	8	20
1,1-Dichloroethene	86	102	65 - 125	18	20
Toluene	89	100	70 - 130	11	20
Trichloroethene	84	94	74 - 134	11	20
Surrogate		LCS % Rec	LCSD %	Rec	Acceptance Limits
4-Bromofluorobenzene		96	97		71 - 139
1,2-Dichloroethane-d4 (Surr)		84	87		62 - 118
Toluene-d8 (Surr)		93	94		73 - 117

Client: Engeo, Inc. Job Number: 720-12803-1

Matrix Spike/ Method: 8260B
Matrix Spike Duplicate Recovery Report - Batch: 720-31331 Preparation: 5030B

Date Prepared:

Date Prepared:

01/30/2008 1313

01/30/2008 1347

MS Lab Sample ID: 720-12761-B-1 MS Analysis Batch: 720-31331 Instrument ID: Varian 3900G

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\saturnws\data\200801\(

Dilution: 1.0 Initial Weight/Volume: 40 mL Date Analyzed: 01/30/2008 1313 Final Weight/Volume: 40 mL

MSD Lab Sample ID: 720-12761-C-1 MSD Analysis Batch: 720-31331 Instrument ID: Varian 3900G

Client Matrix: Water Prep Batch: N/A Lab File ID: c:\saturnws\data\200801\01

Dilution: 1.0 Initial Weight/Volume: 40 mL

Date Analyzed: 01/30/2008 1347 Final Weight/Volume: 40 mL

% Rec. MS MSD **RPD** MS Qual MSD Qual Analyte Limit **RPD Limit** Benzene 89 69 - 129 2 20 91 Chlorobenzene 104 106 61 - 121 2 20 1,1-Dichloroethene 91 89 65 - 125 1 20 2 Toluene 101 99 70 - 130 20 74 - 134 20 Trichloroethene 99 98 MS % Rec MSD % Rec Surrogate Acceptance Limits 103 104 4-Bromofluorobenzene 71 - 139 62 - 118 93 1,2-Dichloroethane-d4 (Surr) 94 73 - 117 Toluene-d8 (Surr) 103 99

Login Sample Receipt Check List

Client: Engeo, Inc.

Job Number: 720-12803-1

Login Number: 12803 List Source: TestAmerica San Francisco

Creator: Bullock, Tracy List Number: 1

Question	T / F/ NA	Comment
Radioactivity either was not measured or, if measured, is at or below background	N/A	
The cooler's custody seal, if present, is intact.	N/A	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
There are no discrepancies between the sample IDs on the containers and the COC.	True	
Samples are received within Holding Time.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	False	NO TIMES
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
VOA sample vials do not have headspace or bubble is <6mm (1/4") in diameter.	True	
If necessary, staff have been informed of any short hold time or quick TAT needs	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	



APPENDIX C

TORRENT LABORATORY, INC.

Laboratory Analytical Report



December 27, 2007

Kelly Krohn Engeo 2010 Crow Canyon Place,#250 San Ramon, CA 94583

TEL: (925) 570-7602 FAX (925) 866-0199

RE: 7584.1.001.01

Dear Kelly Krohn:

Order No.: 0712091

Torrent Laboratory, Inc. received 8 samples on 12/17/2007 for the analyses presented in the following report.

All data for associated QC met EPA or laboratory specification(s) except where noted in the case narrative.

Torrent Laboratory, Inc, is certified by the State of California, ELAP #1991. If you have any questions regarding these tests results, please feel free to contact the Project Management Team at (408)263-5258;ext: 204.

Sincerely,

Patti Sandrock

OA Officer



483 Sinclair Frontage Road * Milpitas, CA * Phone: (408) 2635258 * Fax: (408) 263-8293 Visit us ar www.torrentlab.com email: analysis@torrentlab.com

Report Prepaired For: Kelly Krohn

Engeo

Date Received: Date Reported: 12/17/2007 12/27/2007

Summary Report

SG-11	Toxic Organics in Air	by EPA TO-15		Lab II	D: 0712091-001A
<u>Parameter</u>	Preped	<u>Analyzed</u>	Result	<u>RL</u>	<u>Unit</u>
2-Butanone (MEK)	12/26/2007	12/26/2007	77	0.70	μg/m³
Acetone	12/26/2007	12/26/2007	180	0.90	μg/m³
Benzene	12/26/2007	12/26/2007	3.5	1.4	µg/m³
m,p-Xylene	12/26/2007	12/26/2007	38	0.78	µg/m³
o-xylene	12/26/2007	12/26/2007	11	0.98	µg/m³
Tetrachloroethene	12/26/2007	12/26/2007	64	2.0	µg/m³
Toluene	12/26/2007	12/26/2007	25	0.83	µg/m³
SG-12	Toxic Organics in Air	by EPA TO-15		Lab II	D: 0712091-002A
<u>Parameter</u>	<u>Preped</u>	<u>Analyzed</u>	Result	<u>RL</u>	<u>Unit</u>
2-Butanone (MEK)	12/26/2007	12/26/2007	31	0.69	μg/m³
Acetone	12/26/2007	12/26/2007	70	0.89	µg/m³
Benzene	12/26/2007	12/26/2007	2.5	1.4	µg/m³
m,p-Xylene	12/26/2007	12/26/2007	25	0.77	µg/m³
o-xylene	12/26/2007	12/26/2007	6.4	0.97	µg/m³
Tetrachloroethene	12/26/2007	12/26/2007	10	2.0	µg/m³
Toluene	12/26/2007	12/26/2007	16	0.82	µg/m³
SG-12 (Duplicate)	Toxic Organics in Air	by EPA TO-15		Lab II	D: 0712091-003A
<u>Parameter</u>	<u>Preped</u>	<u>Analyzed</u>	Result	<u>RL</u>	<u>Unit</u>
2-Butanone (MEK)	12/26/2007	12/26/2007	28	0.65	µg/m³
Acetone	12/26/2007	12/26/2007	62	0.85	µg/m³
Benzene	12/26/2007	12/26/2007	2.2	1.3	µg/m³
m,p-Xylene	12/26/2007	12/26/2007	21	0.73	μg/m³
o-xylene	12/26/2007	12/26/2007	5.3	0.92	µg/m³
Tetrachloroethene	12/26/2007	12/26/2007	8.7	1.9	µg/m³
Toluene	12/26/2007	12/26/2007	14	0.78	µg/m³
SG-13	Toxic Organics in Air	by EPA TO-15		Lab II	D: 0712091-004A
<u>Parameter</u>	<u>Preped</u>	<u>Analyzed</u>	Result	<u>RL</u>	<u>Unit</u>
2-Butanone (MEK)	12/26/2007	12/26/2007	42	0.44	µg/m³
Acetone	12/26/2007	12/26/2007	110	0.57	µg/m³



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Report Prepaired For: Kelly Krohn

Engeo

Date Received: Date Reported: 12/17/2007 12/27/2007

Summary Report

		· ·			
SG-13	Toxic Organics in Air	by EPA TO-15		Lab	ID: 0712091-004A
<u>Parameter</u>	Preped	<u>Analyzed</u>	Result	<u>RL</u>	<u>Unit</u>
Benzene	12/26/2007	12/26/2007	3.1	0.89	μg/m³
Hexane	12/26/2007	12/26/2007	27	1.8	μg/m³
m,p-Xylene	12/26/2007	12/26/2007	34	0.49	μg/m³
o-xylene	12/26/2007	δ/2007 12/26/2007 9.2 0.62 μg		μg/m³	
Toluene	12/26/2007	007 12/26/2007 48 0.53		μg/m³	
SG-14	Toxic Organics in Air	by EPA TO-15		Lab	ID: 0712091-005A
<u>Parameter</u>	<u>Preped</u>	<u>Analyzed</u>	Result	<u>RL</u>	<u>Unit</u>
1,2,4-Trimethylbenzene	12/26/2007	12/26/2007	2.5	1.4	μg/m³
2-Hexanone	12/26/2007	12/26/2007	1.7	1.4	μg/m³
4-Ethyl Toluene	12/26/2007	12/26/2007	1.9	1.2	μg/m³
Ethyl Benzene	12/26/2007	12/26/2007	1.7	0.49	μg/m³
m,p-Xylene	12/26/2007	12/26/2007	8.0	0.77	μg/m³
Toluene	12/26/2007	12/26/2007	3.3	0.83	µg/m³
SG-15	Toxic Organics in Air	by EPA TO-15		Lab	ID: 0712091-006A
<u>Parameter</u>	Preped	<u>Analyzed</u>	Result	<u>RL</u>	<u>Unit</u>
2-Butanone (MEK)	12/26/2007	12/26/2007	29	0.65	μg/m³
Acetone	12/26/2007	12/26/2007	200	0.84	μg/m³
Benzene	12/26/2007	12/26/2007	4.0	1.3	μg/m³
Hexane	12/26/2007	12/26/2007	59	2.6	μg/m³
m,p-Xylene	12/26/2007	12/26/2007	39	0.72	μg/m³
o-xylene	12/26/2007	12/26/2007	11	0.91	μg/m³
Toluene	12/26/2007	12/26/2007	68	0.78	μg/m³
SG-16	Toxic Organics in Air	by EPA TO-15		Lab	ID: 0712091-007A
<u>Parameter</u>	Preped	<u>Analyzed</u>	Result	<u>RL</u>	<u>Unit</u>
2-Butanone (MEK)	12/26/2007	12/26/2007	110	0.66	μg/m³
Acetone	12/26/2007	12/26/2007	180	0.86	μg/m³
Benzene	12/26/2007	12/26/2007	6.6	1.3	μg/m³
cis-1,2-dichloroethene	12/26/2007	12/26/2007	7.9	0.83	µg/m³
Ethyl Benzene	12/26/2007	12/26/2007	8.2	0.47	µg/m³



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Report Prepaired For: Kelly Krohn

Engeo

Date Received: Date Reported: 12/17/2007 12/27/2007

Summary Report

SG-16	Toxic Organics in Air	by EPA TO-15		Lab	ID: 0712091-007A
<u>Parameter</u>	<u>Preped</u>	<u>Analyzed</u>	Result	<u>RL</u>	<u>Unit</u>
m,p-Xylene	12/26/2007	12/26/2007	46	0.74	μg/m³
o-xylene	12/26/2007	12/26/2007	13	0.93	μg/m³
Tetrachloroethene	12/26/2007	12/26/2007	15	1.9	μg/m³
Toluene	12/26/2007	12/26/2007	30	0.79	μg/m³
trans-1,2-Dichloroethene	12/26/2007	12/26/2007	8.2	0.83	μg/m³
Trichloroethene	12/26/2007	12/26/2007	22	0.79	μg/m³
SG-10	Toxic Organics in Air by EPA TO-15 Lab ID:				ID: 0712091-008A
	TOXIO OI gaillos III All	by =: /			
<u>Parameter</u>	Preped	Analyzed	Result		<u>Unit</u>
	•	•	Result 58	<u>RL</u>	
<u>Parameter</u>	<u>Preped</u>	Analyzed		<u>RL</u> 0.72	<u>Unit</u>
Parameter 2-Butanone (MEK)	<u>Preped</u> 12/26/2007	<u>Analyzed</u> 12/26/2007	58	<u>RL</u> 0.72	<u>Unit</u> μg/m³
Parameter 2-Butanone (MEK) 2-Hexanone	Preped 12/26/2007 12/26/2007	Analyzed 12/26/2007 12/26/2007	58 5.3	<u>RL</u> 0.72 1.4	<u>Unit</u> μg/m³ μg/m³
Parameter 2-Butanone (MEK) 2-Hexanone Acetone	Preped 12/26/2007 12/26/2007 12/26/2007	Analyzed 12/26/2007 12/26/2007 12/26/2007	58 5.3 100	RL 0.72 1.4 0.93	<u>Unit</u> μg/m³ μg/m³ μg/m³
Parameter 2-Butanone (MEK) 2-Hexanone Acetone Benzene	Preped 12/26/2007 12/26/2007 12/26/2007 12/26/2007	Analyzed 12/26/2007 12/26/2007 12/26/2007	58 5.3 100 2.8	RL 0.72 1.4 0.93 1.5	Unit µg/m³ µg/m³ µg/m³
Parameter 2-Butanone (MEK) 2-Hexanone Acetone Benzene Hexane	Preped 12/26/2007 12/26/2007 12/26/2007 12/26/2007 12/26/2007	Analyzed 12/26/2007 12/26/2007 12/26/2007 12/26/2007 12/26/2007	58 5.3 100 2.8 17	RL 0.72 1.4 0.93 1.5 2.9	Unit µg/m³ µg/m³ µg/m³ µg/m³



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Date Received: 12/17/2007

Lab Sample ID: 0712091-001

Date Prepared:

Report prepared for: Kelly Krohn

Engeo Date Reported: 12/27/2007

Client Sample ID: SG-11

Sample Location: 224 Rickenbacker Circle

Sample Matrix: AIR

Date/Time Sampled 12/17/2007 11:22:00 AM

1,1,1,2-Tetrachloroethane TO-15 12/26/2007 0.687 1.58 1.1 ND µg/m³ R14916 1,1,1-Trichloroethane TO-15 12/26/2007 1.0305 1.58 1.6 ND µg/m³ R14916 1,1,2-Trichloroethane TO-15 12/26/2007 1.0305 1.58 1.6 ND µg/m³ R14916 1,1,2-Trichloroethane TO-15 12/26/2007 1.0305 1.58 1.6 ND µg/m³ R14916 1,1,2-Trichloroethane TO-15 12/26/2007 1.0305 1.58 1.6 ND µg/m³ R14916 1,1,2-Trichloroethane TO-15 12/26/2007 0.6885 1.58 1.1 ND µg/m³ R14916 1,2-4-Trichloroethane TO-15 12/26/2007 0.4984 1.58 0.79 ND µg/m³ R14916 1,2-4-Trimethylbenzene TO-15 12/26/2007 0.8856 1.58 1.7 ND µg/m³ R14916 1,2-4-Trimethylbenzene TO-15 12/26/2007 0.8986 1.58 1.7 ND µg/m³ R14916 1,2-Dichloroethane(Ethylene tiltoromide) 1,2-Dichloroethane TO-15 12/26/2007 0.648 1.58 0.95 ND µg/m³ R14916 1,2-Dichloroethane TO-15 12/26/2007 0.648 1.58 0.95 ND µg/m³ R14916 1,2-Dichloroethane TO-15 12/26/2007 0.648 1.58 0.95 ND µg/m³ R14916 1,2-Dichloroethane TO-15 12/26/2007 0.648 1.58 0.95 ND µg/m³ R14916 1,2-Dichloroethane TO-15 12/26/2007 0.648 1.58 0.05 ND µg/m³ R14916 1,2-Dichloroethane TO-15 12/26/2007 0.648 1.58 0.05 ND µg/m³ R14916 1,3-Dichloroethane(F114) 1.3-Butadiene TO-15 12/26/2007 0.5867 1.58 0.94 ND µg/m³ R14916 1,4-Dichloroethane TO-15 12/26/2007 0.5867 1.58 0.94 ND µg/m³ R14916 1,4-Dichloroethane TO-15 12/26/2007 0.6811 1.58 0.90 ND µg/m³ R14916 1,4-Dichloroethane TO-15 12/26/2007 0.6811 1.58 0.90 ND µg/m³ R14916 1,4-Dichloroethane TO-15 12/26/2007 0.6811 1.58 0.90 ND µg/m³ R14916 1,4-Dichloroethane TO-15 12/26/2007 0.6811 1.58 0.90 ND µg/m³ R14916 1,4-Dichloroethane TO-15 12/26/2007 0.6811 1.58 0.90 ND µg/m³ R14916 1,4-Dichloroethane TO-15 12/26/2007 0.6811 1.58 0.90 ND µg/m³ R14916 1,4-Dichloroethane TO-15 12/26/2007 0.6811 1.58 0.90 ND µg/m³ R14916 1,4-Dichloroethane TO-15 12/26/2007 0.8081 1.80 ND µg/m³ R14916 1,4-Dichloroethane TO-15 12/26/2007 0.8081 1.80 ND µg/m³ R14916 1,4-Dichloroethane TO-15 12/26/2007 0.8081 1.80 ND µg/m³ R14916 1.80 ND µg/m³	Parameters	Analysis Method	Date Analyzed	RL	Dilution Factor	MRL	Result	Units	Analytical Batch
1,1,1-Trichloroethane	1,1 - Dichloroethene	TO-15	12/26/2007	0.794	1.58	1.3	ND	μg/m³	R14916
1.1,1-Trichloroethane	1,1,1,2-Tetrachloroethane	TO-15	12/26/2007	0.687	1.58	1.1	ND	μg/m³	R14916
1,1,2,2-Tetrachloroethane 1,1,2-Tichloroethane 1,1,2-Tichloroethane 1,1,2-Tichloroethane 1,1,2-Tichloroethane 1,1,2-Tichloroethane 1,1,2-Tichloroethane 1,1,2-Tichloroethane 1,1,2-Tichloroethane 1,2,4-Tichloroethane 1,2,4-Tichloroethane 1,2,4-Tichloroethane 1,2,4-Tichloroethane 1,2,4-Tichloroethane 1,2,4-Tichloroethane 1,2,4-Tichloroethane 1,2,4-Tichloroethane(Ethylene 1,2,4-Tichloroethane(Ethylene) 1,2-Dichloroethane(Ethylene) 1,3-S-Timethylbenzene 1,0-15 1,2-Dichloroethane(Ethylene) 1,3-S-Timethylbenzene 1,0-15 1,2-Dichloroethane(Ethylene) 1,3-Dichloroethane(Ethylene) 1,3-Dichloroethane(Ethylen	1,1,1-Trichloroethane	TO-15	12/26/2007	0.819	1.58	1.3	ND		R14916
1,1,2-Trichloroethane	1,1,2,2-Tetrachloroethane	TO-15	12/26/2007						R14916
1,1-Dichloroethane	1,1,2-Trichloroethane	TO-15	12/26/2007	1.0374	1.58	1.6	ND		R14916
1,2,4-Trimethylbenzene TO-15 12/26/2007 0.8856 1.58 1.4 ND µg/m³ R14916 (1,2-Dibromoethane(Ethylene dibromide)	1,1-Dichloroethane	TO-15	12/26/2007	0.6885	1.58	1.1	ND	μg/m³	R14916
1.2.4-Trimethylbenzene TO-15 12/26/2007 0.8856 1.58 1.4 ND µg/m³ R14916 12Dibromoethane(Ethylene dibromide)	1,2,4-Trichlorobenzene	TO-15	12/26/2007	0.4984	1.58	0.79	ND	μg/m³	R14916
dibromide) 1,2-Dichlorobenzene TO-15 12/26/2007 0,601 1,58 0,95 ND µg/m³ R14916 1,2-Dichlorobenzene TO-15 12/26/2007 0,648 1,58 1,0 ND µg/m³ R14916 1,2-Dichloropropane TO-15 12/26/2007 1,0164 1,58 1,6 ND µg/m³ R14916 1,2-Dichloropropane TO-15 12/26/2007 1,0164 1,58 1,6 ND µg/m³ R14916 1,2-Dichloropropane TO-15 12/26/2007 0,9375 1,58 1,5 ND µg/m³ R14916 1,2-Dichlorobenzene TO-15 12/26/2007 0,6888 1,58 1,1 ND µg/m³ R14916 1,3-Dichlorobenzene TO-15 12/26/2007 0,6888 1,58 0,94 ND µg/m³ R14916 1,3-Dichlorobenzene TO-15 12/26/2007 0,6861 1,58 0,94 ND µg/m³ R14916 1,3-Dichlorobenzene TO-15 12/26/2007 0,6611 1,58 0,94 ND µg/m³ R14916 1,4-Dioxane TO-15 12/26/2007 0,504 1,58 0,80 ND µg/m³ R14916 1,4-Dioxane TO-15 12/26/2007 0,504 1,58 0,80 ND µg/m³ R14916 1,4-Dioxane TO-15 12/26/2007 0,504 1,58 0,80 ND µg/m³ R14916 2-Butanone (MEK) TO-15 12/26/2007 0,504 1,58 0,80 ND µg/m³ R14916 2-Hexanone TO-15 12/26/2007 0,504 1,58 0,80 ND µg/m³ R14916 2-Hexanone TO-15 12/26/2007 0,504 1,58 0,80 ND µg/m³ R14916 2-Hexanone TO-15 12/26/2007 0,504 1,58 0,80 ND µg/m³ R14916 2-Hexanone TO-15 12/26/2007 0,504 1,58 0,80 ND µg/m³ R14916 2-Hexanone TO-15 12/26/2007 0,504 1,58 0,80 ND µg/m³ R14916 2-Hexanone TO-15 12/26/2007 0,504 1,58 0,00 ND µg/m³ R14916 2-Hexanone TO-15 12/26/2007 0,504 1,58 0,00 ND µg/m³ R14916 0-Hexthyl-2-Pentanone (MIBK) TO-15 12/26/2007 0,686 1,58 1,0 ND µg/m³ R14916 0-Hexthyl-2-Pentanone (MIBK) TO-15 12/26/2007 0,681 1,58 1,1 ND µg/m³ R14916 0-Hexthyl-2-Pentanone (MIBK) TO-15 12/26/2007 0,691 1,58 1,1 ND µg/m³ R14916 0-Hexanone TO-15 12/26/2007 0,4976 1,58 0,67 ND µg/m³ R14916 0-Hexanone TO-15 12/26/2007 0,4976 1,58 0,67 ND µg/m³ R14916 0-Hexanone TO-15 12/26/2007 0,4976 1,58 0,67 ND µg/m³ R14916 0-Hexanone Chlorobenzene TO-15 12/26/2007 0,4976 1,58 0,67 ND µg/m³ R14916 0-Hexanone Chlorobenzene TO-15 12/26/2007 0,4976 1,58 0,67 ND µg/m³ R14916 0-Hexanone Chlorobenzene TO-15 12/26/2007 0,4976 1,58 0,67 ND µg/m³ R14916 0-Hexanone Chlorobenzene TO-15 12/26/2007 0,4925 1,58 0,67 ND µg/m³ R14916 0-Hexanone Chlorobenzene TO-1	1,2,4-Trimethylbenzene	TO-15	12/26/2007	0.8856	1.58	1.4	ND		R14916
1,2-Dichloroethane TO-15 12/26/2007 0.648 1.58 1.0 ND µg/m³ R14916 1,2-Dichloropropane TO-15 12/26/2007 1.0164 1.58 1.6 ND µg/m³ R14916 1,2-Dichloropropane TO-15 12/26/2007 0.9375 1.58 1.5 ND µg/m³ R14916 1,3-Brutadiene TO-15 12/26/2007 0.9375 1.58 1.5 ND µg/m³ R14916 1,3-Brutadiene TO-15 12/26/2007 0.5967 1.58 0.94 ND µg/m³ R14916 1,3-Dichlorobenzene TO-15 12/26/2007 0.5967 1.58 0.94 ND µg/m³ R14916 1,3-Dichlorobenzene TO-15 12/26/2007 0.5967 1.58 0.94 ND µg/m³ R14916 1,4-Dichlorobenzene TO-15 12/26/2007 0.6611 1.58 1.0 ND µg/m³ R14916 1,4-Dichlorobenzene TO-15 12/26/2007 0.6611 1.58 1.0 ND µg/m³ R14916 1,4-Dichlorobenzene TO-15 12/26/2007 0.504 1.58 0.60 ND µg/m³ R14916 1,4-Dichlorobenzene TO-15 12/26/2007 0.504 1.58 0.80 ND µg/m³ R14916 1,4-Dichlorobenzene TO-15 12/26/2007 0.4425 1.58 0.70 T7 µg/m³ R14916 1,4-Dichlorobenzene TO-15 12/26/2007 0.4425 1.58 0.70 T7 µg/m³ R14916 1,4-Dichlorobenzene TO-15 12/26/2007 0.4425 1.58 0.70 T7 µg/m³ R14916 1,4-Dichlorobenzene TO-15 12/26/2007 0.4425 1.58 0.70 T7 µg/m³ R14916 1,4-Dichlorobenzene TO-15 12/26/2007 0.4861 1.58 1.4 ND µg/m³ R14916 1,4-Dichlorobenzene TO-15 12/26/2007 0.738 1.58 1.2 ND µg/m³ R14916 1,4-Dichlorobenzene TO-15 12/26/2007 0.738 1.58 1.2 ND µg/m³ R14916 1,4-Dichlorobenzene TO-15 12/26/2007 0.5712 1.58 0.90 180 µg/m³ R14916 1,4-Dichloride TO-15 12/26/2007 0.5712 1.58 0.90 180 µg/m³ R14916 1,4-Dichloride TO-15 12/26/2007 0.690 1.58 1.1 ND µg/m³ R14916 1,4-Dichloride TO-15 12/26/2007 0.690 1.58 1.1 ND µg/m³ R14916 1,4-Dichloride TO-15 12/26/2007 0.4776 1.58 1.2 ND µg/m³ R14916 1,4-Dichloride TO-15 12/26/2007 0.4776 1.58 0.60 ND µg/m³ R14916 1,4-Dichloride TO-15 12/26/2007 0.4476 1.58 0.60 ND µg/m³ R14916 1,4-Dichloride TO-15 12/26/2007 0.4976 1.58 0.60 ND µg/m³ R14916 1,4-Dichloride TO-15 12/26/2007 0.4976 1.58 0.60 ND µg/m³ R14916 1,4-Dichloride TO-15 12/26/2007 0.4976 1.58 0.67 ND µg/m³ R14916 1,4-Dichloride TO-15 12/26/2007 0.4976 1.58 0.67 ND µg/m³ R14916 1,4-Dichloride TO-15 12/26/2007 0.4976 1.58 0.60 ND µg/m³ R14916 1,4-Dichloride TO-15 12/26/20	1,2-Dibromoethane(Ethylene dibromide)	TO-15	12/26/2007	1.0752	1.58	1.7	ND	µg/m³	R14916
1,2-Dichloropropane	1,2-Dichlorobenzene	TO-15	12/26/2007	0.601	1.58	0.95	ND	μg/m³	R14916
1,2-	1,2-Dichloroethane	TO-15	12/26/2007	0.648	1.58	1.0	ND	μg/m³	R14916
dichlorotetrafluoroethane(F114) TO-15 12/26/2007 0.6888 1.58 1.1 ND µg/m³ R14916 1,3,5-Trimethylbenzene TO-15 12/26/2007 0.5967 1.58 0.94 ND µg/m³ R14916 1,3-Dichlorobenzene TO-15 12/26/2007 0.3606 1.58 0.57 ND µg/m³ R14916 1,4-Dichlorobenzene TO-15 12/26/2007 0.6611 1.58 1.0 ND µg/m³ R14916 1,4-Dioxane TO-15 12/26/2007 0.504 1.58 0.80 ND µg/m³ R14916 2-Butanone (MEK) TO-15 12/26/2007 0.504 1.58 0.80 ND µg/m³ R14916 2-Hexanone TO-15 12/26/2007 0.861 1.58 1.4 ND µg/m³ R14916 2-Hexanone TO-15 12/26/2007 0.861 1.58 1.4 ND µg/m³ R14916 4-Ethyl Toluene TO-15 12/26/2007 0.738 1.	1,2-Dichloropropane	TO-15	12/26/2007	1.0164	1.58	1.6	ND	μg/m³	R14916
1,3-Butadiene TO-15 12/26/2007 0.5967 1.58 0.94 ND μg/m³ R14916 1,3-Dichlorobenzene TO-15 12/26/2007 0.3606 1.58 0.57 ND μg/m³ R14916 1,4-Dichlorobenzene TO-15 12/26/2007 0.6611 1.58 1.0 ND μg/m³ R14916 1,4-Dicknere TO-15 12/26/2007 0.504 1.58 0.80 ND μg/m³ R14916 1,4-Dioxane TO-15 12/26/2007 0.504 1.58 0.80 ND μg/m³ R14916 2-Butanone (MEK) TO-15 12/26/2007 0.4425 1.58 0.80 ND μg/m³ R14916 2-Butanone (MEK) TO-15 12/26/2007 0.4425 1.58 0.70 77 μg/m³ R14916 2-Butanone (MEK) TO-15 12/26/2007 0.738 1.58 1.4 ND μg/m³ R14916 4-Ethyl Toluene TO-15 12/26/2007 0.738 1.58 1.2 ND μg/m³ R14916 4-Ethyl Toluene TO-15 12/26/2007 0.656 1.58 1.0 ND μg/m³ R14916 4-Ethyl-2-Pentanone (MIBK) TO-15 12/26/2007 0.656 1.58 1.0 ND μg/m³ R14916 Acetone TO-15 12/26/2007 0.5712 1.58 0.90 180 μg/m³ R14916 Benzene TO-15 12/26/2007 0.8932 1.58 1.4 ND μg/m³ R14916 Benzene TO-15 12/26/2007 0.8932 1.58 1.4 ND μg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.891 1.58 1.4 ND μg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.891 1.58 1.1 ND μg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.871 1.58 1.4 ND μg/m³ R14916 Bromomethane TO-15 12/26/2007 0.4976 1.58 1.2 ND μg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 0.79 ND μg/m³ R14916 Carbon Tetrachloride TO-15 12/26/2007 0.4935 1.58 0.63 ND μg/m³ R14916 Carbon Tetrachloride TO-15 12/26/2007 0.4935 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.4232 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.4925 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.4232 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.4232 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.3245 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.3245 1.58 0.63 ND μg/m³ R14916	1,2- dichlorotetrafluoroethane(F114)	TO-15	12/26/2007	0.9375	1.58	1.5	ND	µg/m³	R14916
1,3-Dichlorobenzene TO-15 12/26/2007 0.3606 1.58 0.57 ND μg/m³ R14916 1,4-Dichlorobenzene TO-15 12/26/2007 0.6611 1.58 1.0 ND μg/m³ R14916 1,4-Dichlorobenzene TO-15 12/26/2007 0.504 1.58 0.80 ND μg/m³ R14916 1,4-Dioxane TO-15 12/26/2007 0.504 1.58 0.80 ND μg/m³ R14916 2-Butanone (MEK) TO-15 12/26/2007 0.4425 1.58 0.70 77 μg/m³ R14916 2-Hexanone TO-15 12/26/2007 0.861 1.58 1.4 ND μg/m³ R14916 4-Ethyl Toluene TO-15 12/26/2007 0.738 1.58 1.2 ND μg/m³ R14916 4-Ethyl-2-Pentanone (MIBK) TO-15 12/26/2007 0.656 1.58 1.0 ND μg/m³ R14916 Acetone TO-15 12/26/2007 0.656 1.58 1.0 ND μg/m³ R14916 Benzene TO-15 12/26/2007 0.8932 1.58 1.4 3.5 μg/m³ R14916 Benzene TO-15 12/26/2007 0.8932 1.58 1.4 3.5 μg/m³ R14916 Benzene TO-15 12/26/2007 0.69 1.58 1.1 ND μg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.871 1.58 1.4 ND μg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.871 1.58 1.4 ND μg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.776 1.58 1.4 ND μg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.776 1.58 1.2 ND μg/m³ R14916 Bromomethane TO-15 12/26/2007 0.776 1.58 1.2 ND μg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 0.79 ND μg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4232 1.58 1.5 ND μg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.4232 1.58 0.67 ND μg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007	1,3,5-Trimethylbenzene	TO-15	12/26/2007	0.6888	1.58	1.1	ND	μg/m³	R14916
1,4-Dichlorobenzene TO-15 12/26/2007 0.6611 1.58 1.0 ND µg/m³ R14916 1,4-Dioxane TO-15 12/26/2007 0.504 1.58 0.80 ND µg/m³ R14916 2-Butanone (MEK) TO-15 12/26/2007 0.4425 1.58 0.70 77 µg/m³ R14916 2-Butanone (MEK) TO-15 12/26/2007 0.861 1.58 0.70 77 µg/m³ R14916 2-Hexanone TO-15 12/26/2007 0.861 1.58 1.4 ND µg/m³ R14916 4-Ethyl Toluene TO-15 12/26/2007 0.738 1.58 1.2 ND µg/m³ R14916 4-Methyl-2-Pentanone (MIBK) TO-15 12/26/2007 0.656 1.58 1.0 ND µg/m³ R14916 A-Cetone TO-15 12/26/2007 0.5712 1.58 0.90 180 µg/m³ R14916 Benzene TO-15 12/26/2007 0.8932 1.58 1.4 3.5 µg/m³ R14916 Benzyl Chloride TO-15 12/26/2007 0.8932 1.58 1.1 ND µg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.891 1.58 1.1 ND µg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.871 1.58 1.4 ND µg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.871 1.58 1.4 ND µg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.871 1.58 1.4 ND µg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.871 1.58 1.4 ND µg/m³ R14916 Bromomethane TO-15 12/26/2007 0.776 1.58 1.2 ND µg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 1.2 ND µg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 1.5 ND µg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.4232 1.58 0.67 ND µg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.4232 1.58 0.67 ND µg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chloroform TO-1	1,3-Butadiene	TO-15	12/26/2007	0.5967	1.58	0.94	ND		R14916
1,4-Dioxane TO-15 12/26/2007 0.504 1.58 0.80 ND µg/m³ R14916 2-Butanone (MEK) TO-15 12/26/2007 0.4425 1.58 0.70 77 µg/m³ R14916 2-Hexanone (MEK) TO-15 12/26/2007 0.861 1.58 1.4 ND µg/m³ R14916 4-Ethyl Toluene TO-15 12/26/2007 0.738 1.58 1.2 ND µg/m³ R14916 4-Methyl-2-Pentanone (MIBK) TO-15 12/26/2007 0.656 1.58 1.0 ND µg/m³ R14916 4-Methyl-2-Pentanone (MIBK) TO-15 12/26/2007 0.5712 1.58 0.90 180 µg/m³ R14916 Acetone TO-15 12/26/2007 0.5712 1.58 0.90 180 µg/m³ R14916 Benzene TO-15 12/26/2007 0.8932 1.58 1.4 3.5 µg/m³ R14916 Benzene TO-15 12/26/2007 0.8932 1.58 1.1 ND µg/m³ R14916 Bromofichloromethane TO-15 12/26/2007 0.871 1.58 1.4 ND µg/m³ R14916 Bromoform TO-15 12/26/2007 0.871 1.58 1.4 ND µg/m³ R14916 Bromoform TO-15 12/26/2007 0.776 1.58 1.5 ND µg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 1.2 ND µg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 1.5 ND µg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 1.5 ND µg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 0.67 ND µg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.4232 1.58 0.67 ND µg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.4232 1.58 0.63 ND µg/m³ R14916 Chlorothane TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chlorothane TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chlorothane TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chlorothane TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chlorothane TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chlorothane TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chlorothane TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chlorothane TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chlorothane TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chlorothane TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chlorothane TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chlorothane TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chlorothane TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chlorothane TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chlorothane TO-15 12/26/2007	1,3-Dichlorobenzene	TO-15	12/26/2007	0.3606	1.58	0.57	ND		R14916
2-Butanone (MEK) TO-15 12/26/2007 0.4425 1.58 0.70 77 µg/m³ R14916 2-Hexanone TO-15 12/26/2007 0.861 1.58 1.4 ND µg/m³ R14916 4-Ethyl Toluene TO-15 12/26/2007 0.738 1.58 1.2 ND µg/m³ R14916 4-Methyl-2-Pentanone (MIBK) TO-15 12/26/2007 0.656 1.58 1.0 ND µg/m³ R14916 Acetone TO-15 12/26/2007 0.5712 1.58 0.90 180 µg/m³ R14916 Benzene TO-15 12/26/2007 0.8932 1.58 1.4 3.5 µg/m³ R14916 Benzyl Chloride TO-15 12/26/2007 0.8932 1.58 1.1 ND µg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.8932 1.58 1.1 ND µg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.871 1.58 1.4 ND µg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.871 1.58 1.4 ND µg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.871 1.58 1.58 1.8 ND µg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 1.5 ND µg/m³ R14916 Carbon Tetrachloride TO-15 12/26/2007 0.4976 1.58 0.67 ND µg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.396 1.58 0.63 ND µg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 3.1 ND µg/m³ R14916 Chloromethane TO-15 12/26/2007 0.396 1.58 3.1 ND µg/m³ R14916 Chloromethane TO-15 12/26/2007 0.396 1.58 3.1 ND µg/m³ R14916 Chloromethane TO-15 12/26/2007 0.396 1.58 3.1 ND µg/m³ R14916 Chloromethane TO-15 12/26/2007 0.396 1.58 3.1 ND µg/m³ R14916 Chloromethane TO-15 12/26/2007 0.396 1.58 3.1 ND µg/m³ R14916 Chloromethane	1,4-Dichlorobenzene	TO-15	12/26/2007	0.6611	1.58	1.0	ND	μg/m³	R14916
2-Butanone (MEK) TO-15 12/26/2007 0.4425 1.58 0.70 77 µg/m³ R14916 2-Hexanone TO-15 12/26/2007 0.861 1.58 1.4 ND µg/m³ R14916 4-Ethyl Toluene TO-15 12/26/2007 0.738 1.58 1.2 ND µg/m³ R14916 4-Methyl-2-Pentanone (MIBK) TO-15 12/26/2007 0.656 1.58 1.0 ND µg/m³ R14916 Acetone TO-15 12/26/2007 0.5712 1.58 0.90 180 µg/m³ R14916 Benzene TO-15 12/26/2007 0.8932 1.58 1.4 3.5 µg/m³ R14916 Benzyl Chloride TO-15 12/26/2007 0.8932 1.58 1.1 ND µg/m³ R14916 Benzyl Chloride TO-15 12/26/2007 0.8911 1.58 1.1 ND µg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.871 1.58 1.4 ND µg/m³ R14916 Bromoform TO-15 12/26/2007 0.871 1.58 1.4 ND µg/m³ R14916 Bromomethane TO-15 12/26/2007 0.871 1.58 1.58 1.8 ND µg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 1.58 0.67 ND µg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.4232 1.58 3.1 ND µg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 3.1 ND µg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 3.1 ND µg/m³ R14916 Chloromethane TO-15 12/26/2007 0.396 1.58 3.1 ND µg/m³ R14916 Chloromethane TO-15 12/26/2007 0.396 1.58 3.1 ND µg/m³ R14916 Chloromethane TO-15 12/26/2007 0.396 1.58 3.1 ND µg/m³ R14916 Chloromethane TO-15 12/26/2007 0.396 1.58 3.1 ND µg/m³ R14916 Chloromethane TO-15 12/26/2007 0.396 1.58 3.1 ND µg/m³ R14916 Chloromethane TO-15 12/26/2007 0.396 1.58 3.1 ND µg/m³ R14916 Chloromethane	1,4-Dioxane	TO-15	12/26/2007	0.504	1.58	0.80	ND	μg/m³	R14916
4-Ethyl Toluene TO-15 12/26/2007 0.738 1.58 1.2 ND μg/m³ R14916 4-Methyl-2-Pentanone (MIBK) TO-15 12/26/2007 0.656 1.58 1.0 ND μg/m³ R14916 Acetone TO-15 12/26/2007 0.5712 1.58 0.90 180 μg/m³ R14916 Benzene TO-15 12/26/2007 0.8932 1.58 1.4 3.5 μg/m³ R14916 Benzyl Chloride TO-15 12/26/2007 0.69 1.58 1.1 ND μg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.871 1.58 1.4 ND μg/m³ R14916 Bromoform TO-15 12/26/2007 0.871 1.58 1.4 ND μg/m³ R14916 Bromoform TO-15 12/26/2007 0.871 1.58 1.4 ND μg/m³ R14916 Bromoform TO-15 12/26/2007 0.776 1.58 1.2 ND μg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 0.79 ND μg/m³ R14916 Carbon Tetrachloride TO-15 12/26/2007 0.4976 1.58 1.5 ND μg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.4232 1.58 0.67 ND μg/m³ R14916 Chloroethane TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 1.952 1.58 3.1 ND μg/m³ R14916 Chloromethane	2-Butanone (MEK)	TO-15	12/26/2007	0.4425	1.58	0.70	77	µg/m³	R14916
4-Methyl-2-Pentanone (MIBK) TO-15 12/26/2007 0.656 1.58 1.0 ND μg/m³ R14916 Acetone TO-15 12/26/2007 0.5712 1.58 0.90 180 μg/m³ R14916 Benzene Benzene TO-15 12/26/2007 0.8932 1.58 1.4 3.5 μg/m³ R14916 Benzyl Chloride TO-15 12/26/2007 0.69 1.58 1.1 ND μg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.8971 1.58 1.1 ND μg/m³ R14916 Bromoform TO-15 12/26/2007 1.7578 1.58 2.8 ND μg/m³ R14916 Bromomethane TO-15 12/26/2007 0.7776 1.58 1.2 ND μg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 0.79 ND μg/m³ R14916 Carbon Tetrachloride TO-15 12/26/2007 0.9435 1.58 1.5 ND μg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.4232 1.58 0.67 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 1.952 1.58 3.1 ND μg/m³ R14916 Chloromethane TO-15 12/26/2007 0.7245 1.58 1.1 ND μg/m³ R14916 Chloromethane	2-Hexanone	TO-15	12/26/2007	0.861	1.58	1.4	ND	μg/m³	R14916
Acetone TO-15 12/26/2007 0.5712 1.58 0.90 180 μg/m³ R14916 Benzene TO-15 12/26/2007 0.8932 1.58 1.4 3.5 μg/m³ R14916 Benzyl Chloride TO-15 12/26/2007 0.69 1.58 1.1 ND μg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.871 1.58 1.4 ND μg/m³ R14916 Bromoform TO-15 12/26/2007 1.7578 1.58 2.8 ND μg/m³ R14916 Bromomethane TO-15 12/26/2007 0.776 1.58 1.2 ND μg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 0.79 ND μg/m³ R14916 Carbon Tetrachloride TO-15 12/26/2007 0.9435 1.58 1.5 ND μg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.396 1.58 0.63	4-Ethyl Toluene	TO-15	12/26/2007	0.738	1.58	1.2	ND	μg/m³	R14916
Benzene TO-15 12/26/2007 0.8932 1.58 1.4 3.5 μg/m³ R14916 Benzyl Chloride TO-15 12/26/2007 0.69 1.58 1.1 ND μg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.871 1.58 1.4 ND μg/m³ R14916 Bromoform TO-15 12/26/2007 1.7578 1.58 2.8 ND μg/m³ R14916 Bromomethane TO-15 12/26/2007 0.776 1.58 1.2 ND μg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 0.79 ND μg/m³ R14916 Carbon Tetrachloride TO-15 12/26/2007 0.9435 1.58 1.5 ND μg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.4232 1.58 0.67 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 1.952 1.58 3.1	4-Methyl-2-Pentanone (MIBK)	TO-15	12/26/2007	0.656	1.58	1.0	ND	µg/m³	R14916
Benzyl Chloride TO-15 12/26/2007 0.69 1.58 1.1 ND μg/m³ R14916 Bromodichloromethane TO-15 12/26/2007 0.871 1.58 1.4 ND μg/m³ R14916 Bromoform TO-15 12/26/2007 1.7578 1.58 2.8 ND μg/m³ R14916 Bromomethane TO-15 12/26/2007 0.776 1.58 1.2 ND μg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 0.79 ND μg/m³ R14916 Carbon Tetrachloride TO-15 12/26/2007 0.9435 1.58 1.5 ND μg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.4232 1.58 0.67 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 1.952 1.58 0.63 ND μg/m³ R14916 Chloromethane TO-15 12/26/2007 0.7245 1.58 3.1 </td <td>Acetone</td> <td>TO-15</td> <td>12/26/2007</td> <td>0.5712</td> <td>1.58</td> <td>0.90</td> <td>180</td> <td>µg/m³</td> <td>R14916</td>	Acetone	TO-15	12/26/2007	0.5712	1.58	0.90	180	µg/m³	R14916
Bromodichloromethane TO-15 12/26/2007 0.871 1.58 1.4 ND μg/m³ R14916 Bromoform TO-15 12/26/2007 1.7578 1.58 2.8 ND μg/m³ R14916 Bromomethane TO-15 12/26/2007 0.776 1.58 1.2 ND μg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 0.79 ND μg/m³ R14916 Carbon Tetrachloride TO-15 12/26/2007 0.9435 1.58 1.5 ND μg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.4232 1.58 0.67 ND μg/m³ R14916 Chloroethane TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 1.952 1.58 3.1 ND μg/m³ R14916 Chloromethane TO-15 12/26/2007 0.7245 1.58 1.1 <td>Benzene</td> <td>TO-15</td> <td>12/26/2007</td> <td>0.8932</td> <td>1.58</td> <td>1.4</td> <td>3.5</td> <td>µg/m³</td> <td>R14916</td>	Benzene	TO-15	12/26/2007	0.8932	1.58	1.4	3.5	µg/m³	R14916
Bromoform TO-15 12/26/2007 1.7578 1.58 2.8 ND μg/m³ R14916 Bromomethane TO-15 12/26/2007 0.776 1.58 1.2 ND μg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 0.79 ND μg/m³ R14916 Carbon Tetrachloride TO-15 12/26/2007 0.9435 1.58 1.5 ND μg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.4232 1.58 0.67 ND μg/m³ R14916 Chloroethane TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 1.952 1.58 3.1 ND μg/m³ R14916 Chloromethane TO-15 12/26/2007 0.7245 1.58 1.1 ND μg/m³ R14916	Benzyl Chloride	TO-15	12/26/2007	0.69	1.58	1.1	ND	µg/m³	R14916
Bromomethane TO-15 12/26/2007 0.776 1.58 1.2 ND μg/m³ R14916 Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 0.79 ND μg/m³ R14916 Carbon Tetrachloride TO-15 12/26/2007 0.9435 1.58 1.5 ND μg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.4232 1.58 0.67 ND μg/m³ R14916 Chloroethane TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 1.952 1.58 3.1 ND μg/m³ R14916 Chloromethane TO-15 12/26/2007 0.7245 1.58 1.1 ND μg/m³ R14916	Bromodichloromethane	TO-15	12/26/2007	0.871	1.58	1.4	ND	µg/m³	R14916
Carbon Disulfide TO-15 12/26/2007 0.4976 1.58 0.79 ND μg/m³ R14916 Carbon Tetrachloride TO-15 12/26/2007 0.9435 1.58 1.5 ND μg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.4232 1.58 0.67 ND μg/m³ R14916 Chloroethane TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 1.952 1.58 3.1 ND μg/m³ R14916 Chloromethane TO-15 12/26/2007 0.7245 1.58 1.1 ND μg/m³ R14916	Bromoform	TO-15	12/26/2007	1.7578	1.58	2.8	ND	µg/m³	R14916
Carbon Tetrachloride TO-15 12/26/2007 0.9435 1.58 1.5 ND μg/m³ R14916 Chlorobenzene TO-15 12/26/2007 0.4232 1.58 0.67 ND μg/m³ R14916 Chloroethane TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 1.952 1.58 3.1 ND μg/m³ R14916 Chloromethane TO-15 12/26/2007 0.7245 1.58 1.1 ND μg/m³ R14916	Bromomethane	TO-15	12/26/2007	0.776	1.58	1.2	ND	µg/m³	R14916
Chlorobenzene TO-15 12/26/2007 0.4232 1.58 0.67 ND μg/m³ R14916 Chloroethane TO-15 12/26/2007 0.396 1.58 0.63 ND μg/m³ R14916 Chloroform TO-15 12/26/2007 1.952 1.58 3.1 ND μg/m³ R14916 Chloromethane TO-15 12/26/2007 0.7245 1.58 1.1 ND μg/m³ R14916	Carbon Disulfide	TO-15	12/26/2007	0.4976	1.58	0.79	ND	µg/m³	R14916
$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	Carbon Tetrachloride	TO-15	12/26/2007	0.9435	1.58	1.5	ND	µg/m³	R14916
Chloroform TO-15 12/26/2007 1.952 1.58 3.1 ND μ g/m³ R14916 Chloromethane TO-15 12/26/2007 0.7245 1.58 1.1 ND μ g/m³ R14916	Chlorobenzene	TO-15	12/26/2007	0.4232	1.58	0.67	ND	µg/m³	R14916
Chloromethane TO-15 12/26/2007 0.7245 1.58 1.1 ND μg/m³ R14916	Chloroethane	TO-15	12/26/2007	0.396	1.58	0.63	ND	µg/m³	R14916
Chloromethane TO-15 12/26/2007 0.7245 1.58 1.1 ND μg/m³ R14916	Chloroform	TO-15	12/26/2007	1.952	1.58	3.1	ND		R14916
	Chloromethane	TO-15	12/26/2007	0.7245	1.58	1.1	ND		R14916
	cis-1,2-dichloroethene	TO-15	12/26/2007	0.5544	1.58	0.88	ND		R14916

These analyses were performed according to State of California Environmental Laboratory Accreditation program, Certificate # 1991

Report prepared for: Kelly Krohn

Engeo

Date Received: 12/17/2007

Date Reported: 12/27/2007

Client Sample ID: SG-11

224 Rickenbacker Circle

Sample Matrix: AIR

Sample Location:

Date/Time Sampled 12/17/2007 11:22:00 AM

Lab Sample ID: 0712091-001

Date Prepared:

Parameters	Analysis Method	Date Analyzed	RL	Dilution Factor	MRL	Result	Units	Analytical Batch
cis-1,3-Dichloropropene	TO-15	12/26/2007	0.3632	1.58	0.57	ND	μg/m³	R14916
Dibromochloromethane	TO-15	12/26/2007	0.9372	1.58	1.5	ND	μg/m³	R14916
Dichlorodifluoromethane	TO-15	12/26/2007	0.7425	1.58	1.2	ND	μg/m³	R14916
Ethyl Acetate	TO-15	12/26/2007	0.4248	1.58	0.67	ND	µg/m³	R14916
Ethyl Benzene	TO-15	12/26/2007	0.31062	1.58	0.49	ND	μg/m³	R14916
Freon 113	TO-15	12/26/2007	0.9192	1.58	1.5	ND	μg/m³	R14916
Hexachlorobutadiene	TO-15	12/26/2007	1.8139	1.58	2.9	ND	µg/m³	R14916
Hexane	TO-15	12/26/2007	1.7952	1.58	2.8	ND	μg/m³	R14916
Isopropanol	TO-15	12/26/2007	1.6359	1.58	2.6	ND	μg/m³	R14916
m,p-Xylene	TO-15	12/26/2007	0.492	1.58	0.78	38	μg/m³	R14916
Methylene Chloride	TO-15	12/26/2007	0.6859	1.58	1.1	ND	μg/m³	R14916
MTBE	TO-15	12/26/2007	0.5054	1.58	0.80	ND	μg/m³	R14916
Naphthalene	TO-15	12/26/2007	2.62	1.58	4.1	ND	µg/m³	R14916
o-xylene	TO-15	12/26/2007	0.62062	1.58	0.98	11	μg/m³	R14916
Styrene	TO-15	12/26/2007	0.639	1.58	1.0	ND	μg/m³	R14916
Tetrachloroethene	TO-15	12/26/2007	1.2882	1.58	2.0	64	μg/m³	R14916
Tetrahydrofuran	TO-15	12/26/2007	0.413	1.58	0.65	ND	μg/m³	R14916
Toluene	TO-15	12/26/2007	0.5278	1.58	0.83	25	μg/m³	R14916
trans-1,2-Dichloroethene	TO-15	12/26/2007	0.5544	1.58	0.88	ND	μg/m³	R14916
Trichloroethene	TO-15	12/26/2007	0.52626	1.58	0.83	ND	μg/m³	R14916
Trichlorofluoromethane	TO-15	12/26/2007	0.693	1.58	1.1	ND	μg/m³	R14916
Vinyl Acetate	TO-15	12/26/2007	0.64064	1.58	1.0	ND	μg/m³	R14916
Vinyl Chloride	TO-15	12/26/2007	0.24832	1.58	0.39	ND	μg/m³	R14916
Surr: 4-Bromofluorobenzene	TO-15	12/26/2007	0	1.58	65-135	95.0	%REC	R14916

Note: Due to low initial pressure in a Summa canister, sample was diluted prior to the analysis by addition of argon. Results are reported to the MDL.

Report prepared for: Kelly Krohn **Date Received:** 12/17/2007 Engeo

Date Reported: 12/27/2007

Client Sample ID: SG-12

Sample Location: 224 Rickenbacker Circle

Sample Matrix: AIR

Date/Time Sampled 12/17/2007 1:05:00 PM **Lab Sample ID:** 0712091-002

Date Prepared:

Parameters	Analysis Method	Date Analyzed	RL	Dilution Factor	MRL	Result	Units	Analytical Batch
1,1 - Dichloroethene	TO-15	12/26/2007	0.794	1.56	1.2	ND	μg/m³	R14916
1,1,1,2-Tetrachloroethane	TO-15	12/26/2007	0.687	1.56	1.1	ND	μg/m³	R14916
1,1,1-Trichloroethane	TO-15	12/26/2007	0.819	1.56	1.3	ND	μg/m³	R14916
1,1,2,2-Tetrachloroethane	TO-15	12/26/2007	1.0305	1.56	1.6	ND	µg/m³	R14916
1,1,2-Trichloroethane	TO-15	12/26/2007	1.0374	1.56	1.6	ND	µg/m³	R14916
1,1-Dichloroethane	TO-15	12/26/2007	0.6885	1.56	1.1	ND	µg/m³	R14916
1,2,4-Trichlorobenzene	TO-15	12/26/2007	0.4984	1.56	0.78	ND	µg/m³	R14916
1,2,4-Trimethylbenzene	TO-15	12/26/2007	0.8856	1.56	1.4	ND	µg/m³	R14916
1,2-Dibromoethane(Ethylene dibromide)	TO-15	12/26/2007	1.0752	1.56	1.7	ND	μg/m³	R14916
1,2-Dichlorobenzene	TO-15	12/26/2007	0.601	1.56	0.94	ND	µg/m³	R14916
1,2-Dichloroethane	TO-15	12/26/2007	0.648	1.56	1.0	ND	μg/m³	R14916
1,2-Dichloropropane	TO-15	12/26/2007	1.0164	1.56	1.6	ND	μg/m³	R14916
1,2- dichlorotetrafluoroethane(F114)	TO-15	12/26/2007	0.9375	1.56	1.5	ND	µg/m³	R14916
1,3,5-Trimethylbenzene	TO-15	12/26/2007	0.6888	1.56	1.1	ND	µg/m³	R14916
1,3-Butadiene	TO-15	12/26/2007	0.5967	1.56	0.93	ND	µg/m³	R14916
1,3-Dichlorobenzene	TO-15	12/26/2007	0.3606	1.56	0.56	ND	µg/m³	R14916
1,4-Dichlorobenzene	TO-15	12/26/2007	0.6611	1.56	1.0	ND	μg/m³	R14916
1,4-Dioxane	TO-15	12/26/2007	0.504	1.56	0.79	ND	µg/m³	R14916
2-Butanone (MEK)	TO-15	12/26/2007	0.4425	1.56	0.69	31	µg/m³	R14916
2-Hexanone	TO-15	12/26/2007	0.861	1.56	1.3	ND	µg/m³	R14916
4-Ethyl Toluene	TO-15	12/26/2007	0.738	1.56	1.2	ND	µg/m³	R14916
4-Methyl-2-Pentanone (MIBK)	TO-15	12/26/2007	0.656	1.56	1.0	ND	µg/m³	R14916
Acetone	TO-15	12/26/2007	0.5712	1.56	0.89	70	µg/m³	R14916
Benzene	TO-15	12/26/2007	0.8932	1.56	1.4	2.5 J	µg/m³	R14916
Benzyl Chloride	TO-15	12/26/2007	0.69	1.56	1.1	ND	µg/m³	R14916
Bromodichloromethane	TO-15	12/26/2007	0.871	1.56	1.4	ND	µg/m³	R14916
Bromoform	TO-15	12/26/2007	1.7578	1.56	2.7	ND	µg/m³	R14916
Bromomethane	TO-15	12/26/2007	0.776	1.56	1.2	ND	µg/m³	R14916
Carbon Disulfide	TO-15	12/26/2007	0.4976	1.56	0.78	ND	µg/m³	R14916
Carbon Tetrachloride	TO-15	12/26/2007	0.9435	1.56	1.5	ND	µg/m³	R14916
Chlorobenzene	TO-15	12/26/2007	0.4232	1.56	0.66	ND	µg/m³	R14916
Chloroethane	TO-15	12/26/2007	0.396	1.56	0.62	ND	µg/m³	R14916
Chloroform	TO-15	12/26/2007	1.952	1.56	3.0	ND	µg/m³	R14916
Chloromethane	TO-15	12/26/2007	0.7245	1.56	1.1	ND	µg/m³	R14916
cis-1,2-dichloroethene	TO-15	12/26/2007	0.5544	1.56	0.86	ND	µg/m³	R14916
cis-1,3-Dichloropropene	TO-15	12/26/2007	0.3632	1.56	0.57	ND	µg/m³	R14916
Dibromochloromethane	TO-15	12/26/2007	0.9372	1.56	1.5	ND	μg/m³	R14916
Dichlorodifluoromethane	TO-15	12/26/2007	0.7425	1.56	1.2	ND	μg/m³	R14916
Ethyl Acetate	TO-15	12/26/2007	0.4248	1.56	0.66	ND	μg/m³	R14916
Ethyl Benzene	TO-15	12/26/2007	0.31062	1.56	0.48	ND	μg/m³	R14916
Freon 113	TO-15	12/26/2007	0.9192	1.56	1.4	ND	μg/m³	R14916
Hexachlorobutadiene	TO-15	12/26/2007	1.8139	1.56	2.8	ND	μg/m³	R14916

Report prepared for: Kelly Krohn **Date Received:** 12/17/2007

Engeo **Date Reported:** 12/27/2007

Client Sample ID: SG-12 Lab Sample ID: 0712091-002

Sample Location: 224 Rickenbacker Circle Date Prepared:

Sample Matrix: AIR

Date/Time Sampled 12/17/2007 1:05:00 PM

Parameters	Analysis Method	Date Analyzed	RL	Dilution Factor	MRL	Result	Units	Analytical Batch
Hexane	TO-15	12/26/2007	1.7952	1.56	2.8	ND	μg/m³	R14916
Isopropanol	TO-15	12/26/2007	1.6359	1.56	2.6	ND	μg/m³	R14916
m,p-Xylene	TO-15	12/26/2007	0.492	1.56	0.77	25	μg/m³	R14916
Methylene Chloride	TO-15	12/26/2007	0.6859	1.56	1.1	ND	µg/m³	R14916
MTBE	TO-15	12/26/2007	0.5054	1.56	0.79	ND	µg/m³	R14916
Naphthalene	TO-15	12/26/2007	2.62	1.56	4.1	ND	μg/m³	R14916
o-xylene	TO-15	12/26/2007	0.62062	1.56	0.97	6.4	µg/m³	R14916
Styrene	TO-15	12/26/2007	0.639	1.56	1.0	ND	µg/m³	R14916
Tetrachloroethene	TO-15	12/26/2007	1.2882	1.56	2.0	10	µg/m³	R14916
Tetrahydrofuran	TO-15	12/26/2007	0.413	1.56	0.64	ND	µg/m³	R14916
Toluene	TO-15	12/26/2007	0.5278	1.56	0.82	16	µg/m³	R14916
trans-1,2-Dichloroethene	TO-15	12/26/2007	0.5544	1.56	0.86	ND	µg/m³	R14916
Trichloroethene	TO-15	12/26/2007	0.52626	1.56	0.82	ND	µg/m³	R14916
Trichlorofluoromethane	TO-15	12/26/2007	0.693	1.56	1.1	ND	µg/m³	R14916
Vinyl Acetate	TO-15	12/26/2007	0.64064	1.56	1.0	ND	μg/m³	R14916
Vinyl Chloride	TO-15	12/26/2007	0.24832	1.56	0.39	ND	μg/m³	R14916
Surr: 4-Bromofluorobenzene	TO-15	12/26/2007	0	1.56	65-135	96.2	%REC	R14916

Note: Due to low initial pressure in a Summa canister, sample was diluted prior to the analysis by addition of argon. Results are reported to the MDL.

Report prepared for: Kelly Krohn **Date Received:** 12/17/2007 Engeo

Date Reported: 12/27/2007

Client Sample ID: SG-12 (Duplicate) **Lab Sample ID:** 0712091-003

224 Rickenbacker Circle **Date Prepared: Sample Location:**

Sample Matrix: AIR

Date/Time Sampled 12/17/2007 1:05:00 PM

Parameters	Analysis Method	Date Analyzed	RL	Dilution Factor	MRL	Result	Units	Analytical Batch
1,1 - Dichloroethene	TO-15	12/26/2007	0.794	1.48	1.2	ND	μg/m³	R14916
1,1,1,2-Tetrachloroethane	TO-15	12/26/2007	0.687	1.48	1.0	ND	µg/m³	R14916
1,1,1-Trichloroethane	TO-15	12/26/2007	0.819	1.48	1.2	ND	µg/m³	R14916
1,1,2,2-Tetrachloroethane	TO-15	12/26/2007	1.0305	1.48	1.5	ND	µg/m³	R14916
1,1,2-Trichloroethane	TO-15	12/26/2007	1.0374	1.48	1.5	ND	µg/m³	R14916
1,1-Dichloroethane	TO-15	12/26/2007	0.6885	1.48	1.0	ND	µg/m³	R14916
1,2,4-Trichlorobenzene	TO-15	12/26/2007	0.4984	1.48	0.74	ND	µg/m³	R14916
1,2,4-Trimethylbenzene	TO-15	12/26/2007	0.8856	1.48	1.3	ND	µg/m³	R14916
1,2-Dibromoethane(Ethylene dibromide)	TO-15	12/26/2007	1.0752	1.48	1.6	ND	µg/m³	R14916
1,2-Dichlorobenzene	TO-15	12/26/2007	0.601	1.48	0.89	ND	μg/m³	R14916
1,2-Dichloroethane	TO-15	12/26/2007	0.648	1.48	0.96	ND	μg/m³	R14916
1,2-Dichloropropane	TO-15	12/26/2007	1.0164	1.48	1.5	ND	μg/m³	R14916
1,2- dichlorotetrafluoroethane(F114)	TO-15	12/26/2007	0.9375	1.48	1.4	ND	µg/m³	R14916
1,3,5-Trimethylbenzene	TO-15	12/26/2007	0.6888	1.48	1.0	ND	μg/m³	R14916
1,3-Butadiene	TO-15	12/26/2007	0.5967	1.48	0.88	ND	μg/m³	R14916
1,3-Dichlorobenzene	TO-15	12/26/2007	0.3606	1.48	0.53	ND	μg/m³	R14916
1,4-Dichlorobenzene	TO-15	12/26/2007	0.6611	1.48	0.98	ND	μg/m³	R14916
1,4-Dioxane	TO-15	12/26/2007	0.504	1.48	0.75	ND	μg/m³	R14916
2-Butanone (MEK)	TO-15	12/26/2007	0.4425	1.48	0.65	28	μg/m³	R14916
2-Hexanone	TO-15	12/26/2007	0.861	1.48	1.3	ND	μg/m³	R14916
4-Ethyl Toluene	TO-15	12/26/2007	0.738	1.48	1.1	ND	μg/m³	R14916
4-Methyl-2-Pentanone (MIBK)	TO-15	12/26/2007	0.656	1.48	0.97	ND	μg/m³	R14916
Acetone	TO-15	12/26/2007	0.5712	1.48	0.85	62	μg/m³	R14916
Benzene	TO-15	12/26/2007	0.8932	1.48	1.3	2.22 J	μg/m³	R14916
Benzyl Chloride	TO-15	12/26/2007	0.69	1.48	1.0	ND	μg/m³	R14916
Bromodichloromethane	TO-15	12/26/2007	0.871	1.48	1.3	ND	μg/m³	R14916
Bromoform	TO-15	12/26/2007	1.7578	1.48	2.6	ND	μg/m³	R14916
Bromomethane	TO-15	12/26/2007	0.776	1.48	1.1	ND	μg/m³	R14916
Carbon Disulfide	TO-15	12/26/2007	0.4976	1.48	0.74	ND	µg/m³	R14916
Carbon Tetrachloride	TO-15	12/26/2007	0.9435	1.48	1.4	ND	µg/m³	R14916
Chlorobenzene	TO-15	12/26/2007	0.4232	1.48	0.63	ND	µg/m³	R14916
Chloroethane	TO-15	12/26/2007	0.396	1.48	0.59	ND	µg/m³	R14916
Chloroform	TO-15	12/26/2007	1.952	1.48	2.9	ND	µg/m³	R14916
Chloromethane	TO-15	12/26/2007	0.7245	1.48	1.1	ND	μg/m³	R14916
cis-1,2-dichloroethene	TO-15	12/26/2007	0.5544	1.48	0.82	ND	µg/m³	R14916
cis-1,3-Dichloropropene	TO-15	12/26/2007	0.3632	1.48	0.54	ND	μg/m³	R14916
Dibromochloromethane	TO-15	12/26/2007	0.9372	1.48	1.4	ND	µg/m³	R14916
Dichlorodifluoromethane	TO-15	12/26/2007	0.7425	1.48	1.1	ND	μg/m³	R14916
Ethyl Acetate	TO-15	12/26/2007	0.4248	1.48	0.63	ND	μg/m³	R14916
Ethyl Benzene	TO-15	12/26/2007	0.31062	1.48	0.46	ND	μg/m³	R14916
Freon 113	TO-15	12/26/2007	0.9192	1.48	1.4	ND	μg/m³	R14916
Hexachlorobutadiene	TO-15	12/26/2007	1.8139	1.48	2.7	ND	µg/m³	R14916

Report prepared for: Kelly Krohn **Date Received:** 12/17/2007

Engeo **Date Reported:** 12/27/2007

Client Sample ID: SG-12 (Duplicate) Lab Sample ID: 0712091-003

Sample Location: 224 Rickenbacker Circle Date Prepared:

Sample Matrix: AIR

Date/Time Sampled 12/17/2007 1:05:00 PM

Parameters	Analysis Method	Date Analyzed	RL	Dilution Factor	MRL	Result	Units	Analytical Batch
Hexane	TO-15	12/26/2007	1.7952	1.48	2.7	ND	μg/m³	R14916
Isopropanol	TO-15	12/26/2007	1.6359	1.48	2.4	ND	μg/m³	R14916
m,p-Xylene	TO-15	12/26/2007	0.492	1.48	0.73	21	µg/m³	R14916
Methylene Chloride	TO-15	12/26/2007	0.6859	1.48	1.0	ND	µg/m³	R14916
MTBE	TO-15	12/26/2007	0.5054	1.48	0.75	ND	µg/m³	R14916
Naphthalene	TO-15	12/26/2007	2.62	1.48	3.9	ND	μg/m³	R14916
o-xylene	TO-15	12/26/2007	0.62062	1.48	0.92	5.3	µg/m³	R14916
Styrene	TO-15	12/26/2007	0.639	1.48	0.95	ND	µg/m³	R14916
Tetrachloroethene	TO-15	12/26/2007	1.2882	1.48	1.9	8.7	µg/m³	R14916
Tetrahydrofuran	TO-15	12/26/2007	0.413	1.48	0.61	ND	µg/m³	R14916
Toluene	TO-15	12/26/2007	0.5278	1.48	0.78	14	µg/m³	R14916
trans-1,2-Dichloroethene	TO-15	12/26/2007	0.5544	1.48	0.82	ND	μg/m³	R14916
Trichloroethene	TO-15	12/26/2007	0.52626	1.48	0.78	ND	μg/m³	R14916
Trichlorofluoromethane	TO-15	12/26/2007	0.693	1.48	1.0	ND	μg/m³	R14916
Vinyl Acetate	TO-15	12/26/2007	0.64064	1.48	0.95	ND	μg/m³	R14916
Vinyl Chloride	TO-15	12/26/2007	0.24832	1.48	0.37	ND	μg/m³	R14916
Surr: 4-Bromofluorobenzene	TO-15	12/26/2007	0	1.48	65-135	93.8	%REC	R14916

Note: Due to low initial pressure in a Summa canister, sample was diluted prior to the analysis by addition of argon. Results are reported to the MDL.

Report prepared for: Kelly Krohn **Date Received:** 12/17/2007 Engeo

Date Reported: 12/27/2007

Client Sample ID: SG-13

Sample Location: 224 Rickenbacker Circle

Sample Matrix: AIR

Date/Time Sampled 12/17/2007 2:25:00 PM **Lab Sample ID:** 0712091-004

Date Prepared:

Parameters	Analysis Method	Date Analyzed	RL	Dilution Factor	MRL	Result	Units	Analytical Batch
1,1 - Dichloroethene	TO-15	12/26/2007	0.794	1	0.79	ND	μg/m³	R14916
1,1,1,2-Tetrachloroethane	TO-15	12/26/2007	0.687	1	0.69	ND	μg/m³	R14916
1,1,1-Trichloroethane	TO-15	12/26/2007	0.819	1	0.82	ND	μg/m³	R14916
1,1,2,2-Tetrachloroethane	TO-15	12/26/2007	1.0305	1	1.0	ND	µg/m³	R14916
1,1,2-Trichloroethane	TO-15	12/26/2007	1.0374	1	1.0	ND	µg/m³	R14916
1,1-Dichloroethane	TO-15	12/26/2007	0.6885	1	0.69	ND	μg/m³	R14916
1,2,4-Trichlorobenzene	TO-15	12/26/2007	0.4984	1	0.50	ND	µg/m³	R14916
1,2,4-Trimethylbenzene	TO-15	12/26/2007	0.8856	1	0.89	ND	µg/m³	R14916
1,2-Dibromoethane(Ethylene dibromide)	TO-15	12/26/2007	1.0752	1	1.1	ND	µg/m³	R14916
1,2-Dichlorobenzene	TO-15	12/26/2007	0.601	1	0.60	ND	µg/m³	R14916
1,2-Dichloroethane	TO-15	12/26/2007	0.648	1	0.65	ND	μg/m³	R14916
1,2-Dichloropropane	TO-15	12/26/2007	1.0164	1	1.0	ND	μg/m³	R14916
1,2- dichlorotetrafluoroethane(F114)	TO-15	12/26/2007	0.9375	1	0.94	ND	µg/m³	R14916
1,3,5-Trimethylbenzene	TO-15	12/26/2007	0.6888	1	0.69	ND	µg/m³	R14916
1,3-Butadiene	TO-15	12/26/2007	0.5967	1	0.60	ND	µg/m³	R14916
1,3-Dichlorobenzene	TO-15	12/26/2007	0.3606	1	0.36	ND	µg/m³	R14916
1,4-Dichlorobenzene	TO-15	12/26/2007	0.6611	1	0.66	ND	µg/m³	R14916
1,4-Dioxane	TO-15	12/26/2007	0.504	1	0.50	ND	µg/m³	R14916
2-Butanone (MEK)	TO-15	12/26/2007	0.4425	1	0.44	42	µg/m³	R14916
2-Hexanone	TO-15	12/26/2007	0.861	1	0.86	ND	µg/m³	R14916
4-Ethyl Toluene	TO-15	12/26/2007	0.738	1	0.74	ND	µg/m³	R14916
4-Methyl-2-Pentanone (MIBK)	TO-15	12/26/2007	0.656	1	0.66	ND	µg/m³	R14916
Acetone	TO-15	12/26/2007	0.5712	1	0.57	110	µg/m³	R14916
Benzene	TO-15	12/26/2007	0.8932	1	0.89	3.1	µg/m³	R14916
Benzyl Chloride	TO-15	12/26/2007	0.69	1	0.69	ND	μg/m³	R14916
Bromodichloromethane	TO-15	12/26/2007	0.871	1	0.87	ND	μg/m³	R14916
Bromoform	TO-15	12/26/2007	1.7578	1	1.8	ND	μg/m³	R14916
Bromomethane	TO-15	12/26/2007	0.776	1	0.78	ND	μg/m³	R14916
Carbon Disulfide	TO-15	12/26/2007	0.4976	1	0.50	ND	μg/m³	R14916
Carbon Tetrachloride	TO-15	12/26/2007	0.9435	1	0.94	ND	µg/m³	R14916
Chlorobenzene	TO-15	12/26/2007	0.4232	1	0.42	ND	µg/m³	R14916
Chloroethane	TO-15	12/26/2007	0.396	1	0.40	ND	μg/m³	R14916
Chloroform	TO-15	12/26/2007	1.952	1	2.0	ND	μg/m³	R14916
Chloromethane	TO-15	12/26/2007	0.7245	1	0.72	ND	μg/m³	R14916
cis-1,2-dichloroethene	TO-15	12/26/2007	0.5544	1	0.55	ND	μg/m³	R14916
cis-1,3-Dichloropropene	TO-15	12/26/2007	0.3632	1	0.36	ND	μg/m³	R14916
Dibromochloromethane	TO-15	12/26/2007	0.9372	1	0.94	ND	μg/m³	R14916
Dichlorodifluoromethane	TO-15	12/26/2007	0.7425	1	0.74	ND	μg/m³	R14916
Ethyl Acetate	TO-15	12/26/2007	0.4248	1	0.42	ND	μg/m³	R14916
Ethyl Benzene	TO-15	12/26/2007	0.31062	1	0.31	ND	μg/m³	R14916
Freon 113	TO-15	12/26/2007	0.9192	1	0.92	ND	μg/m³	R14916
Hexachlorobutadiene	TO-15	12/26/2007	1.8139	1	1.8	ND	μg/m³	R14916

These analyses were performed according to State of California Environmental Laboratory Accreditation program, Certificate # 1991

Report prepared for: Kelly Krohn

Engeo

Date Received: 12/17/2007

Date Reported: 12/27/2007

Client Sample ID: SG-13

Lab Sample ID: 0712091-004

Sample Location: 224 Rickenbacker Circle

Date Prepared:

Sample Matrix: AIR

Date/Time Sampled 12/17/2007 2:25:00 PM

Parameters	Analysis Method	Date Analyzed	RL	Dilution Factor	MRL	Result	Units	Analytical Batch
Hexane	TO-15	12/26/2007	1.7952	1	1.8	27	μg/m³	R14916
Isopropanol	TO-15	12/26/2007	1.6359	1	1.6	ND	µg/m³	R14916
m,p-Xylene	TO-15	12/26/2007	0.492	1	0.49	34	µg/m³	R14916
Methylene Chloride	TO-15	12/26/2007	0.6859	1	0.69	ND	µg/m³	R14916
MTBE	TO-15	12/26/2007	0.5054	1	0.51	ND	μg/m³	R14916
Naphthalene	TO-15	12/26/2007	2.62	1	2.6	ND	μg/m³	R14916
o-xylene	TO-15	12/26/2007	0.62062	1	0.62	9.2	µg/m³	R14916
Styrene	TO-15	12/26/2007	0.639	1	0.64	ND	μg/m³	R14916
Tetrachloroethene	TO-15	12/26/2007	1.2882	1	1.3	ND	μg/m³	R14916
Tetrahydrofuran	TO-15	12/26/2007	0.413	1	0.41	ND	μg/m³	R14916
Toluene	TO-15	12/26/2007	0.5278	1	0.53	48	μg/m³	R14916
trans-1,2-Dichloroethene	TO-15	12/26/2007	0.5544	1	0.55	ND	μg/m³	R14916
Trichloroethene	TO-15	12/26/2007	0.52626	1	0.53	ND	μg/m³	R14916
Trichlorofluoromethane	TO-15	12/26/2007	0.693	1	0.69	ND	μg/m³	R14916
Vinyl Acetate	TO-15	12/26/2007	0.64064	1	0.64	ND	μg/m³	R14916
Vinyl Chloride	TO-15	12/26/2007	0.24832	1	0.25	ND	μg/m³	R14916
Surr: 4-Bromofluorobenzene	TO-15	12/26/2007	0	1	65-135	94.2	%REC	R14916

Note: Results reported to the MDL.

Report prepared for: Kelly Krohn **Date Received:** 12/17/2007 Engeo

Date Reported: 12/27/2007

Client Sample ID: SG-14

Sample Location: 224 Rickenbacker Circle

Sample Matrix: AIR

Date/Time Sampled 12/17/2007 11:15:00 AM **Lab Sample ID:** 0712091-005

Date Prepared:

Parameters	Analysis Method	Date Analyzed	RL	Dilution Factor	MRL	Result	Units	Analytical Batch
1,1 - Dichloroethene	TO-15	12/26/2007	0.794	1.57	1.2	ND	μg/m³	R14916
1,1,1,2-Tetrachloroethane	TO-15	12/26/2007	0.687	1.57	1.1	ND	µg/m³	R14916
1,1,1-Trichloroethane	TO-15	12/26/2007	0.819	1.57	1.3	ND	µg/m³	R14916
1,1,2,2-Tetrachloroethane	TO-15	12/26/2007	1.0305	1.57	1.6	ND	µg/m³	R14916
1,1,2-Trichloroethane	TO-15	12/26/2007	1.0374	1.57	1.6	ND	µg/m³	R14916
1,1-Dichloroethane	TO-15	12/26/2007	0.6885	1.57	1.1	ND	µg/m³	R14916
1,2,4-Trichlorobenzene	TO-15	12/26/2007	0.4984	1.57	0.78	ND	µg/m³	R14916
1,2,4-Trimethylbenzene	TO-15	12/26/2007	0.8856	1.57	1.4	2.5 J	µg/m³	R14916
1,2-Dibromoethane(Ethylene dibromide)	TO-15	12/26/2007	1.0752	1.57	1.7	ND	µg/m³	R14916
1,2-Dichlorobenzene	TO-15	12/26/2007	0.601	1.57	0.94	ND	µg/m³	R14916
1,2-Dichloroethane	TO-15	12/26/2007	0.648	1.57	1.0	ND	μg/m³	R14916
1,2-Dichloropropane	TO-15	12/26/2007	1.0164	1.57	1.6	ND	μg/m³	R14916
1,2- dichlorotetrafluoroethane(F114)	TO-15	12/26/2007	0.9375	1.57	1.5	ND	µg/m³	R14916
1,3,5-Trimethylbenzene	TO-15	12/26/2007	0.6888	1.57	1.1	ND	µg/m³	R14916
1,3-Butadiene	TO-15	12/26/2007	0.5967	1.57	0.94	ND	µg/m³	R14916
1,3-Dichlorobenzene	TO-15	12/26/2007	0.3606	1.57	0.57	ND	µg/m³	R14916
1,4-Dichlorobenzene	TO-15	12/26/2007	0.6611	1.57	1.0	ND	µg/m³	R14916
1,4-Dioxane	TO-15	12/26/2007	0.504	1.57	0.79	ND	µg/m³	R14916
2-Butanone (MEK)	TO-15	12/26/2007	0.4425	1.57	0.69	ND	µg/m³	R14916
2-Hexanone	TO-15	12/26/2007	0.861	1.57	1.4	1.7 J	µg/m³	R14916
4-Ethyl Toluene	TO-15	12/26/2007	0.738	1.57	1.2	1.9 J	µg/m³	R14916
4-Methyl-2-Pentanone (MIBK)	TO-15	12/26/2007	0.656	1.57	1.0	ND	µg/m³	R14916
Acetone	TO-15	12/26/2007	0.5712	1.57	0.90	ND	μg/m³	R14916
Benzene	TO-15	12/26/2007	0.8932	1.57	1.4	ND	μg/m³	R14916
Benzyl Chloride	TO-15	12/26/2007	0.69	1.57	1.1	ND	μg/m³	R14916
Bromodichloromethane	TO-15	12/26/2007	0.871	1.57	1.4	ND	μg/m³	R14916
Bromoform	TO-15	12/26/2007	1.7578	1.57	2.8	ND	μg/m³	R14916
Bromomethane	TO-15	12/26/2007	0.776	1.57	1.2	ND	μg/m³	R14916
Carbon Disulfide	TO-15	12/26/2007	0.4976	1.57	0.78	ND	μg/m³	R14916
Carbon Tetrachloride	TO-15	12/26/2007	0.9435	1.57	1.5	ND	μg/m³	R14916
Chlorobenzene	TO-15	12/26/2007	0.4232	1.57	0.66	ND	μg/m³	R14916
Chloroethane	TO-15	12/26/2007	0.396	1.57	0.62	ND	μg/m³	R14916
Chloroform	TO-15	12/26/2007	1.952	1.57	3.1	ND	μg/m³	R14916
Chloromethane	TO-15	12/26/2007	0.7245	1.57	1.1	ND	μg/m³	R14916
cis-1,2-dichloroethene	TO-15	12/26/2007	0.5544	1.57	0.87	ND	μg/m³	R14916
cis-1,3-Dichloropropene	TO-15	12/26/2007	0.3632	1.57	0.57	ND	μg/m³	R14916
Dibromochloromethane	TO-15	12/26/2007	0.9372	1.57	1.5	ND	μg/m³	R14916
Dichlorodifluoromethane	TO-15	12/26/2007	0.7425	1.57	1.2	ND	μg/m³	R14916
Ethyl Acetate	TO-15	12/26/2007	0.4248	1.57	0.67	ND	μg/m³	R14916
Ethyl Benzene	TO-15	12/26/2007	0.31062	1.57	0.49	1.7 J	μg/m³	R14916
Freon 113	TO-15	12/26/2007	0.9192	1.57	1.4	ND	μg/m³	R14916
Hexachlorobutadiene	TO-15	12/26/2007	1.8139	1.57	2.8	ND	μg/m³	R14916

Report prepared for: Kelly Krohn **Date Received:** 12/17/2007

Engeo **Date Reported:** 12/27/2007

Date Prepared:

Client Sample ID: SG-14 Lab Sample ID: 0712091-005

Sample Location: 224 Rickenbacker Circle

Sample Matrix: AIR

Date/Time Sampled 12/17/2007 11:15:00 AM

		1	_	,				
Parameters	Analysis Method	Date Analyzed	RL	Dilution Factor	MRL	Result	Units	Analytical Batch
Hexane	TO-15	12/26/2007	1.7952	1.57	2.8	ND	μg/m³	R14916
Isopropanol	TO-15	12/26/2007	1.6359	1.57	2.6	ND	μg/m³	R14916
m,p-Xylene	TO-15	12/26/2007	0.492	1.57	0.77	8.0	µg/m³	R14916
Methylene Chloride	TO-15	12/26/2007	0.6859	1.57	1.1	ND	µg/m³	R14916
MTBE	TO-15	12/26/2007	0.5054	1.57	0.79	ND	µg/m³	R14916
Naphthalene	TO-15	12/26/2007	2.62	1.57	4.1	ND	µg/m³	R14916
o-xylene	TO-15	12/26/2007	0.62062	1.57	0.97	ND	µg/m³	R14916
Styrene	TO-15	12/26/2007	0.639	1.57	1.0	ND	µg/m³	R14916
Tetrachloroethene	TO-15	12/26/2007	1.2882	1.57	2.0	ND	µg/m³	R14916
Tetrahydrofuran	TO-15	12/26/2007	0.413	1.57	0.65	ND	µg/m³	R14916
Toluene	TO-15	12/26/2007	0.5278	1.57	0.83	3.3	µg/m³	R14916
trans-1,2-Dichloroethene	TO-15	12/26/2007	0.5544	1.57	0.87	ND	µg/m³	R14916
Trichloroethene	TO-15	12/26/2007	0.52626	1.57	0.83	ND	µg/m³	R14916
Trichlorofluoromethane	TO-15	12/26/2007	0.693	1.57	1.1	ND	µg/m³	R14916
Vinyl Acetate	TO-15	12/26/2007	0.64064	1.57	1.0	ND	µg/m³	R14916
Vinyl Chloride	TO-15	12/26/2007	0.24832	1.57	0.39	ND	µg/m³	R14916
Surr: 4-Bromofluorobenzene	TO-15	12/26/2007	0	1.57	65-135	88.4	%REC	R14916

Note: Due to low initial pressure in a Summa canister, sample was diluted prior to the analysis by addition of argon. Results are reported to the MDL.

Report prepared for: Kelly Krohn
Engeo

Date Received: 12/17/2007 **Date Reported:** 12/27/2007

Client Sample ID: SG-15

Lab Sample ID: 0712091-006

Sample Location: 224 Rickenbacker Circle

Date Prepared:

Sample Matrix: AIR

Date/Time Sampled 12/17/2007 1:57:00 PM

Parameters	Analysis Method	Date Analyzed	RL	Dilution Factor	MRL	Result	Units	Analytical Batch
1,1 - Dichloroethene	TO-15	12/26/2007	0.794	1.47	1.2	ND	μg/m³	R14916
1,1,1,2-Tetrachloroethane	TO-15	12/26/2007	0.687	1.47	1.0	ND	μg/m³	R14916
1,1,1-Trichloroethane	TO-15	12/26/2007	0.819	1.47	1.2	ND	μg/m³	R14916
1,1,2,2-Tetrachloroethane	TO-15	12/26/2007	1.0305	1.47	1.5	ND	μg/m³	R14916
1,1,2-Trichloroethane	TO-15	12/26/2007	1.0374	1.47	1.5	ND	μg/m³	R14916
1,1-Dichloroethane	TO-15	12/26/2007	0.6885	1.47	1.0	ND	μg/m³	R14916
1,2,4-Trichlorobenzene	TO-15	12/26/2007	0.4984	1.47	0.73	ND	μg/m³	R14916
1,2,4-Trimethylbenzene	TO-15	12/26/2007	0.8856	1.47	1.3	ND	µg/m³	R14916
1,2-Dibromoethane(Ethylene dibromide)	TO-15	12/26/2007	1.0752	1.47	1.6	ND	µg/m³	R14916
1,2-Dichlorobenzene	TO-15	12/26/2007	0.601	1.47	0.88	ND	μg/m³	R14916
1,2-Dichloroethane	TO-15	12/26/2007	0.648	1.47	0.95	ND	μg/m³	R14916
1,2-Dichloropropane	TO-15	12/26/2007	1.0164	1.47	1.5	ND	μg/m³	R14916
1,2- dichlorotetrafluoroethane(F114)	TO-15	12/26/2007	0.9375	1.47	1.4	ND	μg/m³	R14916
1,3,5-Trimethylbenzene	TO-15	12/26/2007	0.6888	1.47	1.0	ND	µg/m³	R14916
1,3-Butadiene	TO-15	12/26/2007	0.5967	1.47	0.88	ND	μg/m³	R14916
1,3-Dichlorobenzene	TO-15	12/26/2007	0.3606	1.47	0.53	ND	μg/m³	R14916
1,4-Dichlorobenzene	TO-15	12/26/2007	0.6611	1.47	0.97	ND	μg/m³	R14916
1,4-Dioxane	TO-15	12/26/2007	0.504	1.47	0.74	ND	μg/m³	R14916
2-Butanone (MEK)	TO-15	12/26/2007	0.4425	1.47	0.65	29	μg/m³	R14916
2-Hexanone	TO-15	12/26/2007	0.861	1.47	1.3	ND	μg/m³	R14916
4-Ethyl Toluene	TO-15	12/26/2007	0.738	1.47	1.1	ND	µg/m³	R14916
4-Methyl-2-Pentanone (MIBK)	TO-15	12/26/2007	0.656	1.47	0.96	ND	µg/m³	R14916
Acetone	TO-15	12/26/2007	0.5712	1.47	0.84	200	µg/m³	R14916
Benzene	TO-15	12/26/2007	0.8932	1.47	1.3	4.0	µg/m³	R14916
Benzyl Chloride	TO-15	12/26/2007	0.69	1.47	1.0	ND	µg/m³	R14916
Bromodichloromethane	TO-15	12/26/2007	0.871	1.47	1.3	ND	μg/m³	R14916
Bromoform	TO-15	12/26/2007	1.7578	1.47	2.6	ND	μg/m³	R14916
Bromomethane	TO-15	12/26/2007	0.776	1.47	1.1	ND	μg/m³	R14916
Carbon Disulfide	TO-15	12/26/2007	0.4976	1.47	0.73	ND	µg/m³	R14916
Carbon Tetrachloride	TO-15	12/26/2007	0.9435	1.47	1.4	ND	µg/m³	R14916
Chlorobenzene	TO-15	12/26/2007	0.4232	1.47	0.62	ND	µg/m³	R14916
Chloroethane	TO-15	12/26/2007	0.396	1.47	0.58	ND	μg/m³	R14916
Chloroform	TO-15	12/26/2007	1.952	1.47	2.9	ND	μg/m³	R14916
Chloromethane	TO-15	12/26/2007	0.7245	1.47	1.1	ND	μg/m³	R14916
cis-1,2-dichloroethene	TO-15	12/26/2007	0.5544	1.47	0.81	ND	μg/m³	R14916
cis-1,3-Dichloropropene	TO-15	12/26/2007	0.3632	1.47	0.53	ND	μg/m³	R14916
Dibromochloromethane	TO-15	12/26/2007	0.9372	1.47	1.4	ND	μg/m³	R14916
Dichlorodifluoromethane	TO-15	12/26/2007	0.7425	1.47	1.1	ND	μg/m³	R14916
Ethyl Acetate	TO-15	12/26/2007	0.4248	1.47	0.62	ND	μg/m³	R14916
Ethyl Benzene	TO-15	12/26/2007	0.31062	1.47	0.46	ND	μg/m³	R14916
Freon 113	TO-15	12/26/2007	0.9192	1.47	1.4	ND	μg/m³	R14916
Hexachlorobutadiene	TO-15	12/26/2007	1.8139	1.47	2.7	ND	μg/m³	R14916

These analyses were performed according to State of California Environmental Laboratory Accreditation program, Certificate # 1991

Report prepared for: Kelly Krohn

224 Rickenbacker Circle

Engeo

Date Received: 12/17/2007

Date Reported: 12/27/2007

Client Sample ID: SG-15

OU-13

Sample Matrix: AIR

Sample Location:

Date/Time Sampled 12/17/2007 1:57:00 PM

Lab Sample ID: 0712091-006

Date Prepared:

Parameters	Analysis Method	Date Analyzed	RL	Dilution Factor	MRL	Result	Units	Analytical Batch
Hexane	TO-15	12/26/2007	1.7952	1.47	2.6	59	μg/m³	R14916
Isopropanol	TO-15	12/26/2007	1.6359	1.47	2.4	ND	μg/m³	R14916
m,p-Xylene	TO-15	12/26/2007	0.492	1.47	0.72	39	μg/m³	R14916
Methylene Chloride	TO-15	12/26/2007	0.6859	1.47	1.0	ND	μg/m³	R14916
MTBE	TO-15	12/26/2007	0.5054	1.47	0.74	ND	μg/m³	R14916
Naphthalene	TO-15	12/26/2007	2.62	1.47	3.9	ND	μg/m³	R14916
o-xylene	TO-15	12/26/2007	0.62062	1.47	0.91	11	μg/m³	R14916
Styrene	TO-15	12/26/2007	0.639	1.47	0.94	ND	µg/m³	R14916
Tetrachloroethene	TO-15	12/26/2007	1.2882	1.47	1.9	ND	µg/m³	R14916
Tetrahydrofuran	TO-15	12/26/2007	0.413	1.47	0.61	ND	µg/m³	R14916
Toluene	TO-15	12/26/2007	0.5278	1.47	0.78	68	µg/m³	R14916
trans-1,2-Dichloroethene	TO-15	12/26/2007	0.5544	1.47	0.81	ND	µg/m³	R14916
Trichloroethene	TO-15	12/26/2007	0.52626	1.47	0.77	ND	μg/m³	R14916
Trichlorofluoromethane	TO-15	12/26/2007	0.693	1.47	1.0	ND	μg/m³	R14916
Vinyl Acetate	TO-15	12/26/2007	0.64064	1.47	0.94	ND	μg/m³	R14916
Vinyl Chloride	TO-15	12/26/2007	0.24832	1.47	0.37	ND	μg/m³	R14916
Surr: 4-Bromofluorobenzene	TO-15	12/26/2007	0	1.47	65-135	86.3	%REC	R14916

Note: Due to low initial pressure in a Summa canister, sample was diluted prior to the analysis by addition of argon. Results are reported to the MDL.

Report prepared for: Kelly Krohn
Engeo

Date Received: 12/17/2007 **Date Reported:** 12/27/2007

Client Sample ID: SG-16

Sample Location: 224 Rickenbacker Circle

Sample Matrix: AIR

Date/Time Sampled 12/17/2007 12:36:00 PM

Lab Sample ID: 0712091-007

Date Prepared:

Parameters	Analysis Method	Date Analyzed	RL	Dilution Factor	MRL	Result	Units	Analytical Batch
1,1 - Dichloroethene	TO-15	12/26/2007	0.794	1.5	1.2	ND	μg/m³	R14916
1,1,1,2-Tetrachloroethane	TO-15	12/26/2007	0.687	1.5	1.0	ND	μg/m³	R14916
1,1,1-Trichloroethane	TO-15	12/26/2007	0.819	1.5	1.2	ND	μg/m³	R14916
1,1,2,2-Tetrachloroethane	TO-15	12/26/2007	1.0305	1.5	1.5	ND	μg/m³	R14916
1,1,2-Trichloroethane	TO-15	12/26/2007	1.0374	1.5	1.6	ND	μg/m³	R14916
1,1-Dichloroethane	TO-15	12/26/2007	0.6885	1.5	1.0	ND	μg/m³	R14916
1,2,4-Trichlorobenzene	TO-15	12/26/2007	0.4984	1.5	0.75	ND	μg/m³	R14916
1,2,4-Trimethylbenzene	TO-15	12/26/2007	0.8856	1.5	1.3	ND	μg/m³	R14916
1,2-Dibromoethane(Ethylene dibromide)	TO-15	12/26/2007	1.0752	1.5	1.6	ND	µg/m³	R14916
1,2-Dichlorobenzene	TO-15	12/26/2007	0.601	1.5	0.90	ND	μg/m³	R14916
1,2-Dichloroethane	TO-15	12/26/2007	0.648	1.5	0.97	ND	μg/m³	R14916
1,2-Dichloropropane	TO-15	12/26/2007	1.0164	1.5	1.5	ND	μg/m³	R14916
1,2-	TO-15	12/26/2007	0.9375	1.5	1.4	ND	μg/m³	R14916
dichlorotetrafluoroethane(F114)	10 10	12/20/2001	0.0070	1.0	17	NB	рулп	1(14010
1,3,5-Trimethylbenzene	TO-15	12/26/2007	0.6888	1.5	1.0	ND	µg/m³	R14916
1,3-Butadiene	TO-15	12/26/2007	0.5967	1.5	0.90	ND	µg/m³	R14916
1,3-Dichlorobenzene	TO-15	12/26/2007	0.3606	1.5	0.54	ND	μg/m³	R14916
1,4-Dichlorobenzene	TO-15	12/26/2007	0.6611	1.5	0.99	ND	μg/m³	R14916
1,4-Dioxane	TO-15	12/26/2007	0.504	1.5	0.76	ND	μg/m³	R14916
2-Butanone (MEK)	TO-15	12/26/2007	0.4425	1.5	0.66	110	μg/m³	R14916
2-Hexanone	TO-15	12/26/2007	0.861	1.5	1.3	ND	μg/m³	R14916
4-Ethyl Toluene	TO-15	12/26/2007	0.738	1.5	1.1	ND	μg/m³	R14916
4-Methyl-2-Pentanone (MIBK)	TO-15	12/26/2007	0.656	1.5	0.98	ND	μg/m³	R14916
Acetone	TO-15	12/26/2007	0.5712	1.5	0.86	180	μg/m³	R14916
Benzene	TO-15	12/26/2007	0.8932	1.5	1.3	6.6	μg/m³	R14916
Benzyl Chloride	TO-15	12/26/2007	0.69	1.5	1.0	ND	μg/m³	R14916
Bromodichloromethane	TO-15	12/26/2007	0.871	1.5	1.3	ND	μg/m³	R14916
Bromoform	TO-15	12/26/2007	1.7578	1.5	2.6	ND	μg/m³	R14916
Bromomethane	TO-15	12/26/2007	0.776	1.5	1.2	ND	μg/m³	R14916
Carbon Disulfide	TO-15	12/26/2007	0.4976	1.5	0.75	ND	μg/m³	R14916
Carbon Tetrachloride	TO-15	12/26/2007	0.9435	1.5	1.4	ND	μg/m³	R14916
Chlorobenzene	TO-15	12/26/2007	0.4232	1.5	0.63	ND	μg/m³	R14916
Chloroethane	TO-15	12/26/2007	0.396	1.5	0.59	ND	μg/m³	R14916
Chloroform	TO-15	12/26/2007	1.952	1.5	2.9	ND	μg/m³	R14916
Chloromethane	TO-15	12/26/2007	0.7245	1.5	1.1	ND	μg/m³	R14916
cis-1,2-dichloroethene	TO-15	12/26/2007	0.5544	1.5	0.83	7.9	μg/m³	R14916
cis-1,3-Dichloropropene	TO-15	12/26/2007	0.3632	1.5	0.54	ND	μg/m³	R14916
Dibromochloromethane	TO-15	12/26/2007	0.3032	1.5	1.4	ND	μg/m³	R14916 R14916
Dichlorodifluoromethane	TO-15	12/26/2007	0.9372	1.5	1.4	ND	μg/m³	R14916 R14916
	TO-15	12/26/2007	0.7425	1.5	0.64	ND		R14916 R14916
Ethyl Ronzona							μg/m³	
Ethyl Benzene	TO-15	12/26/2007	0.31062	1.5	0.47	8.2 ND	μg/m³	R14916
Freon 113	TO-15	12/26/2007	0.9192	1.5	1.4	ND	μg/m³	R14916
Hexachlorobutadiene	TO-15	12/26/2007	1.8139	1.5	2.7	ND	μg/m³	R14916

Report prepared for: Kelly Krohn

Engeo

Date Received: 12/17/2007

Date Reported: 12/27/2007

Client Sample ID: SG-16

224 Rickenbacker Circle

Sample Matrix: AIR

Sample Location:

Date/Time Sampled 12/17/2007 12:36:00 PM

Lab Sample ID: 0712091-007

Date Prepared:

Parameters	Analysis Method	Date Analyzed	RL	Dilution Factor	MRL	Result	Units	Analytical Batch
Hexane	TO-15	12/26/2007	1.7952	1.5	2.7	ND	μg/m³	R14916
Isopropanol	TO-15	12/26/2007	1.6359	1.5	2.5	ND	μg/m³	R14916
m,p-Xylene	TO-15	12/26/2007	0.492	1.5	0.74	46	μg/m³	R14916
Methylene Chloride	TO-15	12/26/2007	0.6859	1.5	1.0	ND	μg/m³	R14916
MTBE	TO-15	12/26/2007	0.5054	1.5	0.76	ND	μg/m³	R14916
Naphthalene	TO-15	12/26/2007	2.62	1.5	3.9	ND	μg/m³	R14916
o-xylene	TO-15	12/26/2007	0.62062	1.5	0.93	13	μg/m³	R14916
Styrene	TO-15	12/26/2007	0.639	1.5	0.96	ND	µg/m³	R14916
Tetrachloroethene	TO-15	12/26/2007	1.2882	1.5	1.9	15	µg/m³	R14916
Tetrahydrofuran	TO-15	12/26/2007	0.413	1.5	0.62	ND	µg/m³	R14916
Toluene	TO-15	12/26/2007	0.5278	1.5	0.79	30	µg/m³	R14916
trans-1,2-Dichloroethene	TO-15	12/26/2007	0.5544	1.5	0.83	8.2	µg/m³	R14916
Trichloroethene	TO-15	12/26/2007	0.52626	1.5	0.79	22	µg/m³	R14916
Trichlorofluoromethane	TO-15	12/26/2007	0.693	1.5	1.0	ND	µg/m³	R14916
Vinyl Acetate	TO-15	12/26/2007	0.64064	1.5	0.96	ND	µg/m³	R14916
Vinyl Chloride	TO-15	12/26/2007	0.24832	1.5	0.37	ND	μg/m³	R14916
Surr: 4-Bromofluorobenzene	TO-15	12/26/2007	0	1.5	65-135	90.7	%REC	R14916

Note: Due to low initial pressure in a Summa canister, sample was diluted prior to the analysis by addition of argon. Results are reported to the MDL.

Report prepared for: Kelly Krohn **Date Received:** 12/17/2007 Engeo

Date Reported: 12/27/2007

Client Sample ID: SG-10

Sample Location: 224 Rickenbacker Circle

Sample Matrix: AIR

Date/Time Sampled 12/17/2007 2:57:00 PM **Lab Sample ID:** 0712091-008

Date Prepared:

Parameters	Analysis Method	Date Analyzed	RL	Dilution Factor	MRL	Result	Units	Analytical Batch
1,1 - Dichloroethene	TO-15	12/26/2007	0.794	1.63	1.3	ND	μg/m³	R14916
1,1,1,2-Tetrachloroethane	TO-15	12/26/2007	0.687	1.63	1.1	ND	µg/m³	R14916
1,1,1-Trichloroethane	TO-15	12/26/2007	0.819	1.63	1.3	ND	µg/m³	R14916
1,1,2,2-Tetrachloroethane	TO-15	12/26/2007	1.0305	1.63	1.7	ND	µg/m³	R14916
1,1,2-Trichloroethane	TO-15	12/26/2007	1.0374	1.63	1.7	ND	µg/m³	R14916
1,1-Dichloroethane	TO-15	12/26/2007	0.6885	1.63	1.1	ND	μg/m³	R14916
1,2,4-Trichlorobenzene	TO-15	12/26/2007	0.4984	1.63	0.81	ND	µg/m³	R14916
1,2,4-Trimethylbenzene	TO-15	12/26/2007	0.8856	1.63	1.4	ND	μg/m³	R14916
1,2-Dibromoethane(Ethylene dibromide)	TO-15	12/26/2007	1.0752	1.63	1.8	ND	μg/m³	R14916
1,2-Dichlorobenzene	TO-15	12/26/2007	0.601	1.63	0.98	ND	µg/m³	R14916
1,2-Dichloroethane	TO-15	12/26/2007	0.648	1.63	1.1	ND	µg/m³	R14916
1,2-Dichloropropane	TO-15	12/26/2007	1.0164	1.63	1.7	ND	µg/m³	R14916
1,2- dichlorotetrafluoroethane(F114)	TO-15	12/26/2007	0.9375	1.63	1.5	ND	µg/m³	R14916
1,3,5-Trimethylbenzene	TO-15	12/26/2007	0.6888	1.63	1.1	ND	μg/m³	R14916
1,3-Butadiene	TO-15	12/26/2007	0.5967	1.63	0.97	ND	µg/m³	R14916
1,3-Dichlorobenzene	TO-15	12/26/2007	0.3606	1.63	0.59	ND	µg/m³	R14916
1,4-Dichlorobenzene	TO-15	12/26/2007	0.6611	1.63	1.1	ND	µg/m³	R14916
1,4-Dioxane	TO-15	12/26/2007	0.504	1.63	0.82	ND	µg/m³	R14916
2-Butanone (MEK)	TO-15	12/26/2007	0.4425	1.63	0.72	58	µg/m³	R14916
2-Hexanone	TO-15	12/26/2007	0.861	1.63	1.4	5.3	µg/m³	R14916
4-Ethyl Toluene	TO-15	12/26/2007	0.738	1.63	1.2	ND	μg/m³	R14916
4-Methyl-2-Pentanone (MIBK)	TO-15	12/26/2007	0.656	1.63	1.1	ND	μg/m³	R14916
Acetone	TO-15	12/26/2007	0.5712	1.63	0.93	100	µg/m³	R14916
Benzene	TO-15	12/26/2007	0.8932	1.63	1.5	2.8	µg/m³	R14916
Benzyl Chloride	TO-15	12/26/2007	0.69	1.63	1.1	ND	μg/m³	R14916
Bromodichloromethane	TO-15	12/26/2007	0.871	1.63	1.4	ND	μg/m³	R14916
Bromoform	TO-15	12/26/2007	1.7578	1.63	2.9	ND	μg/m³	R14916
Bromomethane	TO-15	12/26/2007	0.776	1.63	1.3	ND	μg/m³	R14916
Carbon Disulfide	TO-15	12/26/2007	0.4976	1.63	0.81	ND	µg/m³	R14916
Carbon Tetrachloride	TO-15	12/26/2007	0.9435	1.63	1.5	ND	µg/m³	R14916
Chlorobenzene	TO-15	12/26/2007	0.4232	1.63	0.69	ND	μg/m³	R14916
Chloroethane	TO-15	12/26/2007	0.396	1.63	0.65	ND	μg/m³	R14916
Chloroform	TO-15	12/26/2007	1.952	1.63	3.2	ND	μg/m³	R14916
Chloromethane	TO-15	12/26/2007	0.7245	1.63	1.2	ND	μg/m³	R14916
cis-1,2-dichloroethene	TO-15	12/26/2007	0.5544	1.63	0.90	ND	μg/m³	R14916
cis-1,3-Dichloropropene	TO-15	12/26/2007	0.3632	1.63	0.59	ND	µg/m³	R14916
Dibromochloromethane	TO-15	12/26/2007	0.9372	1.63	1.5	ND	μg/m³	R14916
Dichlorodifluoromethane	TO-15	12/26/2007	0.7425	1.63	1.2	ND	μg/m³	R14916
Ethyl Acetate	TO-15	12/26/2007	0.4248	1.63	0.69	ND	μg/m³	R14916
Ethyl Benzene	TO-15	12/26/2007	0.31062	1.63	0.51	ND	μg/m³	R14916
Freon 113	TO-15	12/26/2007	0.9192	1.63	1.5	ND	μg/m³	R14916
Hexachlorobutadiene	TO-15	12/26/2007	1.8139	1.63	3.0	ND	μg/m³	R14916

Report prepared for: Kelly Krohn **Date Received:** 12/17/2007

Engeo **Date Reported:** 12/27/2007

Client Sample ID: SG-10 Lab Sample ID: 0712091-008

Sample Location: 224 Rickenbacker Circle Date Prepared:

Sample Matrix: AIR

Date/Time Sampled 12/17/2007 2:57:00 PM

Parameters	Analysis Method	Date Analyzed	RL	Dilution Factor	MRL	Result	Units	Analytical Batch
Hexane	TO-15	12/26/2007	1.7952	1.63	2.9	17	μg/m³	R14916
Isopropanol	TO-15	12/26/2007	1.6359	1.63	2.7	ND	µg/m³	R14916
m,p-Xylene	TO-15	12/26/2007	0.492	1.63	0.80	38	µg/m³	R14916
Methylene Chloride	TO-15	12/26/2007	0.6859	1.63	1.1	ND	μg/m³	R14916
MTBE	TO-15	12/26/2007	0.5054	1.63	0.82	ND	μg/m³	R14916
Naphthalene	TO-15	12/26/2007	2.62	1.63	4.3	ND	μg/m³	R14916
o-xylene	TO-15	12/26/2007	0.62062	1.63	1.0	10	μg/m³	R14916
Styrene	TO-15	12/26/2007	0.639	1.63	1.0	ND	μg/m³	R14916
Tetrachloroethene	TO-15	12/26/2007	1.2882	1.63	2.1	ND	μg/m³	R14916
Tetrahydrofuran	TO-15	12/26/2007	0.413	1.63	0.67	ND	μg/m³	R14916
Toluene	TO-15	12/26/2007	0.5278	1.63	0.86	31	μg/m³	R14916
trans-1,2-Dichloroethene	TO-15	12/26/2007	0.5544	1.63	0.90	ND	μg/m³	R14916
Trichloroethene	TO-15	12/26/2007	0.52626	1.63	0.86	ND	μg/m³	R14916
Trichlorofluoromethane	TO-15	12/26/2007	0.693	1.63	1.1	ND	μg/m³	R14916
Vinyl Acetate	TO-15	12/26/2007	0.64064	1.63	1.0	ND	μg/m³	R14916
Vinyl Chloride	TO-15	12/26/2007	0.24832	1.63	0.40	ND	μg/m³	R14916
Surr: 4-Bromofluorobenzene	TO-15	12/26/2007	0	1.63	65-150	86.8	%REC	R14916

Note: Due to low initial pressure in a Summa canister, sample was diluted prior to the analysis by addition of argon. Results are reported to the MDL.

Definitions, legends and Notes

Note	Description
ug/kg	Microgram per kilogram (ppb, part per billion).
ug/L	Microgram per liter (ppb, part per billion).
mg/kg	Milligram per kilogram (ppm, part per million).
mg/L	Milligram per liter (ppm, part per million).
LCS/LCSD	Laboratory control sample/laboratory control sample duplicate.
MDL	Method detection limit.
MRL	Modified reporting limit. When sample is subject to dilution, reporting limit times dilution factor yields MRL.
MS/MSD	Matrix spike/matrix spike duplicate.
N/A	Not applicable.
ND	Not detected at or above detection limit.
NR	Not reported.
QC	Quality Control.
RL	Reporting limit.
% RPD	Percent relative difference.
а	pH was measured immediately upon the receipt of the sample, but it was still done outside the holding time.
sub	Analyzed by subcontracting laboratory, Lab Certificate #

Date: 27-Dec-07

CLIENT: Engeo **Work Order:** 0712091

Project: 7584.1.001.01

ANALYTICAL QC SUMMARY REPORT

BatchID: R14916

Sample ID mb	SampType:	MBLK	TestCoo	le: TO-15	Units: ppbv		Prep Da	te: 12/26/ 2	2007	RunNo: 149	916	
Client ID: ZZZZZ	Batch ID:	R14916	TestN	lo: TO-15			Analysis Da	te: 12/26/ 2	2007	SeqNo: 214	1608	
Analyte		Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1 - Dichloroethene		ND	0.20									
1,1,1,2-Tetrachloroethane		ND	0.10									
1,1,1-Trichloroethane		ND	0.15									
1,1,2,2-Tetrachloroethane		ND	0.15									
1,1,2-Trichloroethane		ND	0.19									
1,1-Dichloroethane		ND	0.17									
1,2,4-Trichlorobenzene		ND	0.070									
1,2,4-Trimethylbenzene		ND	0.18									
1,2-Dibromoethane(Ethylene dibro	omide	ND	0.14									
1,2-Dichlorobenzene		ND	0.10									
1,2-Dichloroethane		ND	0.16									
1,2-Dichloropropane		ND	0.22									
1,2-dichlorotetrafluoroethane(F11	4)	ND	0.15									
1,3,5-Trimethylbenzene		ND	0.14									
1,3-Butadiene		ND	0.27									
1,3-Dichlorobenzene		ND	0.060									
1,4-Dichlorobenzene		ND	0.11									
1,4-Dioxane		ND	0.14									
2-Butanone (MEK)		ND	0.15									
2-Hexanone		ND	0.21									
4-Ethyl Toluene		ND	0.15									
4-Methyl-2-Pentanone (MIBK)		ND	0.16									
Acetone		ND	0.24									
Benzene		ND	0.28									
Benzyl Chloride		ND	0.12									
Bromodichloromethane		ND	0.13									
Bromoform		ND	0.17									
Bromomethane		ND	0.20									
Carbon Disulfide		ND	0.16									
Carbon Tetrachloride		ND	0.15									

Qualifiers:

Value above quantitation range

ND Not Detected at the Reporting Limit

Holding times for preparation or analysis exceeded

RPD outside accepted recovery limits

Analyte detected below quantitation limits

Spike Recovery outside accepted recovery limits

Page 1 of 6

CLIENT: Engeo **Work Order:**

0712091

Project: 7584.1.001.01

ANALYTICAL QC SUMMARY REPORT

BatchID: R14916

Sample ID mb	SampType: MBLK	TestCod	de: TO-15	Units: ppbv		Prep Da	te: 12/26/ 2	2007	RunNo: 14 9	916	
Client ID: ZZZZZ	Batch ID: R14916	TestN	No: TO-15			Analysis Da	ite: 12/26/ 2	2007	SeqNo: 21	4608	
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
Chlorobenzene	ND	0.092									
Chloroethane	ND	0.15									
Chloroform	ND	0.40									
Chloromethane	ND	0.35									
cis-1,2-dichloroethene	ND	0.14									
cis-1,3-Dichloropropene	ND	0.080									
Dibromochloromethane	ND	0.11									
Dichlorodifluoromethane	ND	0.15									
Ethyl Acetate	ND	0.12									
Ethyl Benzene	ND	0.093									
Freon 113	ND	0.12									
Hexachlorobutadiene	ND	0.17									
Hexane	ND	0.51									
Isopropanol	ND	0.40									
m,p-Xylene	ND	0.12									
Methylene Chloride	ND	0.19									
MTBE	ND	0.14									
Naphthalene	ND	0.50									
o-xylene	ND	0.14									
Styrene	ND	0.15									
Tetrachloroethene	ND	0.19									
Tetrahydrofuran	ND	0.14									
Toluene	ND	0.14									
trans-1,2-Dichloroethene	ND	0.14									
Trichloroethene	ND	0.098									
Trichlorofluoromethane	ND	0.14									
Vinyl Acetate	ND	0.18									
Vinyl Chloride	ND	0.097									
Surr: 4-Bromofluorobenzene	18.14	0	20	0	90.7	65	135				

Value above quantitation range Qualifiers:

ND Not Detected at the Reporting Limit

Holding times for preparation or analysis exceeded

RPD outside accepted recovery limits

Analyte detected below quantitation limits

Spike Recovery outside accepted recovery limits

Page 2 of 6

CLIENT: Engeo

Work Order: 0712091

Project: 7584.1.001.01

ANALYTICAL QC SUMMARY REPORT

BatchID: R14916

Sample ID LCS	SampType: LCS	TestCo	de: TO-15	Units: ppbv		Prep Da	te: 12/26/2	2007	RunNo: 14	916	
Client ID: ZZZZZ	Batch ID: R14916	Test	No: TO-15			Analysis Da	te: 12/26/2	2007	SeqNo: 21	4609	
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1 - Dichloroethene	20.56	0.20	20	0	103	65	135				
1,1,1,2-Tetrachloroethane	23.18	0.10	20	0	116	65	135				
1,1,1-Trichloroethane	20.37	0.15	20	0	102	65	135				
1,1,2,2-Tetrachloroethane	23.48	0.15	20	0	117	65	135				
1,1,2-Trichloroethane	23.08	0.19	20	0	115	65	135				
1,1-Dichloroethane	20.54	0.17	20	0	103	65	135				
1,2,4-Trichlorobenzene	20.15	0.070	20	0	101	65	135				
1,2,4-Trimethylbenzene	22.24	0.18	20	0	111	65	135				
1,2-Dibromoethane(Ethylene dibron	nide 22.31	0.14	20	0	112	65	135				
1,2-Dichlorobenzene	21.85	0.10	20	0	109	65	135				
1,2-Dichloroethane	23.54	0.16	20	0	118	65	135				
1,2-Dichloropropane	24.14	0.22	20	0	121	65	135				
1,2-dichlorotetrafluoroethane(F114)	19.79	0.15	20	0	99.0	65	135				
1,3,5-Trimethylbenzene	23.15	0.14	20	0	116	65	135				
1,3-Butadiene	20.86	0.27	20	0	104	65	135				
1,3-Dichlorobenzene	21.84	0.060	20	0	109	65	135				
1,4-Dichlorobenzene	21.84	0.11	20	0	109	65	135				
1,4-Dioxane	23.83	0.14	20	0	119	65	135				
2-Butanone (MEK)	22.23	0.15	20	0	111	65	135				
2-Hexanone	24.32	0.21	20	0	122	65	135				
4-Ethyl Toluene	22.61	0.15	20	0	113	65	135				
4-Methyl-2-Pentanone (MIBK)	23.39	0.16	20	0	117	65	135				
Acetone	17.46	0.24	20	0	87.3	65	135				
Benzene	19.99	0.28	20	0	100	65	135				
Benzyl Chloride	22.24	0.12	20	0	111	65	135				
Bromodichloromethane	23.53	0.13	20	0	118	65	135				
Bromoform	24.06	0.17	20	0	120	65	135				
Bromomethane	20.49	0.20	20	0	102	65	135				
Carbon Disulfide	21.53	0.16	20	0	108	65	135				
Carbon Tetrachloride	20.09	0.15	20	0	100	65	135				
Chlorobenzene	23.12	0.092	20	0	116	65	135				

Value above quantitation range Qualifiers:

ND Not Detected at the Reporting Limit

Holding times for preparation or analysis exceeded

RPD outside accepted recovery limits

Analyte detected below quantitation limits

Spike Recovery outside accepted recovery limits

Page 3 of 6

CLIENT: Engeo

Work Order: 0712091

Project: 7584.1.001.01

ANALYTICAL QC SUMMARY REPORT

BatchID: R14916

Sample ID LCS	SampType: LCS	TestCod	e: TO-15	Units: ppbv	Prep Date: 12/26/2007			007	RunNo: 14	916	
Client ID: ZZZZZ	Batch ID: R14916	TestN	o: TO-15			Analysis Date	12/26/2	007	SeqNo: 21	4609	
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit I	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
Chloroethane	20.33	0.15	20	0	102	65	135				
Chloroform	20.46	0.40	20	0	102	65	135				
Chloromethane	25.88	0.35	20	0	129	65	135				
cis-1,2-dichloroethene	20.11	0.14	20	0	101	65	135				
cis-1,3-Dichloropropene	22.81	0.080	20	0	114	65	135				
Dibromochloromethane	23.07	0.11	20	0	115	65	135				
Ethyl Acetate	22.04	0.12	20	0	110	65	135				
Ethyl Benzene	22.36	0.093	20	0	112	65	135				
Freon 113	21.48	0.12	20	0	107	65	135				
Hexachlorobutadiene	19.66	0.17	20	0	98.3	65	135				
Hexane	20.52	0.51	20	0	103	65	135				
Isopropanol	22.21	0.40	20	0	111	65	135				
m,p-Xylene	44.02	0.12	40	0	110	65	135				
Methylene Chloride	20.88	0.19	20	0	104	65	135				
MTBE	20.82	0.14	20	0	104	65	135				
Naphthalene	20.49	0.50	20	0	102	65	135				
o-xylene	21.47	0.14	20	0	107	65	135				
Styrene	22.29	0.15	20	0	111	65	135				
Tetrachloroethene	22.83	0.19	20	0	114	65	135				
Toluene	22.69	0.14	20	0	113	65	135				
trans-1,2-Dichloroethene	20.04	0.14	20	0	100	65	135				
Trichloroethene	21.26	0.098	20	0	106	65	135				
Trichlorofluoromethane	21.60	0.14	20	0	108	65	135				
Vinyl Acetate	21.20	0.18	20	0	106	65	135				
Vinyl Chloride	20.96	0.097	20	0	105	65	135				
Surr: 4-Bromofluorobenzene	20.20	0	20	0	101	65	135				
Sample ID LCSD	SampType: LCSD	TestCod	e: TO-15	Units: ppbv		Prep Date	: 12/26/2	007	RunNo: 14	916	
Client ID: ZZZZZ	Batch ID: R14916	TestN	o: TO-15	• •		Analysis Date			SeqNo: 21		
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit I	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual

Value above quantitation range Qualifiers:

ND Not Detected at the Reporting Limit

Holding times for preparation or analysis exceeded

RPD outside accepted recovery limits

Spike Recovery outside accepted recovery limits

Page 4 of 6

Analyte detected below quantitation limits

CLIENT: Engeo **Work Order:** 0712091

ANALYTICAL QC SUMMARY REPORT BatchID: R14916

Project: 7584.1.001.01

Sample ID LCSD Sam	pType: LCSD	TestCod	de: TO-15	Units: ppbv		Prep Date	e: 12/26/2	2007	RunNo: 149	916	
Client ID: ZZZZZ Ba	tch ID: R14916	TestN	No: TO-15			Analysis Date	e: 12/26/2	2007	SeqNo: 214	4610	
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
1,1 - Dichloroethene	19.90	0.20	20	0	99.5	65	135	20.56	3.26	30	
1,1,1,2-Tetrachloroethane	23.06	0.10	20	0	115	65	135	23.18	0.519	30	
1,1,1-Trichloroethane	20.76	0.15	20	0	104	65	135	20.37	1.90	30	
1,1,2,2-Tetrachloroethane	22.36	0.15	20	0	112	65	135	23.48	4.89	30	
1,1,2-Trichloroethane	22.78	0.19	20	0	114	65	135	23.08	1.31	30	
1,1-Dichloroethane	19.60	0.17	20	0	98.0	65	135	20.54	4.68	30	
1,2,4-Trichlorobenzene	19.48	0.070	20	0	97.4	65	135	20.15	3.38	30	
1,2,4-Trimethylbenzene	21.93	0.18	20	0	110	65	135	22.24	1.40	30	
1,2-Dibromoethane(Ethylene dibromide	22.50	0.14	20	0	112	65	135	22.31	0.848	30	
1,2-Dichlorobenzene	20.72	0.10	20	0	104	65	135	21.85	5.31	30	
1,2-Dichloroethane	22.50	0.16	20	0	112	65	135	23.54	4.52	30	
1,2-Dichloropropane	23.93	0.22	20	0	120	65	135	24.14	0.874	30	
1,2-dichlorotetrafluoroethane(F114)	23.36	0.15	20	0	117	65	135	19.79	16.5	30	
1,3,5-Trimethylbenzene	22.69	0.14	20	0	113	65	135	23.15	2.01	30	
1,3-Butadiene	20.43	0.27	20	0	102	65	135	20.86	2.08	30	
1,3-Dichlorobenzene	20.80	0.060	20	0	104	65	135	21.84	4.88	30	
1,4-Dichlorobenzene	20.80	0.11	20	0	104	65	135	21.84	4.88	30	
1,4-Dioxane	22.93	0.14	20	0	115	65	135	23.83	3.85	30	
2-Butanone (MEK)	20.92	0.15	20	0	105	65	135	22.23	6.07	30	
2-Hexanone	23.18	0.21	20	0	116	65	135	24.32	4.80	30	
4-Ethyl Toluene	22.17	0.15	20	0	111	65	135	22.61	1.97	30	
4-Methyl-2-Pentanone (MIBK)	22.64	0.16	20	0	113	65	135	23.39	3.26	30	
Acetone	20.64	0.24	20	0	103	65	135	17.46	16.7	30	
Benzene	21.22	0.28	20	0	106	65	135	19.99	5.97	30	
Benzyl Chloride	21.71	0.12	20	0	109	65	135	22.24	2.41	30	
Bromodichloromethane	22.12	0.13	20	0	111	65	135	23.53	6.18	30	
Bromoform	24.49	0.17	20	0	122	65	135	24.06	1.77	30	
Bromomethane	20.44	0.20	20	0	102	65	135	20.49	0.244	30	
Carbon Disulfide	20.76	0.16	20	0	104	65	135	21.53	3.64	30	
Carbon Tetrachloride	20.32	0.15	20	0	102	65	135	20.09	1.14	30	
Chlorobenzene	21.94	0.092	20	0	110	65	135	23.12	5.24	30	

Qualifiers:

Value above quantitation range

ND Not Detected at the Reporting Limit

Holding times for preparation or analysis exceeded

RPD outside accepted recovery limits

Analyte detected below quantitation limits

Spike Recovery outside accepted recovery limits

Page 5 of 6

CLIENT: Engeo

Work Order: 0712091

Project: 7584.1.001.01

ANALYTICAL QC SUMMARY REPORT

BatchID: R14916

Sample ID LCSD	SampType: LCSD	TestCo	de: TO-15	Units: ppbv		Prep Dat	te: 12/26/2	007	RunNo: 149	916	
Client ID: ZZZZZ	Batch ID: R14916	Test	No: TO-15			Analysis Dat	te: 12/26/2	007	SeqNo: 214	1610	
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	RPD Ref Val	%RPD	RPDLimit	Qual
Chloroethane	20.49	0.15	20	0	102	65	135	20.33	0.784	30	
Chloroform	17.45	0.40	20	0	87.2	65	135	20.46	15.9	30	
Chloromethane	20.80	0.35	20	0	104	65	135	25.88	21.8	30	
cis-1,2-dichloroethene	19.53	0.14	20	0	97.6	65	135	20.11	2.93	30	
cis-1,3-Dichloropropene	22.60	0.080	20	0	113	65	135	22.81	0.925	30	
Dibromochloromethane	23.35	0.11	20	0	117	65	135	23.07	1.21	30	
Ethyl Acetate	20.81	0.12	20	0	104	65	135	22.04	5.74	30	
Ethyl Benzene	22.22	0.093	20	0	111	65	135	22.36	0.628	30	
Freon 113	20.68	0.12	20	0	103	65	135	21.48	3.80	30	
Hexachlorobutadiene	19.40	0.17	20	0	97.0	65	135	19.66	1.33	30	
Hexane	20.89	0.51	20	0	104	65	135	20.52	1.79	30	
Isopropanol	20.38	0.40	20	0	102	65	135	22.21	8.59	30	
m,p-Xylene	43.48	0.12	40	0	109	65	135	44.02	1.23	30	
Methylene Chloride	21.34	0.19	20	0	107	65	135	20.88	2.18	30	
MTBE	21.10	0.14	20	0	106	65	135	20.82	1.34	30	
Naphthalene	18.88	0.50	20	0	94.4	65	135	20.49	8.18	30	
o-xylene	21.75	0.14	20	0	109	65	135	21.47	1.30	30	
Styrene	21.97	0.15	20	0	110	65	135	22.29	1.45	30	
Tetrachloroethene	22.64	0.19	20	0	113	65	135	22.83	0.836	30	
Toluene	22.70	0.14	20	0	114	65	135	22.69	0.0441	30	
trans-1,2-Dichloroethene	19.45	0.14	20	0	97.3	65	135	20.04	2.99	30	
Trichloroethene	21.78	0.098	20	0	109	65	135	21.26	2.42	30	
Trichlorofluoromethane	20.92	0.14	20	0	105	65	135	21.6	3.20	30	
Vinyl Acetate	20.69	0.18	20	0	103	65	135	21.2	2.43	30	
Vinyl Chloride	22.19	0.097	20	0	111	65	135	20.96	5.70	30	
Surr: 4-Bromofluorobenzene	20.24	0	20	0	101	65	135	0	0	30	

Value above quantitation range Qualifiers:

ND Not Detected at the Reporting Limit

Holding times for preparation or analysis exceeded

RPD outside accepted recovery limits

Analyte detected below quantitation limits

Spike Recovery outside accepted recovery limits

Page 6 of 6

CHAIN OF CUSTODY RECORD

PROJECT NUMBER 7584.1.001.01		PROJECT NA 224 Ricken		cle						_			T										
PROJECT MANAGE	XATURE/PRINT)	ICH MU) (CA	W) a Ro									İ										
PROJECT MANAGE Kelly Krohn	ER:		- 0				TO-15																REMARKS REQUIRED DETECTION LIMITS
ROUTING: E-MAIL	kkrohn@engeo.c	om		Hard Copy	Keliy Krohn		-			,					ł		1						
SAMPLE NUMBER	DATE	TIME	MATRIX	NUMBER OF CONTAINERS	CONTAINER SIZE	PRESERVATIVE																	
56-11	17-Dec-07	11:22	Air	1	6 L Summa	None	×																0 A .
SG-12	17-Dec-07	13:05	Air	1		1	Х																0 2 A
SG-121Dupli	Celta 7-Dec-07	13:05	Air	1			х	L				\Box											03A
39-13	17-Dec-07	14:25	Air	. 1			х																04 A
59-14	17-Dec-07	11:12	Air	1			Х	L			,		•										05A
54-15	17-Dec-07	13:57	Air	1			х							\Box									061
36-10	17-Dec-07	12:36	Air	11			Х															_	072
5410	17-Dec-07	14:57	Air	1	: 45	4	х														<u> </u>	_	064
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44	2			12/12/1	17:40	2011	1		5									14	1/2	12	44	ф <u></u>	
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ENGEO SAN RAMON, CALIFOR (925) 866-9000 FAX (92																							
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10,2/17

Kelly Krohn

From: Project Management [pm@torrentlab.com]

Sent: Monday, July 21, 2008 12:50 PM

To: Kelly Krohn

Subject: RE: Initial and Final Pressures for 0712091 Canister

```
HI Kelly,
```

I am not the best person to answer the 1st question because I do not know the mathimatical conversion factors necessary to move from PSI to Hg inches but I have passed on the question to Mukesh. I know there is no direct conversion because the Hg inches are a much broader range than the definitive PSI units, but that's all I know!

As far as the dilution factor affecting results - the factor is applied to the raw value determined so the final result reported is the true value, not a manipulated value. The biggest effect is in the initial reporting limit which is raised by the dilution factor so that samples that ar ND have a raised ND value. This happens because all of the detection limits are based on an undiluted volume of spiked sample introduced onto the column. As with waters and soils, the DLs are "perfect world" limits so anything outside of that scenario is affected up or down: a decresed DL (concentration of a sample) or increased DL (dilution or prep factor).

Give me a call if I need to clear up some mud in that response!

Patti

```
On Mon, July 21, 2008 11:10 am, Kelly Krohn wrote:
> Hi Patti,
> We have a few follow up questions regarding the information you have
> provided below.
> 1. Laboratory Pressure Readings - please verify the reported units on
> the previously provided pdf. As stated below, the units are psi;
> however, we cannot resolve the conversion to mm Hg. 2. A dilution
> factor has been applied to all of the samples. Has this dilution
> factor in any way biased the final results?
> We very much appreciate your help.
> Thank you,
> Kelly
> Kelly Krohn, PE
> Project Engineer
>
>
>
> ----Original Message----
> From: Project Management [mailto:pm@torrentlab.com]
> Sent: Monday, June 02, 2008 2:05 PM
> To: Kelly Krohn
> Subject: RE: Initial and Final Pressures for 0712091 Canister
>
```

```
> No, those is actually PSI values. In order to introduce 500 mL of
> sample onto the instrument, the HG inches must be as close to -5" as
> possible. FOr low level analysis, we need to introduce a full Liter
> onto the instrument and then the Hg inches must be between "0" and
> "+5" inches. As a result, Anytime we receive the canister with less
> that -5" of Mercury, it ia Torrent's practice to add Argon to the
> canister to raise the pressure to 0" (plus or minues 5"). When two or
> more analyses are requested (TO-15, TO-3, ASTM1946D, etc), we raise
> the Hg inches to slightly above 0 in order to prevent the necessity of
> adding additional Argon later for the multiple tests - or, if ESLs are
> required for most compounds but some are way over calibration range,
> multiple injections of different volume may be necessay.
> I looked at the tags attached to the canisters that were submitted by
> En GEO and almost all of them had -7" or less. I could copy those
> tags and send them to you if you would like. Just let me kow.
> On Mon, June 2, 2008 1:38 pm, Kelly Krohn wrote:
>> This looks great. I assume that the initial pressure would be that
>> at the time of receipt. May I ask why it is needed to increase the
> pressure?
>> Are the units mm Hq?
>>
>>
>>
>> Thanks,
>> Kelly
>>
>>
>>
>>
>> Kelly Krohn, PE
>>
>>
>>
>> Project Engineer
>>
>>
>>
>>
>>
>> ----Original Message----
>> From: Project Management [mailto:pm@torrentlab.com]
>> Sent: Monday, June 02, 2008 1:11 PM
>> To: Kelly Krohn
>> Subject: Initial and Final Pressures for 0712091 Canister
>>
>>
>>
>> Hi Kelly,
>>
>>
>> Attached is a copy from our intital/final pressure logbook that is
>>
> kept
>> at the instrument. Is this what you were looking for? Let me know!
>> Patti
>>
>>
```

> Hi Kelly,

```
>>
>>
>> Project Management Team
>> Torrent Laboratory, Inc.
>> 483 Sinclair Frontage Rd
>> Milpitas, CA 95035
>> PH:(408)263-5258; Nutan x209, Patti x208, Stacy 707-206-0216
>> Fax: (408) 263-8293
>> Email: pm@torrentlab.com
>> www.torrentlab.com
>>
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>> Visit us at our newly re-developed website!! www.torrentlab.com
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>> If you are not the intended recipient of this message and its
>>
> contents,
>> please contact us immediately at (408) 263-5258 and delete the
>> message
> and
>> its contents.
>>
>>
> Warm regards,
> Project Management Team
> Torrent Laboratory, Inc.
> 483 Sinclair Frontage Rd
> Milpitas, CA 95035
> PH: (408)263-5258; Nutan x209, Patti x208, Stacy 707-206-0216
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APPENDIX D

Monitoring Well Sampling Data Sheets

ENGEO INCORPORATED GROUNDWATER MONITORING WELL SAMPLING DATA

Job Name: 224 Rickenbacker Circle

Location: 224 Rickenbacker Circle, Livermore, CA

Date: January 28, 2008

Client: Mr. Robert Strong

By: Richard Gandolfo

WELL INFORMATION

Well Number: MW-1

Diameter (in): 2

Total Depth (feet below top of casing): 34.5

Screen Interval (feet bgs): 10 to 35

Depth to Water (ft): 25.1

Casing Volume (gal): 1.59

PURGING INFORMATION

Bailer: N/A Pump: 12-V Rate (gpm): 0.15 Time (initial/finish): 9:46/10:00

Volume Removed (gal): 5 No. of Casing Volumes: 3

Time	Volume (gal.)	DO (mg/l)	Temp. (°C)	Conductivity (mS/cm)	Salinity (%)	pН
9:46	1/2	7.49	16.5	0.910	0.04	7.70
9:49	1	7.52	17.4	0.940	0.04	7.54
9:50	1.5	7.70	17.7	0.813	0.03	7.49
9:53	2	7.71	17.7	0.807	0.03	7.50
9:55	4	7.73	17.9	0.809	0.03	7.45
9:57	4.5	7.84	18.2	0.806	0.03	7.40
10:00	5.0	7.97	18.2	0.807	0.03	7.40

SAMPLE INFORMATION AND LABORATORY TESTING PARAMETERS

Bailer: N/A Pump: 12-V Rate (gpm): 0.15 Decon: TSP/Alcononx

Sample ID	Sample Date/Time	Field Blank Sample ID (if applicable)	Duplicate Sample ID (if applicable)
MW-1	1/28/2008 10:00	N/A	N/A

Quantity of Containers	Size and Preservative	Test Method	Analytes
3	40-mL VOA, HCl preserved	EPA 8260B	Volatile Organic Compounds

ENGEO INCORPORATED GROUNDWATER MONITORING WELL SAMPLING DATA

Job Name: 224 Rickenbacker CircleJob Number: 7584.100.101Location: 224 Rickenbacker Circle, Livermore, CADate: January 28, 2008Client: Mr. Robert StrongBy: Richard Gandolfo

WELL INFORMATION

Well Number: MW-2

Total Depth (feet below top of casing): 34.3

Depth to Water (ft): 25.23

Diameter (in): 2

Screen Interval (feet bgs): 10 to 35

Casing Volume (gal): 1.56

PURGING INFORMATION

Bailer: N/A Pump: 12-V Rate (gpm): 0.15 Time (initial/finish): 11:26/11:45

Volume Removed (gal): 5.5 No. of Casing Volumes: 3.5

Time	Volume (gal.)	DO (mg/l)	Temp. (°C)	Conductivity (mS/cm)	Salinity (%)	рН
11:26	1/2	8.11	15.9	0.774	0.04	7.58
11:29	1	8.08	17.2	0.801	0.03	7.48
11:31	2	7.84	18.1	0.807	0.03	7.39
11:33	3	7.90	18.1	0.810	0.03	7.41
11:34	3.5	8.10	18.2	0.805	0.03	7.41
11:36	4	8.12	18.3	0.807	0.03	7.40
11:37	4.5	8.13	18.2	0.803	0.03	7.41
11:40	5	8.05	17.8	0.795	0.03	7.39
11:45	5.5	7.85	18.4	0.800	0.03	7.39

SAMPLE INFORMATION AND LABORATORY TESTING PARAMETERS

Bailer: N/A Pump: 12-V Rate (gpm): 0.15 Decon: TSP/Alcononx

Sample ID	Sample Date/Time	Field Blank Sample ID (if applicable)	Duplicate Sample ID (if applicable)
MW-2	1/28/2008 11:45	N/A	N/A

Quantity of Containers	Size and Preservative	Test Method	Analytes
3	40-mL VOA, HCl preserved	EPA 8260B	Volatile Organic Compounds

ENGEO INCORPORATED GROUNDWATER MONITORING WELL SAMPLING DATA

Job Name: 224 Rickenbacker CircleJob Number: 7584.100.101Location: 224 Rickenbacker Circle, Livermore, CADate: January 28, 2008Client: Mr. Robert StrongBy: Richard Gandolfo

WELL INFORMATION

Well Number: MW-3

Diameter (in): 2

Total Depth (feet below top of casing): 34.6

Screen Interval (feet bgs): 10 to 35

Depth to Water (ft): 25.25

Casing Volume (gal): 1.58

PURGING INFORMATION

Bailer: N/A Pump: 12-V Rate (gpm): 0.15 Time (initial/finish): 8:30 / 8:57

Volume Removed (gal): 5 No. of Casing Volumes: 3

Time	Volume (gal.)	DO (mg/l)	Temp. (°C)	Conductivity (mS/cm)	Salinity (%)	рН
8:30	1/2	7.54	15.2		0.03	7.62
8:40	1.5	8.02	17.2	0.779	0.03	7.45
8:42	2	6.88	17.7	0.788	0.03	7.50
8:44	2.5	7.44	17.7	0.790	0.03	7.49
8:48	3	7.37	17.9	0.784	0.03	7.46
8:50	3.5	7.66	17.4	0.785	0.03	7.48
8:52	4	7.68	18.1	0.788	0.03	7.40
8:54	4.5	7.77	18.0	0.787	0.03	7.43
8:57	5	7.93	18.0	0.785	0.03	7.40

SAMPLE INFORMATION AND LABORATORY TESTING PARAMETERS

Bailer: N/A Pump: 12-V Rate (gpm): 0.15 Decon: TSP/Alcononx

Sample ID	Sample Date/Time	Field Blank Sample ID (if applicable)	Duplicate Sample ID (if applicable)
MW-3	1/28/2008 8:57	N/A	N/A

Quantity of Containers	Size and Preservative	Test Method	Analytes
3	40-mL VOA, HCl preserved	EPA 8260B	Volatile Organic Compounds