Memorandum

Date:	March 19, 2009
То:	Ms. Kristi Doll, AFCEE/EXEC
From:	Mr. Walt Gee, Environmental Scientist
Subject:	Long-Term Monitoring Activities and Soil Gas Investigation at Air Force Plant 59, Johnson City, New York

Distribution: Mr. George Walters, ASC; Mr. Brian Jankauskas, NYSDEC

In 2008, Earth Tech AECOM completed the Long-Term Monitoring (LTM) activities, United States Geological Survey (USGS) monitoring well sampling, and soil gas investigation at Air Force Plant 59 (AFP 59) in Johnson City, New York. The Air Force Center for Engineering and the Environment (AFCEE) contracted Earth Tech AECOM to complete the LTM activities, USGS monitoring well sampling, and soil gas investigation based on the findings from the *Final Soil-Gas and Groundwater Monitoring Report* (Earth Tech, 2007).

The objectives of this memorandum are to summarize:

- The purpose, collection procedures, and results of the LTM activities;
- The purpose, collection procedures, and results of the USGS monitoring well sampling activities;
- The purpose, collection procedures, and results of the soil gas investigation; and
- The conclusions and recommendations based on the results of the LTM activities, USGS monitoring well sampling activities, and soil gas investigation.

Long-Term Monitoring Activities

Purpose of the Long-Term Monitoring Activities

Based on the conclusions presented in the *Final Remedial Investigation Report* (Earth Tech, 1996) and recommendations made by the New York State Department of Environmental Conservation (NYSDEC), it was determined that volatile organic compounds (VOCs) represent the only chemicals of potential concern (COPCs) in groundwater at AFP 59. As a result, the *Record of Decision* (Earth Tech, 1999b) for AFP 59 describes the remedial alternative (i.e., the upgrade of the Camden Street Well Field groundwater treatment system) chosen as most appropriate for treating the VOCs in groundwater at AFP 59. As part of the requirements defined in the *Record of Decision* (Earth Tech, 1999b), a LTM program was established for AFP 59. The LTM program, which is defined in the April 27, 1999 letter to the NYSDEC (Earth Tech, 1999a), was concluded with the November 2004 sampling event. The LTM included sampling the following monitoring wells: SW1, DW1, SW3, DW3, SW4, and SW7. Monitoring wells SW1 and DW1 represent upgradient (background) wells, and monitoring wells SW3 and DW3 represent downgradient wells. Monitoring wells SW4 and SW7 have historically had the highest concentrations of VOCs.

A soil pile containing trichloroethene (TCE) contamination in the East Basement of the AFP 59 facility was excavated and removed in July 2005. The soil pile was upgradient of monitoring wells SW3, DW3, SW4, and SW7. This sampling event was designed to observe what effect this removal action might have on groundwater concentrations of VOCs.

Procedures Used for the Long-Term Monitoring Activities

Sampling activities followed protocols presented in the Final Work Plan for Groundwater Monitoring at AFP 59 (Earth Tech, 1998) and the Final Work Plan Addendum (Earth Tech, 2008).

Earth Tech collected two rounds of groundwater samples during the LTM activities. Groundwater samples were collected in June 2008 at six on-site monitoring wells (SW1, DW1, SW3, DW3, SW4, and SW7) and two off-site monitoring wells (URS_2S and URS_2D), and analyzed for VOCS by USEPA Method SW8260B. Five of the on-site monitoring wells (SW1, SW3, DW3, SW4, and SW7) were also analyzed for 1,4-dioxane using United States Environmental Protection Agency (USEPA) Method SW8260SIM. In November 2008, Earth Tech collected groundwater samples from four monitoring wells (SW3, DW3, SW4, and SW7) and analyzed the samples for VOCs by USEPA Method 8260B.

Analytical Results from the Long-Term Monitoring Activities

The following paragraphs discuss the VOCs that were detected in the groundwater samples, including those samples collected from both on-site and background monitoring wells and off-site monitoring wells. The analytical results for groundwater samples collected from monitoring wells installed in the shallow and deep zones of the aquifer are discussed below. The VOCs detected in groundwater samples are illustrated on Figure 1. The analytical results for all groundwater samples collected during the June and November 2008 sampling events are summarized in Table 1. Note: Sample 59SW7WG1 served as the matrix spike/matrix spike duplicate (MS/MSD). Recoveries of cis-1,2-dichloroethene (cis-1,2-DCE) and vinyl chloride exceeded the upper control limits of MS/MSD samples. As a result, cis-1,2-DCE and vinyl chloride for sample 59SW7WG1 were qualified "M" (Matrix effect: the concentration is estimated due to a matrix effect).

Shallow Zone of the Aquifer. VOCs were detected in the groundwater samples collected from on-site monitoring wells SW3, SW4, and SW7, and the off-site monitoring well URS_2S (Refer to Figure 1). Chlorinated hydrocarbons were the only detected VOCs in the samples collected from the shallow zone of the aquifer. VOCs were not detected in the groundwater sample collected from monitoring well SW1.

The following maximum concentrations were detected in the regular and duplicate groundwater samples collected from monitoring well SW4 during both the June and November 2008 events: 1,1,1-trichloroethane (1,1,1-TCA) at 2.98 micrograms per liter (μ g/L); 1,1-dichloroethane (1,1-DCA) at 1.51 μ g/L; 1,1-DCE at 0.751 J μ g/L; cis-1,2-DCE at 4.35 μ g/L; tetrachloroethene (PCE) at 0.965 J μ g/L; trans-1,2-DCE at 0.364 J μ g/L; and TCE at 17.8 μ g/L. The following maximum concentrations were detected in the groundwater sample collected from monitoring well SW7 during both the June and November 2008 events: 1,1,1-TCA at 2.5 μ g/L; 1,1-DCA at 5.04 μ g/L; PCE at 0.843 J μ g/L; cis-1,2-DCE at 35.3 M μ g/L; trans-1,2-DCE at 0.302 J μ g/L; vinyl chloride at 1.21 M μ g/L; and TCE at 8.15 μ g/L.

During the June 2008 sampling effort, 1,4-dioxane was sampled in four on-site shallow monitoring wells. 1,4-Dioxane was detected in monitoring wells SW4, SW4 duplicate sample, and SW7 at concentrations of 8.18 μ g/L; 7.2 μ g/L; and 4.66 μ g/L, respectively. 1,4-Dioxane was not detected in monitoring wells SW1 and SW3.

The following maximum concentrations were detected in the regular and duplicate groundwater samples collected from off-site monitoring well URS_2S: 1,1,1-TCA at 2.25 μ g/L; 1,1-DCA at 0.585 J μ g/L; cis-1,2-DCE at 0.996 J μ g/L; and TCE at 2.22 μ g/L.

Deep Zone of the Aquifer. VOCs were detected in the groundwater samples collected from on-site monitoring well DW3 and the off-site monitoring well URS_2D (Refer to Figure 1). Chlorinated hydrocarbons were the only VOCs detected in the samples collected from the deep zone of the aquifer. VOCs were not detected in the groundwater sample collected from monitoring well DW1.

The only VOC detected in monitoring well DW3 was cis-1,2-DCE at 73.1 μ g/L. The following maximum concentrations were detected in the groundwater sample collected from the off-site monitoring well URS_2D: 1,1-DCA at 0.239 J μ g/L; cis-1,2-DCE at 71.9 μ g/L; and vinyl chloride at 0.354 J μ g/L. Cis-1,2-DCE exceeded the New York State Drinking Water Standard of 5 μ g/L.

During the June 2008 sampling event, 1,4-dioxane was sampled in two on-site deep monitoring wells (DW1 and DW3). 1,4-Dioxane was only detected in monitoring well DW3 (14.3 μ g/L).

Trend Analysis

Table 2 presents concentrations of the most commonly detected chlorinated hydrocarbons in groundwater at AFP 59 over time. Only monitoring wells that were sampled as part of the groundwater monitoring program are included in the table.

In the groundwater samples collected from the shallow monitoring wells during the November 2008 sampling event, concentrations of the chlorinated hydrocarbons in monitoring well SW3 remained relatively constant (TCE and 1,1,1-TCA) or decreased to non-detect (ND) concentrations (cis-1,2-DCE) when compared to the previous sampling event in October 2005. Concentrations of TCA, TCE, cis-1,2-DCE; and 1,1-DCA increased from the October 2005 sampling event.

The concentrations of the chlorinated hydrocarbons in monitoring well SW4 remained relatively constant, with only moderate variation in TCE concentrations when compared to the October 2005 sampling event. TCE concentrations decreased from 43 μ g/L in October 2005 to 12.7 μ g/L in November 2008. The concentrations of TCA (2.2 μ g/L to 0.513 J μ g/L); 1,1-DCE (1 μ g/L to ND); cis-1,2-DCE (6.3 μ g/L to 3.38 μ g/L); and 1,1-DCA (1.7 μ g/L to 0.825 J μ g/L) each decreased during the November 2008 sampling event. The concentration of trans-1,1-DCE (ND to 0.364 J μ g/L) slightly increased compared to the November 2008 sampling event.

Concentrations of TCA in monitoring well SW7 increased from 0.73 J μ g/L in October 2005 to 2.5 μ g/L in June 2008 and then decreased to 1.88 μ g/L in November 2008. TCE decreased slightly from 3.1 μ g/L in October 2005 to 2.94 μ g/L in June 2008 and more than doubled to 8.15 μ g/L in November 2008. Cis-1,2-DCE decreased from 12 μ g/L in October 2005 to 6.34 μ g/L in June 2008 before increasing to 35.3 M μ g/L in November 2008. Trans-1,2-DCE increased from ND in October 2005 and June 2008 to 0.302 J μ g/L in November 2008. 1,1-DCA increased slightly from 1.4 μ g/L to 1.59 μ g/L in June 2008 and more than tripled to 5.04 μ g/L in November 2008.

In the groundwater sample collected from deep monitoring well DW3 during the June 2008 sampling event, the concentrations of chlorinated hydrocarbons remained at ND with the exception of cis-1,2-DCE, which increased from 3 μ g/L in October 2005 to 73.1 μ g/L in June 2008. Cis-1,2-DCE decreased slightly to 67.3 μ g/L in November 2008. Additionally, 1,1-DCA and trans-1,2-DCE were detected at concentrations of 0.41 J μ g/L and 1.18 μ g/L, respectively, during the November 2008 sampling event. VOCs were not detected in the groundwater sample collected from deep monitoring well DW1 and shallow monitoring well SW1. These results are consistent with previous sampling events.

USGS Monitoring Well Sampling

Purpose of the United States Geological Survey Monitoring Well Sampling

The USGS, in cooperation with the United States Air Force (USAF), installed eight monitoring wells in April 1995 in conjunction with an investigation of the hydrogeology and water quality of the Camden Street-Ballpark aquifer near Johnson City, New York. As part of the site closure activities for AFP 59,

the USAF needed to evaluate the purpose for the eight USGS monitoring wells in the groundwater sampling program for the site (Refer to Figure 2 for monitoring well locations). Earth Tech AECOM was tasked with locating the eight monitoring wells, assess the integrity of the monitoring wells, collect groundwater samples from each monitoring well, and make a recommendation on monitoring well abandonment.

Procedures Used to Collect the United States Geological Survey Monitoring Well Samples

The USGS monitoring well sampling was completed using the procedures found in the *Final Work Plan* Addendum for the LTM and Soil Gas Investigation at AFP 59 (Earth Tech, 2008).

Analytical Results from the United States Geological Survey Monitoring Well Samples

Of the eight USGS monitoring wells, only six monitoring wells were located during this investigation. Two monitoring wells (GS_9501S and GS_9501D) were not located and are believed to be buried. Monitoring well GS_9503 was located, but the surface completion was missing including the well cap. This monitoring well was not sampled due to the infiltration of surface water. The monitoring well will be difficult to access with a drill rig when the decision to abandon the monitoring well is made. Monitoring well GS_9504 was also located, but the surface completion had been destroyed with surface water and fill material entering the well casing.

Four monitoring wells (GS_9502S, GS_9502D, GS_9505, and GS_9506) were sampled in June 2008 for VOCs using USEPA Method 8260B. VOCs were not detected in monitoring well GS_9502D. 1,1,1-TCA was detected in monitoring wells GS_9502S and GS_9505 at concentrations of 9.04 μ g/L and 4.49 μ g/L, respectively. The 1,1,1-TCA detection in monitoring well GS_9502S exceeds the New York State Drinking Water Standard of 5 μ g/L. 1,1-DCA was detected in monitoring well GS_9505 at a concentration of 0.389 J μ g/L. The 1,1-DCA detection was below the New York State Drinking Water Standard of 5 μ g/L. PCE was detected at a concentration of 27.2 μ g/L in monitoring well GS_9506, which exceeds the New York State Drinking Water Standard of 5 μ g/L.

Soil Gas Investigation

Purpose of the Soil Gas Investigation

Two soil gas samples were collected in November 2004 to evaluate the potential off-site migration of soil gas downgradient of the chlorinated hydrocarbon plume. Elevated concentrations of chlorinated hydrocarbons were detected. Additional soil gas samples were collected in October 2006, and elevated concentrations of chlorinated hydrocarbons were once again detected. Based on the results from these previous investigations and the data gaps that existed after the previous investigations, the NYSDEC requested that additional soil-gas sampling be initiated on the western side of AFP 59.

Procedures Used to Collect the Soil Gas Samples

The soil gas sampling protocol was in accordance with the NY State Department of Health (NYSDOH) *Guidance for Evaluating Soil Vapor Intrusion in the State of New York* (NYSDOH, 2006) and included the following:

- A hole was drilled to a depth corresponding to the depth of residential basements in the area (approximately 3 feet below ground surface [bgs] in the parking lot, approximately 9 feet bgs on the embankment along the property boundary, and up to 9 feet bgs at the off-site monitoring well location).
- A piece of laboratory or food grade Teflon-lined tubing was inserted through the sampling rods. Modeling clay was used to seal the tubing at the surface and ensure ambient air was not entering the sample container. An adapter was utilized to connect the Teflon-lined tubing to the vapor probe to prevent ambient air within the drill rods from impacting the sample interval.

- To further ensure ambient air was not entering the sample container, a domed enclosure filled with a tracer gas covered the sample points. When conducting the tracer gas test, the dome covered the borehole, drill rods, and as many connections as possible to confirm that ambient air is not impacting the sample.
- A vacuum pump was used to purge between one and three volumes of the tubing prior to sample collection. Purge rates were less than 200 milliliters per minute (mL/min). The purged air was containerized, screened with a photoionization detector, and released.
- After purging was complete, samples were collected in a Summa® canister. These samples were collected at a flow rate of less than 200 mL/min.
- One duplicate sample was collected from a soil gas sampling point (SG-38). The duplicate sample was taken from the same sample point using a new Summa® canister.

Analytical Results from the Soil Gas Samples

During the June 2008 soil gas investigation, three sample points were advanced on the western plant boundary to evaluate soil gas migration off-site. The samples were advanced to a depth of approximately 9 feet bgs, the approximate depth of the residential basements on the adjacent property. Additionally, three soil gas locations (two samples near the reservoir on the western part of the plant boundary and one sample on the eastern plant boundary) were advanced to 3 feet bgs near sampling locations where soil gas sampling was attempted, but samples were unable to be collected. Finally, one soil gas location was advanced to 9 feet bgs, adjacent to an off-site monitoring well where a groundwater sample was collected to attempt to relate groundwater and soil gas concentrations. The analytical results, as well as the locations of the soil gas samples, are illustrated on Figure 2. Table 3 summarizes the analytical data from the soil gas sampling event.

Chlorinated hydrocarbons were detected in the three samples collected on the western plant boundary (SG-33, SG-34, and SG-35). 1,1,1-TCA concentrations ranged from 2.1 micrograms per cubic meter (μ g/m³) in soil gas sample SG-34 to 26 μ g/m³ in soil gas sample SG-33. Cis-1,2-DCE was only detected in soil gas sample SG-34 at a concentration of 0.44 J μ g/m³. PCE was detected in all three samples at a maximum concentration of 20 μ g/m³ (soil gas sample SG-34). TCE was detected in two of the soil gas samples at a maximum concentration of 29 μ g/m³ (soil gas sample SG-33).

Elevated VOCs were detected in the two soil gas samples collected near the fire suppression reservoir on the western side of AFP 59 (SG-36 and SG-37). The following concentrations were detected in soil gas sample SG-36: 1,1,1-TCA at 9,000 μ g/m³; 1,1-DCA at 120 μ g/m³; acetone at 1,800 μ g/m³; benzene at 35 μ g/m³; cis-1,2-DCE at 0.97 μ g/m³; PCE at 13 μ g/m³; TCE at 1,900 μ g/m³; and vinyl chloride at 6.8 μ g/m³.

The following concentrations were detected in soil gas sample SG-37: 1,1,1-TCA at 110,000 μ g/m³; 1,1,2-TCA at 20 μ g/m³; 1,1-DCA at 4,900 μ g/m³; 1,1-DCE at 200 μ g/m³; 1,2-DCA at 9,600 μ g/m³; acetone at 1,000 J μ g/m³; benzene at 16 μ g/m³; carbon tetrachloride at 8.8 μ g/m³; cis-1,2-DCE at 3,100 μ g/m³; PCE at 38 μ g/m³; trans-1,2-DCE at 80 μ g/m³; TCE at 42,000 μ g/m³; and vinyl chloride at 5.1 μ g/m³.

One soil gas sample (and one duplicate sample) was collected adjacent to shallow monitoring well URS_2S to compare soil gas and groundwater concentrations (SG-38). The following maximum concentrations were detected in the regular and duplicate soil gas samples: 1,1,1-TCA at 11 J μ g/m³; benzene at 9.4 J μ g/m³; cis-1,2-DCE at 1.0 J μ g/m³; PCE at 14 μ g/m³; and TCE at 11 μ g/m³. In monitoring well URS_2S, the above VOC concentrations were as follows: 1,1,1-TCA at 2.25 μ g/L, cis-1,2-DCE at 0.996 J μ g/L, and TCE at 2.22 μ g/L. PCE was not detected in the monitoring well URS_2S.

One soil gas sample was collected east of the TCE-contaminated soil pile removed in July 2005 (SG-39). The following concentrations were detected in soil gas sample SG-39: 1,1,1-TCA at 1.9 μ g/m³; acetone at 2,400 μ g/m³; benzene at 29 μ g/m³; cis-1,2-DCE at 0.69 μ g/m³; and PCE at 13 μ g/m³. TCE was not detected in the soil gas sample at this location.

Additional Fire Suppression Reservoir Investigation

Based on the results of the June 2008 soil gas investigation, an additional investigation was conducted to determine the nature and extent of the elevated soil gas concentrations at soil gas samples SG-36 and SG-37. Efforts to collect soil gas samples surrounding soil gas sample SG-36 were unsuccessful. After numerous attempts to advance a hand auger, one soil sample was collected at SG-36 at 3 feet bgs. The following VOCs were detected in soil gas sample SG-36: methylene chloride at 3.87 micrograms per kilogram (μ g/kg); 1,1,1-TCA at 2.53 J μ g/kg; and TCE at 0.637 J μ g/kg. Table 4 summarizes the results of the soil sampling event.

Conclusions and Recommendations

Long-Term Monitoring Activities

Although VOC concentrations in the shallow monitoring wells have generally decreased since October 2005, concentrations of TCE and 1,1-DCA exceed the New York State Drinking Water Standard of 5 μ g/L. Additionally, the concentration of cis-1,2-DCE exceeded the New York State Drinking Water Standard of 5 μ g/L during the June 2008 sampling event in monitoring well SW7. Groundwater concentrations detected in off-site shallow monitoring well URS_2S did not exceed the New York State Drinking Water Drinking Water Standard of 5 μ g/L for chlorinated compounds.

In the deep monitoring wells, cis-1,2-DCE was the only contaminant that had concentrations exceeding the New York State Drinking Water Standard of 5 μ g/L. Monitoring well DW3, located on the AFP 59 boundary downgradient of the suspected source, and monitoring well URS_2D, located at a downgradient, off-site location, exceeded the New York State Drinking Water Standard for cis-1,2-DCE in both groundwater sampling events.

Based on the results of the LTM activities, groundwater exceeding the New York State Drinking Water Standards is migrating off of AFP 59 property in the deep monitoring wells. Additional groundwater monitoring is recommended to monitor the migration of contaminants off-site.

United States Geological Survey Monitoring Well Sampling

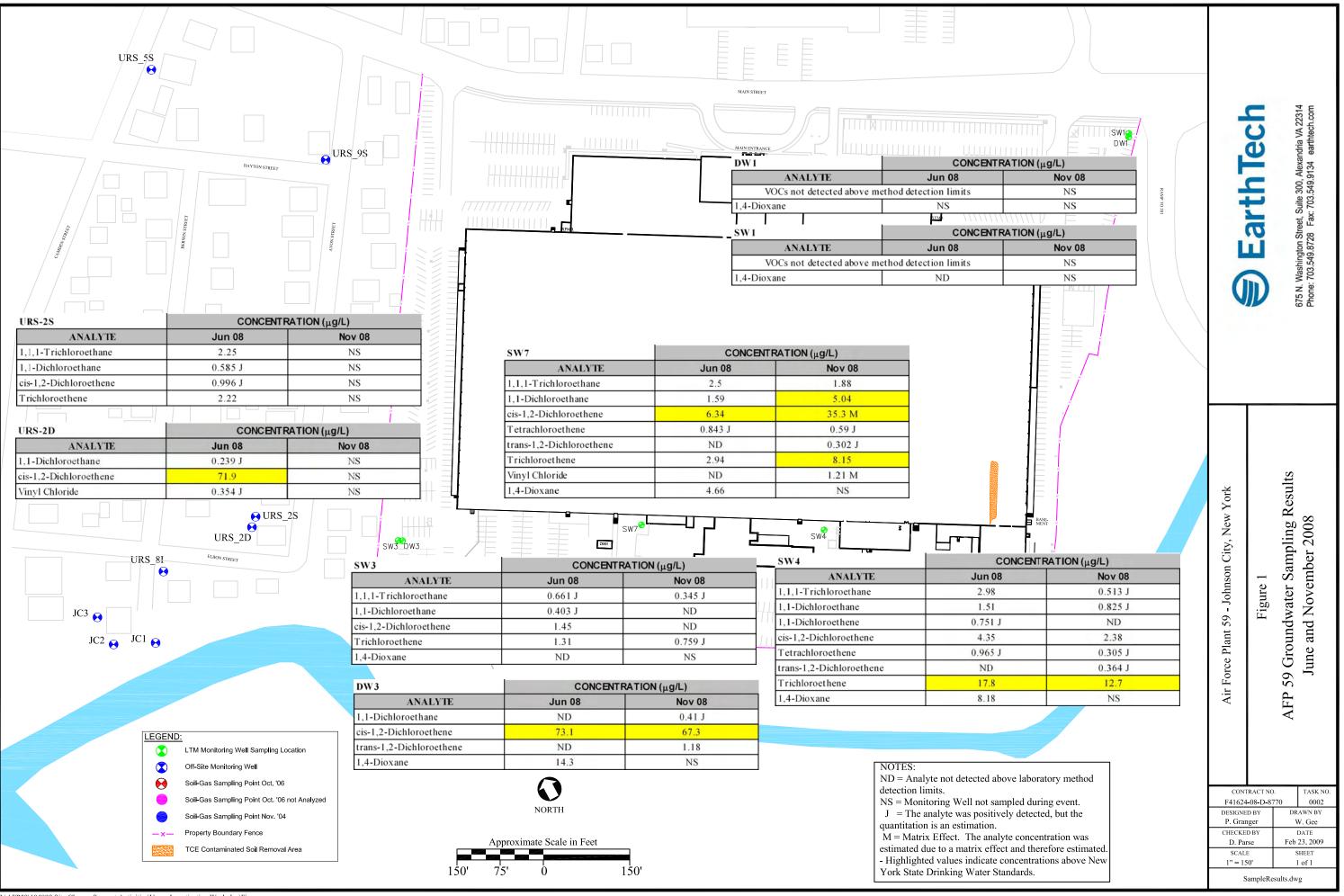
Concentrations of VOCs detected in the USGS monitoring wells were well below the New York State Drinking Water Standards, with the exception of PCE at monitoring well GS_9506 and 1,1,1-TCA at monitoring well GS_9502D. Monitoring wells GS_9502S and GS_9506 are located upgradient and do not contribute to the groundwater plume migrating on-site. Two of the eight monitoring wells could not be sampled due to surface water infiltration as a result of damage to the wellhead. Also, two of the remaining six monitoring wells could not be located. Therefore, it is recommended to abandon the USGS monitoring wells since the monitoring wells are no longer needed as part of the AFP 59 groundwater monitoring program. The USGS monitoring wells should be abandoned in accordance with New York State guidelines.

Soil Gas Investigation

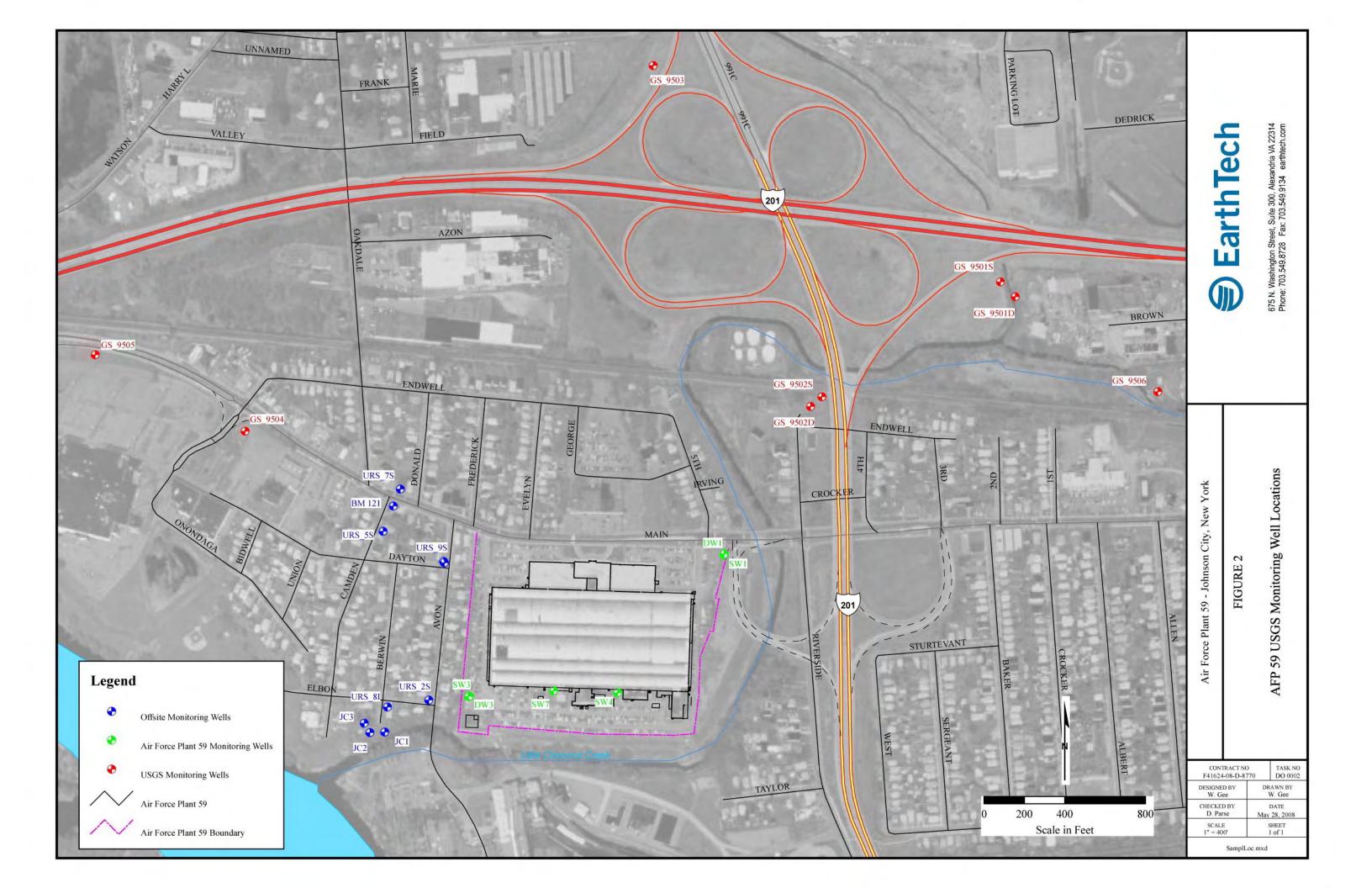
Contaminants were detected in the soil gas samples collected during the soil gas investigation. At the western AFP 59 boundary, VOCs appear to be migrating off-site into the neighborhood. VOCs were detected at low concentrations adjacent to monitoring well URS_2S. Additionally, high VOC concentrations were detected around the fire suppression reservoir. Based on the results of the soil gas investigation, additional sampling should be conducted around the fire suppression reservoir and in the neighborhood.

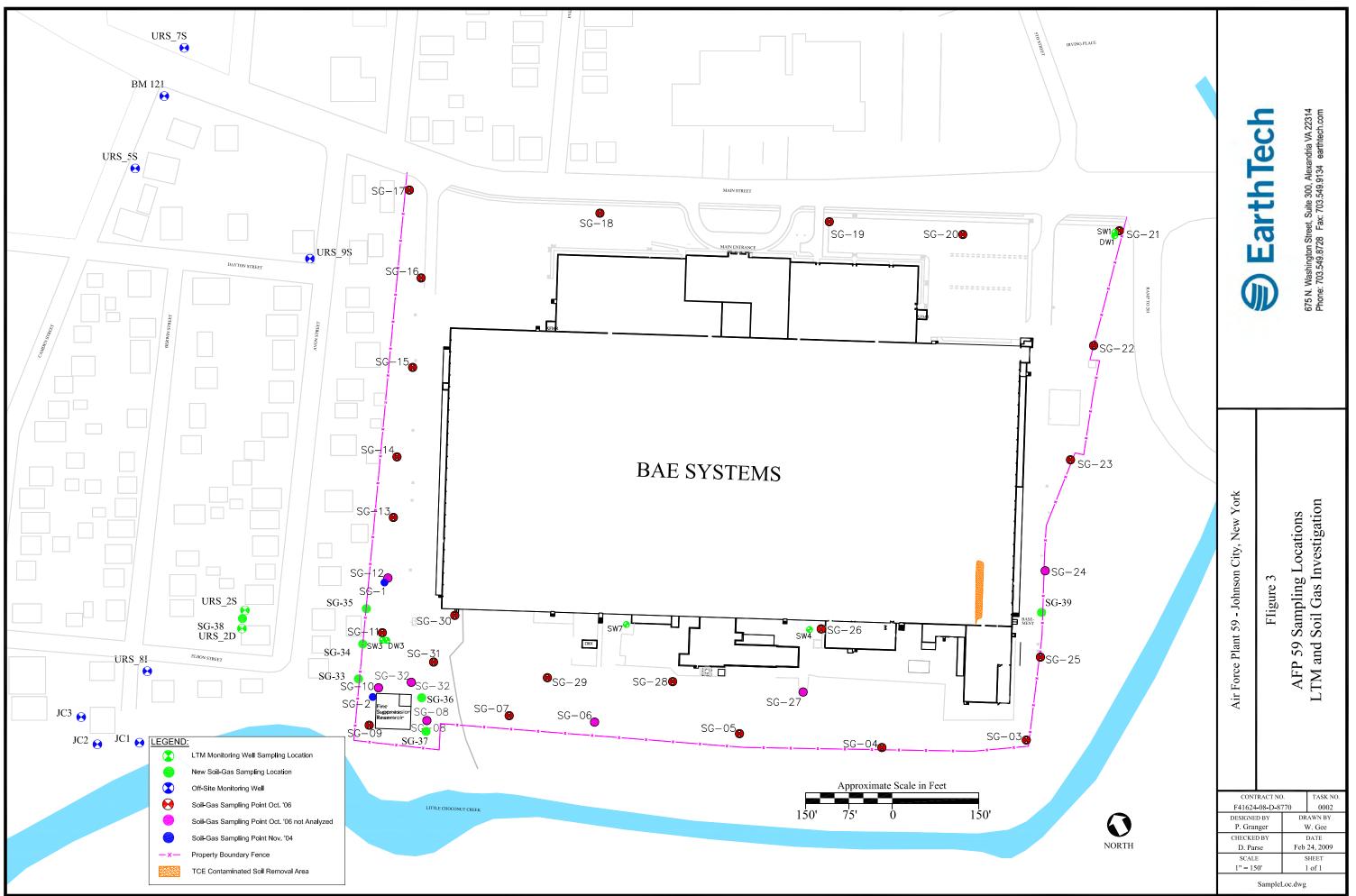
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Figures



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Tables

	Sum	nary of Detecte	Table 1 d VOCs in Mon June 2008	itoring Well Sa	amples		
Location ID Date Sampled	DW1 6/14/2008	DW3 6/15/2008	SW1 6/14/2008	SW3 6/15/2008	SW4 6/16/2008	SW4 (DUP) 6/16/2008	SW7 6/15/2008
Analyte		Volatiles by I	EPA SW-846 Meth	od 8260 (ug/L)			
1,1,1-Trichloroethane	1 U	1 U	1 U	0.661 J	2.98	2.87	2.5
1,1-Dichloroethane	1 U	1 U	1 U	0.403 J	1.51	1.48	1.59
1,1-Dichloroethene	1 U	1 U	1 U	1 U	0.751 J	0.712 J	1 U
1,4-Dioxane	100 U	100 R	100 R	100 R	100 R	100 R	100 R
Chloroform	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
Chloromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U
cis-1,2-Dichloroethene	1 U	73.1	1 U	1.45	4.35	4.13	6.34
Tetrachloroethene	1 U	1 U	1 U	1 U	0.959 J	0.965 J	0.843 J
trans-1,2-Dichloroethene	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichloroethene	1 U	1 U	1 U	1.31	17	17.8	2.94
Trichlorofluoromethane	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Vinyl chloride	1 UJ	1 U	1 U	1 U	1 U	1 U	1 U
Analyte		Volatiles by EP.	A SW-846 Method	8260 SIM (ug/L)			
,4-Dioxane	-	14.3	2 U	2 U	8.18	7.2	4.66

Table 1	
Summary of Detected VOCs in Monitoring Well Samples (continued)	

June 2008

Location ID Date Sampled	GS-9502S 6/15/2008	GS-9505 6/15/2008	GS-9506 6/16/2008
Analyte	Volatiles by EPA S	SW-846 Method 8260 (ug/L)	
1,1,1-Trichloroethane	9.04	4.49	IU
I, I-Dichloroethane	10	0.389 J	ΙU
1,1-Dichloroethene	IU	1 U	ΙU
1,4-Dioxane	100 U	100 U	100 U
Chloroform	0.3 U	0.3 U	0.3 U
Chloromethane	1 U	1 U	10
cis-1,2-Dichloroethene	1 U	10	1 U
Tetrachloroethene	10	10	27.2
trans-1,2-Dichloroethene	1 U	IU	1 U
Trichloroethene	1 U	ιυ	1 U
Trichlorofluoromethane	ιU	10	1.0
Vinyl chloride	I UJ	I UJ	1 UJ

Su	•	Table 1 Monitoring Well Samples (continu une 2008	ed)
Location ID Date Sampled	URS-2D 6/16/2008	URS-2S 6/16/2008	URS-2S (DUP) 6/16/2008
Analyte	Volatiles by EPA S	W-846 Method 8260 (ug/L)	
1,1,1-Trichloroethane	1 U	2.2	2.25
1,1-Dichloroethane	0.239 J	0.569 J	0.585 J
1,1-Dichloroethene	1 U	1 U	1 U
1,4-Dioxane	100 U	100 U	100 U
Chloroform	0.3 U	0.204 J	0.197 J
Chloromethane	1 U	1 U	1 U
cis-1,2-Dichloroethene	71.9	0.996 J	0.966 J
Tetrachloroethene	1 U	1 U	1 U
trans-1,2-Dichloroethene	1 U	1 U	1 U
Trichloroethene	1 U	2.19	2.22
Trichlorofluoromethane	1 U	1 U	1 U
Vinyl chloride	0.354 J	1 UJ	1 UJ

	Summary of Dete	Table 1 ected VOCs in Monito November 2	oring Well Samples (continued)	
Location ID Date Sampled	DW3 11/11/2008	DW3 (DUP) 11/11/2008	SW3 11/11/2008	SW4 11/11/2008	SW7 11/11/2008
Analyte	v	olatiles by EPA SW-846 N	1ethod 8260 (ug/L)		
1,1,1-Trichloroethane	I U	1 ()	0.345 J	0.513 J	1.88
I,1-Dichloroethane	10	0.41 J	10	0.825 J	5.04
1,1-Dichloroethene	1 U	τU	10	10	1 U
1,4-Dioxane	100 U	100 U	100 U	100 U	100 M
Chloroform	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
Chloromethane	IU	1 U	1 U	IU	1 M
cis-1,2-Dichloroethene	67.3	65.2	10	2.38	35.3 M
Tetrachloroethene	10	1 U	1 U	0.305 J	0.59 J
trans-1,2-Dichloroethene	1.18	1.09	1 U	0.364 J	0.302 J
Frichloroethene	IU	IU.	0.759 J	12.7	8.15
Frichlorofluoromethane	1 U	ΙU	ΙU	0.651 J	۱U
Vinyl chloride	10	1U	IU	ιU	1.21 M

Key:

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= The analyte was not analyzed for in the sample.

The analyte was positively identified, but the quantitation is an estimation.

UI = The analyte was analyzed for but was not detected. The associated value is an estimate and may be inaccurate or imprecise.

R = The associated quality control indicates that the sample results are not useable.

U = The analyte was analyzed for, but not detected. The associated numerical value is at or below the method detection limit (MDL).

M = Matrix effect: The concentration is estimated due to a matrix effect.

(DUP) = Duplicate sample taken in the field.

Notes:

Bolded values indicate the analyte was detected above the associated MDL.

		Trend Ana	Table Iysis of VO		ndwater		
	Date	C	oncentratio	on of Analy	te in Ground	dwater (ug/	
Well ID	Sampled	TCA	TCE	VC	1,1-DCE	1,2-DCE	-, 1,1-DCA
SW1	Sept. 1986 ¹						
	Jan. 1992 ²	0.5					
	Dec. 1994 ³						
	Nov. 1999 ³						
	May 2000 ³						
	Nov. 2000 ³						
	May 2001 ³						
	Nov. 2001 ³	0.11 J					
	May 2002 ³						
	May 2003 ³						
	Nov. 2003 ³						
	Jun. 2004 ³						
	Nov. 2004 ³						
	Oct. 2005 ³						
	Jun. 2008						
	Nov. 2008	NS	NS	NS	NS	NS	NS
DW1	Jan. 1992 ²	0.6					
	Dec. 1994 ³					1.8 (c)	
	Nov. 1999 ³						
	May 2000 ³						
	Nov. 2000 ³						
	May 2001 ³						
	Nov. 2001 ³						
	May 2002 ³						
	May 2003 ³						
	Nov. 2003 ³						
	Jun. 2004 ³						
	Nov. 2004 ³						
	Oct. 2005 ³						
	Jun. 2008						
	Nov. 2008	NS	NS	NS	NS	NS	NS

	_		Table			N	
	Irenc	a Analysis c	of VOCs in C	Foundwate	er (Continue	ed)	
	Date	C	oncentratio	on of Analy	te in Groun	dwater (µg/l	_)
Well ID	Sampled	TCA	TCE	VC	1,1-DCE	1,2-DCE	, 1,1-DCA
SW3	Sept. 1986 ¹		6				
	Jan. 1992 ²	12	9				5
	Dec. 1994 ³	0.50	1.8				
	Dec. 1995 ³	0.86	2.8			0.44 (c)	
	July 1997 ⁴		1				
	Nov. 1998 ³	0.22	0.81			0.10 (c)	
	Apr. 1999 ³	0.51	0.71			0.17 (c)	
	Nov. 1999 ³	0.29	0.9			0.39 (c)	
	May 2000 ³	0.69	1			1.29 (c)	0.55
	Nov. 2000 ³	0.43	0.9			0.22 (c)	
	May 2001 ³	0.46	0.8			1.29 (c)	0.32
	Nov. 2001 ³	0.32 J	0.5 J				
	May 2002 ³	0.42 J	0.8 J			0.46 J	
	May 2003 ³	0.584 J	0.893 J			1.37 J (c)	0.302 J
	Nov. 2003 ³	0.398 J	0.856 J			0.511 J (c)	
	Jun. 2004 ³	0.9 J	0.94 J			3.7 (c)	0.95 J
	Nov. 2004 ³	0.52 J	1.0	0.26 J		1.5 (c)	0.38 J
	Oct. 2005 ³	0.47 J	0.86 J			0.55 J (c)	
	Jun. 2008	0.661 J	1.31			1.45 (c)	0.403 J
	Nov. 2008	0.345 J	0.759 J				
DW3	Jan. 1992 ²	0.3					0.3
	Dec. 1994 ³			0.28		36 (c)	0.26
	Dec. 1995 ³					5.2 (c)	
	April 1997 ⁴					41 (c)	
	July 1997 ⁴					49 (c)	
	Nov. 1998 ³			0.35		66 (c)	0.34
	Apr. 1999 ³			0.28	0.11	67 (c)	0.35
	Nov 1999 ³						0.11
	May 2000 ³					0.25 (t) 24.98 (c)	0.16
	Nov. 2000 ³					16.85 (c)	
	May 2001 ³					13.29 (c)	

	Trend	Analysis of	Table VOCs in G		r (Continue	d)		
	Date	с С	oncentratio	on of Analy	te in Groun	dwater (ug/		
Well ID	Date Concentration of Analyte in Groundwate Sampled TCA TCE VC 1,1-DCE 1,2-							
DW3 (cont'd)	Nov. 2001 ³					13.58 (c)	1,1-DCA 	
	May 2002 ³					21.08 (c)	0.1 J	
	May 2003 ³							
	Nov. 2003 ³					1.18 J (c)		
	Jun. 2004 ³					1.3 (c)		
	Nov. 2004 ³					2.1 (c)		
	Oct. 2005 ³					3 (c)		
	Jun. 2008 ³					73.1 (c)		
	Nov. 2008 ³					67.3 (c)	0.41 J	
SW4	Jan. 1992 ²	2	97		0.3		0.6	
	Dec. 1994 ³	20	370		2.1	19 (c)	8.5	
	Dec. 1995 ³	34	1200		4.9	2.1 (t) 34 (c)	6.9	
	April 1997 ⁴					71 (c)	7.1	
	July 1997 ⁴	23	290			15 (c)		
	Nov. 1998 ³	8.0	46	0.42	0.82	10 (c)	9.0	
	Apr. 1999 ³	1.9	9.53			1.85 (c)	0.87	
	Nov. 1999 ³	2.13	9.5		0.18	7.15·(c)	7.7	
	May 2000 ³	2.88	8	0.11	0.21	0.49 (t) 4.3 (c)	1.67	
	Nov. 2000 ³	1.14	15.2	1.49	0.29	11.18 (c)	15.25	
	May 2001 ³	3.35	34		0.36	0.38 (t) 3.19 (c)	1.3	
	Nov. 2001 ³	0.88	5.7	0.43 J	0.12 J	5.27 (c)	7.18	
	May 2002 ³	2.54	21.63		0.34 J	2.07 (c)	0.79 J	
	May 2003 ³	3.05 J	9.09 J			3.36 J (c)	1.44 J	
	Nov. 2003 ³	2.03	4.63			1.93 (c)	0.93	
	Jun. 2004 ³	2.8	41		0.57 J	0.11 (t) 3.3 (c)	1.3	

	Trend	Analysis of	Table VOCs in G		r (Continue	d)				
					(0000000	~)				
Well ID	Date	С	Concentration of Analyte in Groundwater (μg/				Concentration of Analyte in Groundwater (µg/L)			L)
Weil IB	Sampled	TCA	TCE	VC	1,1-DCE	1,2-DCE	1,1-DCA			
SW4 (cont'd)	Nov. 2004 ³	3.1	56		0.88 J	0.19 J (t) 4.1 (c)	1.4			
	Oct. 2005 ³	2.2	43		1	6.3 (c)	1.7			
	Jun. 2008 ³	2.98	17.8		0.751 J	4.35 (c)	1.51			
	Nov. 2008 ³	0.513 J	12.7			3.38 (c) 0.364 J (t)	0.825 J			
	Dec. 1994 ³	4.6	15	6.2	1	0.3(t) 150(c)	33			
	Dec. 1995 ³	2.2	7.9	6.8	0.80	130 (c)	20			
	July 1997 ⁴		4			2 (c)				
SW7	Nov. 1998 ³	2.5	11	3.4	0.65	0.28 (t) 82 (c)	12			
	Apr. 1999 ³	1.23	3.95			5.25 (c)	1.46			
	Nov. 1999 ³	1.01	5.7		0.19	18.8·(c)	3.38			
	May 2000 ³	0.67	1.5			0.12 (t) 2.43 (c)	0.71			
	Nov. 2000 ³	0.91	3.8	0.52	0.15	16.06 (c)	3.48			
	May 2001 ³	1.18	1.9			1.46 (c)	0.47			
	Nov. 2001 ³	0.8 J	4.7	0.85 J	0.19 J	0.13 J (t) 25.89 (c)	3.02			
	May 2002 ³	0.87 J	1.65			2.79 (c)	0.47 J			
	May 2003 ³	1.5 J	1.44 J			1.43 J (c)	0.409 J			
	Nov. 2003 ³	0.674 J	1.64			2.76 (c)	0.509			
	Jun. 2004 ³	1	1			1.1 (c)	0.3 J			
	Nov. 2004 ³	1.5	2.1	0.47 J	0.25 J	10 J (c)	1.5 J			
	Oct. 2005 ³	0.73 J	3.1			12 (c)	1.4			
	Jun. 2008 ³	2.5	2.94			6.34 (c)	1.59			

		Trend	Analysis of	Table VOCs in G	_	r (Continue	d)		
Well ID Date Concentration of Analyte in Groundwater (µg/L)									
wei		Sampled	TCA	TCE	VC	1,1-DCE	1,2-DCE	1,1-DCA	
SW7 (cont		Nov. 2008 ^{3,4}	1.88	.88 8.15 1.21 M			0.302 J (t) 35.3 M (c)	5.04	
Key:μg/L=Micrograms per literVC=Vinyl chloride(c)=cis-1,2-Dichloroethene1,1-DCE=1,1-Dichloroethene(t)=trans-1,2-Dichloroethene1,2-DCE=1,2-DichloroetheneTCA=1,1,1-Trichloroethane1,1-DCA=1,1-DichloroethaneTCE=TrichloroetheneDPW=Deep production well(1)=Fred C. Hart Associates(3)=Earth Tech(2)=Argonne National Laboratories(4)=United States Geological Se							ical Services		

Notes: 1. At monitoring well locations where a duplicate groundwater sample was collected, the higher analytical value between the normal and duplicate samples is reported in this table.

2. For 1992 data, the maximum value of either round A or B of sampling was used.

3. Concentrations in bold font exceed the New York State Drinking Water Standard for the associated compound.

4. M = Matrix Effect. The concentration is estimated due to a matrix effect.

		Sur	T nmary of Detect	able 3 ed VOCs in Air (Samples			
			•	ne 2008	I and			
Location ID Date Sampled	SG-33 6/13/2008	SG-34 6/13/2008	SG-35 6/13/2008	SG-36 6/13/2008	SG-37 6/13/2008	SG-38 6/13/2008	SG-38 (DUP) 6/13/2008	SG-39 6/13/2008
Analyte			V	olatiles by EPA M	ethod TO15 (ug/n	13)	·	
1,1,1-Trichloroethane	26	2.1	9.8	9,000	110,000	11 J	2.2 J	1.9
1,1,2-Trichloroethane	0.83 U	0.83 U	0.83 U	0.83 U	20	0.83 U	0.83 U	0.83 U
1,1-Dichloroethane	0.62 U	0.62 U	0.62 U	120	4,900	0.62 U	0.62 U	0.62 U
1,1-Dichloroethene	0.60 U	0.60 U	0.60 U	0.60 U	200	0.60 U	0.60 U	0.60 U
1,2,4-Trimethylbenzene	1.7 U	1.6 U	1.1 U	6.1	13	1.9 U	2.6 U	21
1,2-Dichloroethane	0.62 U	0.62 U	0.62 U	0.62 U	9,600	0.62 U	0.62 U	0.62 U
1,3,5-Trimethylbenzene	1.2	0.75 U	1.5	6.0	14	2.2	2.5	14
1,4-Dichlorobenzene	2.5 U	2.5 U	2.4 U	5.2 U	12	4.5 U	7.4 U	61
2,2,4-Trimethylpentane	6.6	1.2	2.1	3.2	13	0.71 U	1.1	0.71 U
2-Butanone	34	26	33	80	38	6.3 J	16 J	150
2-Hexanone	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	1.2 U	82
4-Ethyltoluene	0.60 J	0.75 U	0.75 U	2.5	4.2	0.55 J	0.90	4.7
4-Methyl-2-pentanone	32	29	51	81	1.2 U	12 J	25 J	1.2 U
Acetone	420 U	130 U	220 U	1,800	1,000 J	40 U	47 U	2,400
Benzene	12	12	18	35	16	6.7 J	9.4 J	29
Carbon disulfide	2.2	1.5	3.1	20	13	1.4	1.5	19
Carbon tetrachloride	0.96 U	0.96 U	0.96 U	0.96 U	8.8	0.96 U	0.96 U	0.96 U
Chloroethane	0.40 U	0.40 U	0.40 U	0.40 U	0.40 U	0.40 U	0.40 U	0.40 U
Chloroform	120	0.74 U	0.74 U	14	230	0.74 U	0.74 U	0.74 U
Chloromethane	0.31 U	0.31 U	0.31 U	0.31 U	0.31 U	0.31 U	0.31 U	0.31 U
cis-1,2-Dichloroethene	0.60 U	0.44 J	0.60 U	0.97	3,100	1.0 J	0.77 J	0.69
Cyclohexane	29	24	23	34	5.4	15	38	0.52 U
Dichlorodifluoromethane	2.8 U	2.8 U	2.7 U	3.4 U	3.5 U	2.3 U	2.4 U	2.3 U
Ethyl Acetate	0.92 U	0.92 U	0.92 U	0.92 U	0.92 U	0.92 U	0.92 U	0.92 U
Ethylbenzene	1.5	1.3	1.7	6.6	12	2.4	2.6	6.4
Freon 113	2.4	1.2 U	1.2 U	6.9	61	1.2 U	1.9	3.0
m,p-Xylene	5.4	4.2	4.6 U	18	33	6.4	7.5	16

Table 3 Summary of Detected VOCs in Air Samples (continued) June 2008											
Location ID Date Sampled	SG-33 6/13/2008	SG-34 6/13/2008	SG-35 6/13/2008	SG-36 6/13/2008	SG-37 6/13/2008	SG-38 6/13/2008	SG-38 (DUP) 6/13/2008	SG-39 6/13/2008			
Analyte			V	olatiles by EPA M	ethod TO15 (ug/m	i 3)					
Methylene chloride	0.53 U	0.53 U	0.53 U	0.53 U	1.6 U	1.1 U	0.78 U	0.53 U			
n-Heptane	32	29	53	87	35	12 J	26 J	170			
n-Hexane	52	42	74	110	60	14 J	29 J	100			
o-Xylene	2.5	1.3	1.3	6.4	19	1.9	2.3	6.5			
Styrene	0.65 U	0.65 U	0.65 U	3.6	0.65 U	0.65 U	1.5	0.65 U			
Tetrachloroethene	14	20	9.9	13	38	12	14	13			
Tetrahydrofuran	0.45 U	0.45 U									
Toluene	13	13	17	30	27	7.7	9.6	31			
trans-1,2-Dichloroethene	0.60 U	0.60 U	0.60 U	0.60 U	80	0.60 U	0.60 U	0.60 U			
Trichloroethene	29	3.3 U	14	1,900	42,000	11	5.7	4.3 U			
Trichlorofluoromethane	1.7 U	1.8 U	1.7 U	2.5 U	2.7 U	1.5 U	1.5 U	1.8 U			
Vinyl acetate	0.54 U	0.54 U	0.54 U	6.8	5.1	0.54 U	0.54 U	0.54 U			

Table 4 Summary of Detected VOCs in Soil Samples November 2008							
Location ID Date Sampled	598G36-3.0-SO1 11 12 2008						
Analyte	Volatiles by EPA Method 8260 (ug/kg)						
1,1,1-Trichloroethane	2.53 J						
Methylene chloride	3.87						
Trichloroethene	0.637 J						

Key: J = The analyte was positively identified, but the quatitation is an estimation.

Notes: Bolded values indicate the analyte was detected above the associated MDL.

Data Validation Report

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1.0 INTRODUCTION

This data quality review pertains to air samples collected in June 2008 at Air Force Plant 59 (AFP-59). Parameters evaluated included the total concentration of volatile organic constituent (VOC). The samples were analyzed by Centek Laboratories, Syracuse, New York.

Data quality review is an after-the-fact technical review of analytical data whereby the quality and usability of the data are determined based on a set of predefined criteria. These criteria depend upon the type of data involved and the purpose for which those data were collected. Data quality review assesses whether and to what extent specified criteria were met, and places restrictions on data use based on quality parameters. The data quality review process can range from a cursory review used to detect out-of-control situations to a detailed evaluation, depending on the analytical protocol, the associated quality control samples collected, and the intended data use.

Specific criteria for data quality review may include, but are not limited to: technical holding times, analysis of blanks, surrogate spike recovery, analysis of duplicates, and reported practical quantitation limits (PQLs). Where applicable, the recommendations of USEPA SW-846 Test Methods for Evaluating Solid Waste (Third Edition, December 1996) or USEPA Methods for Chemical Analysis of Water and Wastes (Revised March 1983) analytical method requirements, USEPA CLP National Functional Guidelines for Organic and Inorganic Data Review (February 1994, Functional Guidelines) data review guidance, and professional judgment. In addition, USEPA Region IX Volatile Organic Compounds in Air (Ambient Air/Soil Vapor/Stack Gas) Samples Collected in Specially-Prepared Canisters and Analyzed by Gas Chromatography/Mass Spectrometry (GC/MS), January 1999, was utilized for this effort.

Table 1 presents the data qualifiers applied during this review effort and their meanings.

Qualifier	Description
J	This is an estimated value.
U	The analyte was analyzed for, but not detected. The associated numerical value is at or below the MDL.

Table 1Data Qualifiers

Table 2 provides a cross-reference list for field sample IDs and lab sample IDs.

Table 2
Field Sample ID/Lab Sample ID Cross Reference
Lot C0801020

Field Sample ID	Lab Sample ID	Field Sample ID	Lab Sample ID
SG33-061308	C0806018-001A	SG38-061308	C0806018-006A
SG34-061308	C0806018-002A	SG39-061308	C0806018-007A
SG35-061308	C0806018-003A	SG38-061308DUP	C0806018-008A
SG36-061308	C0806018-004A	AB-061308	C0806018-009A
SG37-061308	C0806018-005A		

During the data quality review process, laboratory qualified and unqualified data are verified against all available supporting documentation. Based on this review, qualifier codes may be added, deleted, or modified by the validator. (Note: In those cases where the laboratory added a "N.D." flag indicated a non-detect result, and the validator agrees with this flag, and replaces it with a "U" qualifier as noted on the corresponding Results Summary Forms.)Final results are therefore either qualified or unqualified. Changes to the data are reflected on the Results of Analysis Form in Appendix A.

2.0 VOLATILE ORGANIC CONSTITUENTS

Volatile organic constituents were analyzed using EPA Method TO-15. Results were reported in units of ug/m3 and ppbV.

2.1 Holding Times

The samples were analyzed within the prescribed hold time. No qualification is needed.

2.2 Calibration

One Initial Calibration (6/19/2008) was required for this effort. Standards were analyzed at 0.50, 1.0, 5.0, 25.0, 50.0, and 100.0 ppbv in support of the Initial Calibration. The %RSD was less than 25% for all target constituents.

The %D value associated with the continuing calibration standard was less than 20%. No qualification is necessary.

Finally, unless qualified otherwise, the validator qualifies U those values reported ND and agrees with the J qualifier by the laboratory.

2.3 Laboratory Control Samples

Constituent recoveries from the associated laboratory control sample were within control limits for all target constituents. No qualification is needed.

2.4 Blanks

No positive results were observed above the laboratory reporting limits in the associated method blanks. No qualification is needed.

AB-061308 exhibited 0.160 ppbv of 1,2,4-trimethylbenzene, 0.240 ppbv of 1,4-dichlorobenzene, 31 ppbv of acetone, 0.190 ppbv of benzene, 0.480 ppbv of chloromethane, 0.150 ppbv of ethyl acetate, 0.280 ppbv of Freon 11, 0.590 ppbv of Freon 12, 0.120 ppbv of hexane, 0.170 ppbv of m&p-xylene, 0.540 ppbv of methyl ethyl ketone, 0.240 ppbv of methylene chloride, 0.390 ppbv of toluene, and 0.170 ppbv of trichloroethene.

The validator qualifies U any positive result less than or equal to 0.8 ppbv in 1,2,4-trimethylbenzene; 1.2 ppbv in 1,4-dichlorobenzene, 310 ppbv in acetone, 0.95 ppbv in benzene, 2.4 ppbv in chloromethane, 0.75 ppbv in ethyl acetate, 1.4 ppbv in Freon 11, 2.95 ppbv in Freon 12, 0.60 ppbv in hexane, 0.85 ppbv in m&p-xylene, 5.40 ppbv in methyl ethyl ketone, 2.4 ppbv in methylene chloride, 1.95 ppbv in toluene, and 0.85 ppbv in trichlorethene.

2.5 Matrix Spike/Matrix Spike Duplicate

Site-specific MS/MSD samples were not included with either batch. No qualification is needed based on this information.

2.6 Surrogate Recovery

Surrogate bromofluorobenzene was high and did not meet criteria for samples SG37-061308. This is most likely due to matrix effects.

The sample was analyzed further as serial dilutions with results meeting criteria.

2.7 Internal Standards

Internal standard was high and did not meet criteria for samples SG36-061308, SG37-061308, SG38-061308, SG39-061308 were analyzed with further dilutions with results meeting criteria. No qualification is necessary.

2.8 Duplicates

A field duplicate was collected for sample SG38-061308. One of two criteria was followed when evaluating field duplicates, depending on the amount detected. If the amount detected was greater than five times the reporting limit (RL), then the relative percent difference (RPD) should have been less than 25 percent. If the amount was less than five times the RL, then the difference between the duplicate and the sample concentrations should have been less than the RL. Agreement is excellent and no qualification is needed except for 1,1,1-trichloroethane, 1,2,4-trimethylbenzene, 1,4-dichlorobenzene, 4-ethyltoluene, benzene, cis-1,2-dichloroethene, cyclohexane, heptane, hexane, methyl ethyl ketone, methyl isobutyl ketone, methylene chloride, and trichloroethene . The validator qualifies J these analytes.

A comparison of field sample and duplicate is presented in Table 3.

Analyte	Reporting Limit	SG38-061308	SG38-061308 DUP	RPD
1,1,1-	0.15	2.0	0.40	133%
Trichloroethane				
1,2,4-	0.15	0.38	0.52	31%
Trimethylbenzene				
1,3,5-	0.15	0.45	0.50	11%
Trimethylbenzene		2		
1,4-	0.15	0.73	1.2	49%
Dichlorobenzene				
2,2,4-	0.15	ND	0.24	Not calculated
Trimethylpentane				
4-ethyltoluene	0.15	0.11J	0.18	48%
Acetone	3.0	16	20	22%
Benzene	0.15	2.1	2.9	32%
Carbon disulfide	0.15	0.44	0.46	4.4%
Cis-1,2-	0.15	0.25	0.19	30%
Dichloroethene				
Cyclohexane	1.5	4.2	11	89%

Table 3 Duplicate Comparison (ppbv)

Data Validation Review June 2008 Air Sampling Air Force Plant 59(AFP-59)

Ethylbenzene	0.15	0.55	0.60	8.7%
Freon 11	0.15	0.26	0.27	3.8%
Freon 113	0.15	ND	0.25	Not calculated
Freon 12	0.15	0.45	0.47	4.3%
Heptane	1.5	3.0	6.2	69%
Hexane	1.5	3.9	8.1	70%
m&p-xylene	0.30	1.4	1.7	19%
Methyl ethyl ketone	0.30	2.1	5.3	86%
Methyl isobutyl	3.0	3.0	6.1	67%
ketone				
Methylene chloride	0.15	0.32	0.22	37%
o-xylene	0.15	0.43	0.51	17%
Styrene	0.15	ND	0.35	Not calculated
Tetrachloroethylene	0.15	1.8	2.0	10%
Toluene	1.5	2.0	2.5	22%
Trichloroethene	0.15	2.1	1.0	71%

2.9 Summary

The data are acceptable with validator-assigned qualifiers.

Date: 20-Jun-08

CLIENT:	Earth Tech		(Client Sample ID:		
Lab Order:	C0806018			Tag Number:	363, 4	47
Project:	AFB 59 (BAE)			Collection Date:	6/13/2	008
Lab ID:	C0806018-001A			Matrix:	AIR	
Analyses		Result	Limit Qua	Units	DF	Date Analyzed
FIELD PARAM	ETERS		FLD			Analyst:
Vacuum Readir	ng "Hg	-2		"Hg		6/13/2008
HELIUM LEAK	TEST		GC			Analyst: RJP
Helium		ND	1.0	%	1	6/19/2008
1UG/M3 BY ME	ETHOD TO15		TO-15			Analyst: RJP
1,1,1-Trichloroe	ethane	4.6	3.0	ppbV	20	6/19/2008 1:56:00 AM
1,1,2,2-Tetrach	loroethane	ND V	0.15	ppbV	1	6/18/2008 11:39:00 AM
1,1,2-Trichloroe	ethane	ND	0.15	ppbV	1	6/18/2008 11:39:00 AM
1,1-Dichloroeth	ane	ND	0.15	ppbV	1	6/18/2008 11:39:00 AM
1,1-Dichloroeth	iene	ND	0.15	ppbV	1	6/18/2008 11:39:00 AM
1,2,4-Trichlorot	penzene	ND 🗸	0.15	ppbV	1	6/18/2008 11:39:00 AM
1,2,4-Trimethyl	benzene	0.34 U	0.15	ppbV	1	6/18/2008 11:39:00 AM
1,2-Dibromoeth	ane	ND V	0.15	ppbV	1	6/18/2008 11:39:00 AM
1,2-Dichlorober	nzene	ND i	0.15	ppbV	1	6/18/2008 11:39:00 AM
1,2-Dichloroeth	ane	ND	0.15	ppbV	1	6/18/2008 11:39:00 AM
1,2-Dichloropro	pane	ND 🗸	0.15	ppbV	1	6/18/2008 11:39:00 AM
1,3,5-Trimethyl	benzene	0.25	0.15	ppbV	1	6/18/2008 11:39:00 AM
1,3-butadiene		ND U	0.15	ppbV	1	6/18/2008 11:39:00 AM
1,3-Dichlorober	nzene	ND \vee	0.15	ppbV	1	6/18/2008 11:39:00 AM
1,4-Dichlorober	nzene	0.41 \) 0.15	ppbV	1	6/18/2008 11:39:00 AM
1,4-Dioxane		ND V	0.30	ppbV	1	6/18/2008 11:39:00 AM
2,2,4-trimethylp	pentane	1.4	0.15	ppbV	1	6/18/2008 11:39:00 AM
4-ethyltoluene		0.12	0.15 J	ppbV	1	6/18/2008 11:39:00 AN
Acetone		170 V	36	ppbV	120	6/19/2008 9:42:00 PM
Allyl chloride		ND 🗸	0.15	ppbV	1	6/18/2008 11:39:00 AN
Benzene		3.8	3.0	ppbV	20	6/19/2008 1:56:00 AM
Benzyl chloride	3	ND V	0.15	ppbV	1	6/18/2008 11:39:00 AN
Bromodichloro		ND	0.15	ppbV	1	6/18/2008 11:39:00 AM
Bromoform		ND	0.15	ppbV	1	6/18/2008 11:39:00 AM
Bromomethane	2	ND 🔱	0.15	ppbV	1	6/18/2008 11:39:00 AM
Carbon disulfid		0.70	0.15	ppbV	1	6/18/2008 11:39:00 AM
Carbon tetrach		ND V		ppbV	1	6/18/2008 11:39:00 AM
Chlorobenzene		ND	0.15	ppbV	1	6/18/2008 11:39:00 AM
Chloroethane	-	ND	/ 0.15	ppbV	1	6/18/2008 11:39:00 AM
Chloroform		23	3.0	ppbV	20	6/19/2008 1:56:00 AM
Chloromethane	2	ND U		ppbV	1	6/18/2008 11:39:00 AM
cis-1.2-Dichlore		ND	0.15	ppbV	1	6/18/2008 11:39:00 AM
cis-1,3-Dichlor		ND	0.15	ppbV	1	6/18/2008 11:39:00 AM
Cyclohexane	+F. = \$ 2	8.2	3.0	ppbV	20	6/19/2008 1:56:00 AM
Dibromochloro	methane	ND L		ppbV	1	6/18/2008 11:39:00 AM

Qualifiers:

B Analyte detected in the associated Method Blank
 Holding times for preparation or analysis exceeded

E Value above quantitation range

J Analyte detected at or below quantitation limits

JN Non-routine analyte. Quantitation estimated.

S Spike Recovery outside accepted recovery limits

ND Not Detected at the Reporting Limit

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1/22/03

Date: 20-Jun-08

CLIENT:	Earth Tech				C	Client Sample ID:	SG33-	-061308
Lab Order:	C0806018					Tag Number:	363, 4	47
Project:	AFB 59 (BAE)					Collection Date:	6/13/2	800
Lab ID:	C0806018-001A					Matrix:	AIR	
Analyses		Result		Limit	Qual	Units	DF	Date Analyzed
1UG/M3 BY ME	ETHOD TO15			то)-15			Analyst: RJP
Ethyl acetate		ND	U	0.25		ppbV	1	6/18/2008 11:39:00 AM
Ethylbenzene		0.35		0.15		ppbV	1	6/18/2008 11:39:00 AM
Freon 11		0.30	\vee	0.15		ppbV	1	6/18/2008 11:39:00 AM
Freon 113		0.31		0.15		ppbV	1	6/18/2008 11:39:00 AM
Freon 114		ND	\mathcal{O}	0.15		ppbV	1	6/18/2008 11:39:00 AN
Freon 12		0.55	υ	0.15		ppbV	1	6/18/2008 11:39:00 AM
Heptane		7.8		3.0		ppbV	20	6/19/2008 1:56:00 AM
Hexachioro-1,3	-butadiene	ND	\mathcal{V}	0.15		ppbV	1	6/18/2008 11:39:00 AN
Hexane		14		3.0		ppbV	20	6/19/2008 1:56:00 AM
Isopropyl alcoh	ol	ND	\mathcal{O}	0.15		ppbV	1	6/18/2008 11:39:00 AN
m&p-Xylene		1.2		0.30		ppbV	1	6/18/2008 11:39:00 AM
Methyl Butyl Ke	etone	ND	\mathcal{O}	0.30		ppbV	1	6/18/2008 11:39:00 AM
Methyl Ethyl Ke	etone	11		6.0		ppbV	20	6/19/2008 1:56:00 AM
Methyl Isobutyl	Kelone	7.6		6.0		ppbV	20	6/19/2008 1:56:00 AM
Methyl tert-buty	/l ether	ND	\mathcal{O}^{-}	0.15		ppbV	1	6/18/2008 11:39:00 AM
Methylene chlo	ride	ND	U	0.15		ppbV	1	6/18/2008 11:39:00 AM
o-Xylene		0.56		0.15		ppbV	1	6/18/2008 11:39:00 AM
Propylene		ND	\cup	0.15		ppbV	1	6/18/2008 11:39:00 AM
Styrene		ND	J	0.15		ppbV	1	6/18/2008 11:39:00 AM
Tetrachloroethy	ylene	2.0		0.15		ppbV	1	6/18/2008 11:39:00 AM
Tetrahydrofura	n	ND	V	0.15		ppbV	1	6/18/2008 11:39:00 AM
Toluene		3.4		3.0		ppbV	20	6/19/2008 1:56:00 AM
trans-1,2-Dichl	oroethene	ND	J	0.15		ppbV	1	6/18/2008 11:39:00 AM
trans-1,3-Dichl		ND	\mathcal{O}	0.15		ppbV	1	6/18/2008 11:39:00 AM
Trichtoroethene		5.4		3.0		ppbV	20	6/19/2008 1:56:00 AM
Vinyl acetate		ND	\cup	0.15		ppbV	1	6/18/2008 11:39:00 AM
Vinyl Bromide		ND	i	0.15		ppbV	1	6/18/2008 11:39:00 AM
Vinyl chloride		ND		0.15		ppbV	1	6/18/2008 11:39:00 AM
•	fluorobenzene	115	-	70-130		%REC	1	6/18/2008 11:39:00 AM

Qualifiers:

B Analyte detected in the associated Method Blank

.

- 11 Holding times for preparation or analysis exceeded
- JN Non-routine analyte. Quantitation estimated.
- S Spike Recovery outside accepted recovery limits

E Value above quantitation range

J Analyte detected at or below quantitation limits

ND Not Detected at the Reporting Limit

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1/22/08

Date: 20-Jun-08

CLIENT:	Earth Tech				C	lient Sample ID:	SG34-	061308
Lab Order:	C0806018					Tag Number:	242, 2	79
Project:	AFB 59 (BAE)					Collection Date:		
						Matrix:		
Lab ID:	C0806018-002A					Matrix.		
Analyses		Result	I	Jimit Q)ual	Units	ÐF	Date Analyzed
FIELD PARAM	ETERS			FLC)			Analyst:
Vacuum Readir	ng "Hg	-2				"Нд		6/13/2008
HELIUM LEAK	TEST			GC				Analyst: RJP
Helium		ND		1.0		%	1	6/19/2008
1UG/M3 BY ME	THOD TO15			TO-1	5			Analyst: RJP
1,1,1-Trichloroe	thane	0.37		0.15		ppbV	1	6/18/2008 12:16:00 PM
1,1,2,2-Tetrach	loroethane	ND ¹	J	0.15		ppbV	1	6/18/2008 12:16:00 PM
1,1,2-Trichloroe	thane	ND	1	0.15		ppbV	1	6/18/2008 12:16:00 PM
1,1-Dichloroeth		ND		0.15		ppbV	1	6/18/2008 12:16:00 PM
1.1-Dichloroeth	ene	ND		0.15		ppbV	1	6/18/2008 12:16:00 PM
1,2,4-Trichlorob		ND	L	0.15		ppbV	1	6/18/2008 12:16:00 PM
1,2.4-Trimethyll		0.32	V	0.15		ppbV	1	6/18/2008 12:16:00 PM
1,2-Dibromoeth			V	0.15		ppbV	1	6/18/2008 12:16:00 PM
1,2-Dichlorober		ND	1	0.15		ppbV	1	6/18/2008 12:16:00 PM
1,2-Dichloroeth		ND		0.15		ppbV	1	6/18/2008 12:16:00 PM
1,2-Dichloropro		ND		0.15		ppbV	1	6/18/2008 12:16:00 PM
1,3,5-Trimethyl		ND		0.15		ppbV	1	6/18/2008 12:16:00 PM
1,3-butadiene		ND		0.15		ppbV	1	6/18/2008 12:16:00 PM
1,3-Dichlorober	ozene		Ţ.	0.15		ppbV	1	6/18/2008 12:16:00 PM
1,4-Dichlorober			V	0.15		ppbV	1	6/18/2008 12:16:00 PM
1.4-Dioxane			V	0.30		ppbV	1	6/18/2008 12:16:00 PM
2,2;4-trimethylp	entane	0.25		0.15		ppb∨	1	6/18/2008 12:16:00 PM
4-ethyltoluene	Si dano		V İ	0.15		ppbV	1	6/18/2008 12:16:00 PM
Acetone			0	12		ppbV	40	6/19/2008 8:32:00 PM
Alfyl chloride		ND	· .	0.15		ppbV	1	6/18/2008 12:16:00 PM
Benzene		3.7	-	1.5		ppbV	10	6/19/2008 2:31:00 AM
Benzyl chloride		ND	.1	0.15		ppbV	1	6/18/2008 12:16:00 PM
Bromodichloror		ND		0.15		ppbV	1	6/18/2008 12:16:00 PM
Bromoform	nesitane	ND		0.15		ppbV	1	6/18/2008 12:16:00 PM
Bromomethane	3	ND		0.15		ppbV	1	6/18/2008 12:16:00 PM
Carbon disulfid		0.48	\bigvee	0.15		ppbV	1	6/18/2008 12:16:00 PM
Carbon tetrach		ND	V.	0.15		ppbV	1	6/18/2008 12:16:00 PM
Chlorobenzene		ND		0.15		ppbV	1	6/18/2008 12:16:00 PM
Chloroethane	•	ND		0.15		ppbV	1	6/18/2008 12:16:00 PM
Chloroform		ND		0.15		ppbV	1	6/18/2008 12:16:00 PM
Chloromethane	2		<u>_</u>	0.15		ppbV	1	6/18/2008 12:16:00 PM
cis-1,2-Dichlor		0.11	~	0.15	J	ppbV	1	6/18/2008 12:16:00 PM
cis-1.3-Dichlor		ND	0	0.15	Ý	ppbV	1	6/18/2008 12:16:00 PM
Cyclohexane	phohene	6.9	14	1,5		ppbV	10	6/19/2008 2:31:00 AM
Dibromochloro	mathana	ND	. ,	0.15		ppbv pdq	1	6/18/2008 12:16:00 PM

Qualifiers: B Analyte detected in the associated Method Blank

31

E Value above quantitation range

J Analyte detected at or below quantitation limits

JN Non-routine analyte. Quantitation estimated.

Holding times for preparation or analysis exceeded

S Spike Recovery outside accepted recovery limits

ND Not Detected at the Reporting Limit

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1122/08 oc

Date: 20-Jun-08

CLIENT:	Earth Tech				(Client Sample ID:	SG34-	061308
Lab Order:	C0806018					Tag Number:	242, 2	79
Project:	AFB 59 (BAE)					Collection Date:	6/13/2	008
Lab ID:	C0806018-002A					Matrix:	AIR	
Analyses		Result		Limit	Qual	Units	DF	Date Analyzed
1UG/M3 BY ME	ETHOD TO15		-	то	-15			Analyst: RJP
Ethyl acetate		ND	U	0.25		ppbV	1	6/18/2008 12:16:00 PM
Ethylbenzene		0.29		0.15		ppbV	1	6/18/2008 12:16:00 PM
Freon 11		0.31	∇	0.15		ppbV	1	6/18/2008 12:16:00 PM
Freon 113		ND	\mathcal{O}	0.15		ppbV	1	6/18/2008 12:16:00 PM
Freon 114		ND	J	0.15		ppbV	1	6/18/2008 12:16:00 PM
Freon 12		0.55	Ū	0.15		ppbV	1	6/18/2008 12:16:00 PN
Heptane		7.0		1.5		ppbV	10	6/19/2008 2:31:00 AM
Hexachloro-1,3	-butadiene	ND	\cup	0.15		ppbV	1	6/18/2008 12:16:00 PN
Hexane		12		1.5		ppbV	10	6/19/2008 2:31:00 AM
Isopropyl alcoh	ol	ND	V	0.15		ppbV	1	6/18/2008 12:16:00 PN
m&p-Xylene		0.95		0.30		ppbV	1	6/18/2008 12:16:00 PM
Methyl Butyl Ke	etone	ND	V	0.30		ppbV	1	6/18/2008 12:16:00 PN
Methyl Ethyl Ke	etone	8.7		3.0		ppbV	10	6/19/2008 2:31:00 AM
Methyl Isobutyl	Ketone	7.0		3.0		ppbV	10	6/19/2008 2:31:00 AM
Methyl tert-buty	/l ether	ND	Ú	0.15		ppbV	1	6/18/2008 12:16:00 PN
Methylene chlo	ride	ND	J	0.15		ppbV	1	6/18/2008 12:16:00 PN
o-Xylene		0.30		0.15		ppbV	1	6/18/2008 12:16:00 PM
Propylene		ND	J	0.15		ppbV	1	6/18/2008 12:16:00 PM
Styrene		ND	\downarrow	0.15		ppbV	1	6/18/2008 12:16:00 PM
Tetrachloroethy	ylene	2.9		1.5		ppbV	10	6/19/2008 2:31:00 AM
Tetrahydrofurai	n	ND	J	0.15		Vdqq	1	6/18/2008 12:16:00 PM
Toluene		3.5		1.5		ppbV	10	6/19/2008 2:31:00 AM
trans-1,2-Dichle	oroethene	ND	V	0.15		ppbV	1	6/18/2008 12:16:00 PM
trans-1,3-Dichle	oropropene	ND	\cup	0.15		ppbV	1	6/18/2008 12:16:00 PM
Trichloroethene	9	0.61	\mathbf{U}	0.15		ppbV	1	6/18/2008 12:16:00 PM
Vinyl acetate		ND	\overline{U}	0.15		ppb∨	1	6/18/2008 12:16:00 PN
Vinyl Bromide		ND	I	0.15		ppbV	1	6/18/2008 12:16:00 PM
Vinyl chloride		ND	1	0.15		Vđqq	1	6/18/2008 12:16:00 PM
,	fluorobenzene	116	~	70-130		%REC	1	6/18/2008 12:16:00 PM

Qualifiers:

H

JN

S

B Analyte detected in the associated Method Blank

Spike Recovery outside accepted recovery limits

Non-routine analyte. Quantitation estimated.

- Holding times for preparation or analysis exceeded
- E Value above quantitation range
- Analyte detected at or below quantitation limits 1
- ND Not Detected at the Reporting Limit

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7/22/08

Date: 20-Jun-08

CLIENT:	Earth Tech			C	lient Sample ID:	SG35-	-061308
Lab Order:	C0806018				Tag Number:	369, 1	79
Project:	AFB 59 (BAE)				Collection Date:	6/13/2	008
Lab ID:	C0806018-003A				Matrix:	AIR	
LAD ID:	0000010-0057						
Analyses		Result		Limit Qual	Units	DF	Date Analyzed
FIELD PARAMI	ETERS			FLD			Analyst:
Vacuum Readir	ng "Hg	-3			"Hg		6/13/2008
HELIUM LEAK	TEST			GC			Analyst: RJP
Helium		ND		1.0	%	1	6/19/2008
1UG/M3 BY ME	ETHOD TO15			TO-15			Analyst: RJF
1,1,1-Trichloroe	ethane	1.8		0.15	ppbV	1	6/18/2008 1:30:00 PM
1,1,2,2-Tetrach	ioroethane	ND	\mathcal{V}	0.15	ppbV	1	6/18/2008 1:30:00 PM
1,1,2-Trichloroe	ethane	ND	ì	0.15	ppbV	1	6/18/2008 1:30:00 PM
1,1-Dichloroeth	ane	ND		0.15	ppbV	1	6/18/2008 1:30:00 PM
1,1-Dichloroeth	ene	ND		0.15	ppbV	1	6/18/2008 1:30:00 PM
1,2,4-Trichlorot	benzene	ND	\checkmark	0.15	ppbV	1	6/18/2008 1:30:00 PM
1,2,4-Trimethyll	benzene	0.23	v	0.15	ppbV	1	6/18/2008 1:30:00 PM
1,2-Dibromoeth	ane	ND	V	0.15	ppbV	1	6/18/2008 1:30:00 PM
1,2-Dichlorober	nzene	ND	i	0.15	ppbV	1	6/18/2008 1:30:00 PM
1.2-Dichloroeth	ane	ND		0.15	Vdqq	1	6/18/2008 1:30:00 PM
1,2-Dichloropro	pane	ND	Į,	0.15	ppbV	1	6/18/2008 1:30:00 PM
1,3,5-Trimethyll	benzene	0.31		0.15	ppbV	1	6/18/2008 1:30:00 PM
1,3-butadiene		ND	\cup	0.15	ppbV	1	6/18/2008 1:30:00 PM
1,3-Dichlorober	nzene	ND	\lor	0.15	ppbV	1	6/18/2008 1:30:00 PM
1,4-Dichlorober	nzene	0.39	\mathcal{V}	0.15	ppbV	1	6/18/2008 1:30:00 PM
1,4-Dioxane		ND	V	0.30	ppbV	1	6/18/2008 1:30:00 PM
2,2,4-trimethylp	entane	0.45		0.15	ppbV	1	6/18/2008 1:30:00 PM
4-ethyltoluene		ND	U	0.15	ppbV	1	6/18/2008 1:30:00 PM
Acetone		90	\mathcal{O}	12	ppbV	40	6/19/2008 9:07:00 PM
Allyl chloride		ND	U	0.15	ppbV	1	6/18/2008 1:30:00 PM
Benzene	··· .	5.4		1.5	ppbV	10	6/19/2008 3:41:00 AM
Benzyl chloride	2	ND	\mathcal{V}	0.15	ppbV	1	6/18/2008 1:30:00 PM
Bromodichloror		ND	i	0.15	ppbV	1	6/18/2008 1:30:00 PM
Bromoform		ND		0.15	ppbV	1	6/18/2008 1:30:00 PM
Bromomethane	2	ND	V	0.15	ppbV	1	6/18/2008 1:30:00 PM
Carbon disulfid		0.99		0.15	ppbV	1	6/18/2008 1:30:00 PM
Carbon tetrach		ND	\mathbf{V}^{-}	0.15	ppbV	1	6/18/2008 1:30:00 PM
Chlorobenzene		ND	1	0.15	ppbV	1	6/18/2008 1:30:00 PM
Chloroethane		ND		0.15	ppbV	1	6/18/2008 1:30:00 PM
Chloroform		ND		0.15	ppbV	1	6/18/2008 1:30:00 PM
Chloromethane	è	ND		0.15	ppbV	1	6/18/2008 1:30:00 PM
cis-1,2-Dichloro		ND		0.15	ppbV	1	6/18/2008 1:30:00 PM
cis-1,3-Dichlor		ND	$\stackrel{!}{\checkmark}$	0.15	ppbV	1	6/18/2008 1:30:00 PN
Cyclohexane	-1 1	6.6		1.5	ppbV	10	6/19/2008 3:41:00 AM
Dibromochloro	methane	ND	V	0.15	ppbV	1	6/18/2008 1:30:00 PM

Qualifiers:

H

B Analyte detected in the associated Method Blank Holding times for preparation or analysis exceeded E Value above quantitation range

ND Not Detected at the Reporting Limit

Analyte detected at or below quantitation limits J

JN Non-routine analyte. Quantitation estimated.

Spike Recovery outside accepted recovery limits S

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1122108

Date: 20-Jun-08

CLIENT:	Earth Tech				(lient Sample ID:	SG35-	061308
Lab Order:				Tag Number:	369, 179			
Project:	AFB 59 (BAE)					Collection Date:	6/13/2	.008
Lab ID:	C0806018-003A					Matrix:	AIR	
Analyses		Result		Limit	Qual	Units	DF	Date Analyzed
1UG/M3 BY ME	ETHOD TO15			то	-15			Analyst: RJI
Ethyl acetate		ND	Ú	0.25		ppbV	1	6/18/2008 1:30:00 PM
Ethylbenzene		0.38		0.15		ppbV	1	6/18/2008 1:30:00 PM
Freon 11		0.30	V.	0.15		ppbV	1	6/18/2008 1:30:00 PM
Freon 113		ND	V.	0.15		ppbV	1	6/18/2008 1:30:00 PM
Freon 114		ND	\mathcal{V}	0.15		ppbV	1	6/18/2008 1:30:00 PM
Freon 12		0.54	\mathcal{V}	0.15		ppbV	1	6/18/2008 1:30:00 PM
Heptane		13		1.5		ppbV	10	6/19/2008 3:41:00 AM
Hexachloro-1,3-butadiene		NĎ	\mathcal{V}	0.15		ppbV	1	6/18/2008 1:30:00 PM
Hexane		21		1.5		ppbV	10	6/19/2008 3:41:00 AM
Isopropyl alcohol		ND	\mathcal{O}	0.15		ppbV	1	6/18/2008 1:30:00 PM
m&p-Xylene		1.0	2	0.30		ppbV	1	6/18/2008 1:30:00 PM
Methyl Butyl Ketone		ND	\mathcal{V}	0.30		ppbV	1	6/18/2008 1:30:00 PM
Methyl Ethyl Ketone		11		3.0		ppbV	10	6/19/2008 3:41:00 AM
Methyl Isobutyl Ketone		12		3.0		ppbV	10	6/19/2008 3:41:00 AN
Methyl tert-butyl ether		ND	\mathcal{O}	0.15		ppbV	1	6/18/2008 1:30:00 PN
Methylene chloride		ND	\mathcal{O}	0.15		ppbV	1	6/18/2008 1:30:00 PM
o-Xylene		0.30		0.15		ppbV	1	6/18/2008 1:30:00 PM
Propylene		ND	\mathcal{I}	0.15		ppbV	1	6/18/2008 1:30:00 PM
Styrene		ND	\cup	0.15		ppbV	1	6/18/2008 1:30:00 PN
Tetrachloroethylene		1.4		0.15		ppbV	1	6/18/2008 1:30:00 PM
Tetrahydrofuran		ND	\cup	0.15		ppbV	1	6/18/2008 1:30:00 PM
Toluene	Toluene			1.5		ppbV	10	6/19/2008 3:41:00 AM
trans-1,2-Dichle	trans-1,2-Dichloroethene		\cup	0.15		ppbV	1	6/18/2008 1:30:00 PN
trans-1,3-Dichle	trans-1,3-Dichloropropene		\mathcal{O}	0.15		ppbV	1	6/18/2008 1:30:00 PM
Trichloroethene		2.6		1.5		ppbV	10	6/19/2008 3:41:00 AN
Vinyl acetate		ND	\mathcal{U}	0.15		vdqq	1	6/18/2008 1:30:00 PN
Vinyl Bromide		ND	Ì	0.15		ppbV	1	6/18/2008 1:30:00 PM
Vinyl chloride		ND	$\langle \cdot \rangle$	0.15		ppbV	1	6/18/2008 1:30:00 PM
Surr: Bromofluorobenzene		117		70-130		%REC	1	6/18/2008 1:30:00 PM

Qualifiers:

П

- B Analyte detected in the associated Method Blank
- - Holding times for preparation or analysis exceeded
- Non-routine analyte. Quantitation estimated. JN
- Spike Recovery outside accepted recovery limits S

E Value above quantitation range

Analyte detected at or below quantitation limits J

ND Not Detected at the Reporting Limit

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1122/08

Date: 20-Jun-08

CLIENT: Lab Order: Project: Lab ID:	Earth Tech C0806018 AFB 59 (BAE) C0806018-004A		(Client Sample ID: Tag Number: Collection Date: Matrix:	356, 18 6/13/20	31
Analyses		Result	Limit Qual	Units	DF	Date Analyzed
		-2	FLD	"Hg		Analyst: 6/13/2008
Vacuum Readir	iy riy	-2		19		
HELIUM LEAK	TEST	ND	GC 1.0	%	1	Analyst: RJP 6/19/2008
Helium		140	1.0	70		010/2000
1UG/M3 BY ME	THOD TO15		TO-15			Analyst: RJP
1,1,1-Trichloroe	thane	1600	240	ppbV	1620	6/19/2008 10:17:00 PM
1,1,2,2-Tetrach		ND V	0.15	ppbV	1	6/18/2008 2:07:00 PM
1,1,2-Trichloroe	lhane	ND 🗸		ppbV	1	6/18/2008 2:07:00 PM
1,1-Dichloroeth	ane	29	1.5	ppbV	10	6/19/2008 4:51:00 AM
1,1-Dichloroeth	ene	ND 🤳	0.15	ppbV	1	6/18/2008 2:07:00 PM
1,2,4-Trichlorob	enzene	ND V	0.15	ppbV	1	6/18/2008 2:07:00 PM
1,2,4-Trimethyll	benzene	1,2	0.15	ppbV	1	6/18/2008 2:07:00 PM
1,2-Dibromoeth	ane	ND 🗸	0.15	ppbV	1	6/18/2008 2:07:00 PM
1,2-Dichlorober	nzene	ND	0.15	opbV	1	6/18/2008 2:07:00 PM
1,2-Dichloroeth	ane	ND	0.15	ppbV	1	6/18/2008 2:07:00 PM
1,2-Dichloropro	pane	ND 🕹	0.15	ppbV	1	6/18/2008 2:07:00 PM
1,3,5-Trimethyll	benzene	1.2	0.15	ppbV	1	6/18/2008 2:07:00 PM
1,3-butadiene		ND V		ppbV	1	6/18/2008 2:07:00 PM
1,3-Dichlorober	nzene	ND	0.15	ppbV	1	6/18/2008 2:07:00 PM
1,4-Dichlorober	izene	0.85 🕔		ppbV	1	6/18/2008 2:07:00 PM
1,4-Dioxane		ND U	i 0.30	ppbV	1	6/18/2008 2:07:00 PM
2.2.4-trimethylp	entane	0.68	0.15	ppbV	1	6/18/2008 2:07:00 PM
4-ethyltoluene		0.50	0.15	ppbV	1	6/18/2008 2:07:00 PM
Acetone		760	490	ppbV	1620	6/19/2008 10:17:00 PM
Allyl chloride		ND V	0.15	vdqq	1	6/18/2008 2:07:00 PM
Benzene	• "	11	1.5	ppbV	10	6/19/2008 4:51:00 AM
Benzyl chloride	,	ND V	0.15	ppbV	1	6/18/2008 2:07:00 PM
Bromodichloror	methane	ND (0.15	Vdqq	1	6/18/2008 2:07:00 PM
Bromoform		ND	0.15	ppbV	1	6/18/2008 2:07:00 PM
Bromomethane	;	ND 💐	/ 0.15	ppbV	1	6/18/2008 2:07:00 PM
Carbon disulfid	e	6.3	1.5	ppbV	10	6/19/2008 4:51:00 AM
Carbon tetrach	loride	ND 🔍	× 0.15	ppbV	1	6/18/2008 2:07:00 PM
Chlorobenzene		ND	0.15	ppbV	1	6/18/2008 2:07:00 PM
Chloroethane		ND	/ 0.15	ppbV	1	6/18/2008 2:07:00 PM
Chloroform		2.9	1.5	ppbV	10	6/19/2008 4:51:00 AM
Chioromethane	9	ND 🕔	/ 0.15	ppbV	1	6/18/2008 2:07:00 PM
cis-1,2-Dichloro	pethene	0.24	0.15	ppbV	1	6/18/2008 2:07:00 PM
cis-1.3-Dichloro	opropene	ND V	0.15	ppbV	1	6/18/2008 2:07:00 PM
Cyclohexane		9.6	1.5	ppbV	10	6/19/2008 4:51:00 AM
Dibromochloro	methane	ND U	<i>i</i> 0.15	ppbV	1	6/18/2008 2:07:00 PM

Qualifiers:

JN

B Analyte detected in the associated Method Blank
 H Holding times for preparation or analysis exceeded

E Value above quantitation range

ND Not Detected at the Reporting Limit

J Analyte detected at or below quantitation limits

Non-routine analyte, Quantitation estimated.

S Spike Recovery outside accepted recovery limits

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7122/03

Date: 20-Jun-08

CLIENT:	Earth Tech				C	lient Sample ID:	SG36-	061308
Lab Order:	C0806018					Tag Number:	356, 18	31
Project:	AFB 59 (BAE)					Collection Date:	6/13/20	008
Lab ID:	C0806018-004A					Matrix:	AIR	
Analyses		Result		Limit	Qual	Units	DF	Date Analyzed
1UG/M3 BY ME	ETHOD TO15			тс	-15			Analyst: RJP
Ethyl acetate		ND	U	0.25		ppbV	1	6/18/2008 2:07:00 PM
Ethylbenzene		1.5		0.15		ppbV	1	6/18/2008 2:07:00 PM
Freon 11		0.44	Ú.	0.15		ppbV	1	6/18/2008 2:07:00 PM
Freon 113		0.88		0.15		ppbV	1	6/18/2008 2:07:00 PM
Freon 114		ND	J	0.15		ppbV	1	6/18/2008 2:07:00 PM
Freon 12		0.68	V.	0.15		ppbV	1	6/18/2008 2:07:00 PM
Heptane		21		1.5		ppbV	10	6/19/2008 4:51:00 AM
Hexachloro-1,3	-butadiene	ND	\mathcal{O}	0.15		ppbV	1	6/18/2008 2:07:00 PM
Hexane		32		1.5		ppbV	10	6/19/2008 4:51:00 AM
Isopropyl alcoh	ol	ND	\mathcal{O}	0.15		ppbV	1	6/18/2008 2:07:00 PM
m&p-Xylene		4.1		0.30		ppbV	1	6/18/2008 2:07:00 PM
Methyl Butyl Ke	etone	ND	\mathcal{O}	0.30		ppbV	1	6/18/2008 2:07:00 PM
Methyl Ethyl Ke	etone	27		3.0		ppbV	10	6/19/2008 4:51:00 AM
Methyl Isobutyl	Ketone	19		3.0		ppbV	10	6/19/2008 4:51:00 AM
Methyl tert-buly	/l ether	ND	U	0.15		ppbV	1	6/18/2008 2:07:00 PM
Methylene chlo	ride	ND	\cup	0.15		ppbV	1	6/18/2008 2:07:00 PM
o-Xylene		1.4		0.15		ppbV	1	6/18/2008 2:07:00 PM
Propylene		ND	\mathcal{O}	0.15		ppbV	1	6/18/2008 2:07:00 PM
Styrene		0.83		0.15		ppbV	1	6/18/2008 2:07:00 PM
Tetrachloroethy	ylene	1,9		1.5		ppbV	10	6/19/2008 4:51:00 AM
Tetrahydrofurai	n	ND	J	0.15		ppbV	1	6/18/2008 2:07:00 PM
Toluene	19	7.8		1.5		ppbV	10	6/19/2008 4:51:00 AM
trans-1,2-Dichle	oroethene	ND	\mathcal{O}	0.15		ppbV	1	6/18/2008 2:07:00 PM
trans-1,3-Dichle		ND	\cup	0.15		ppbV	1	6/18/2008 2:07:00 PM
Trichloroethene	9	360		240		ppbV	1620	6/19/2008 10:17:00 PM
Vinyl acetate		1.9		1.5		ppbV	10	6/19/2008 4:51:00 AM
Vinyl Bromide		ND	U	0.15		ppbV	1	6/18/2008 2:07:00 PM
Vinyl chloride		ND	Ū.	0.15		ppbV	1	6/18/2008 2:07:00 PM
,	fluorobenzene	118		70-130		%REC	1	6/18/2008 2:07:00 PM

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Qualifiers:

- B Analyte detected in the associated Method Blank
- H Holding times for preparation or analysis exceeded
- JN Non-routine analyte. Quantitation estimated.
- Spike Recovery outside accepted recovery limits S
- E Value above quantitation range
- Analyte detected at or below quantitation limits J
- ND Not Detected at the Reporting Limit

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1122/08 OC

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Date: 20-Jun-08

CLIENT:	Earth Tech			4	Client Sample ID:		
Lab Order:	C0806018				Tag Number:		
Project:	AFB 59 (BAE)				Collection Date:	6/13/20	800
Lab ID:	C0806018-005A				Matrix:	AIR	
Analyses		Result		Limit Qua	l Units	DF	Date Analyzed
FIELD PARAMI				FLD			Analyst:
Vacuum Readir	ng "Hg	-2			"Hg		6/13/2008
HELIUM LEAK	TEST			GC			Analyst: RJP
Helium		ND		1.0	%	1	6/19/2008
1UG/M3 BY ME	ETHOD TO15			TO-15			Analyst: RJP
1,1,1-Trichloroe	ethane	20000		2200	ppbV	14580	6/20/2008 6:50:00 AM
1,1,2,2-Tetrachi	ioroethane	ND	U	0.15	Vdqq	1	6/18/2008 2:43:00 PM
1,1,2-Trichloroe	elhane	3.6		1.5	ppbV	10	6/19/2008 6:03:00 AM
1,1-Dichloroeth	ane	1200		240	Vdqq	1620	6/19/2008 10:51:00 PM
1,1-Dichloroeth	ene	51		1.5	ppbV	10	6/19/2008 6:03:00 AM
1,2,4-Trichlorob	benzene	ND	$\mathcal{O}_{\mathbb{C}}$	0.15	ppbV	1	6/18/2008 2:43:00 PM
1,2,4-Trimethyll	benzene	2.6		1.5	ppbV	10	6/19/2008 6:03:00 AM
1,2-Dibromoeth	ane	ND	U	0.15	ppbV	1	6/18/2008 2:43:00 PM
1,2-Dichlorober	nzene	ND	\cup	0.15	ppbV	1	6/18/2008 2:43:00 PM
1,2-Dichloroeth	ane	2300		240	ppbV	1620	6/19/2008 10:51:00 PM
1,2-Dichloropro	pane	ND	\mathcal{V}	0.15	ppbV	1	6/18/2008 2:43:00 PM
1,3.5-Trimethyll	benzene	2.9		1.5	ppbV	10	6/19/2008 6:03:00 AM
1,3-butadiene		ND	0	0.15	ppbV	1	6/18/2008 2:43:00 PM
1,3-Dichlorober	nzene	ND	\mathcal{V}	0.15	ppbV	1	6/18/2008 2:43:00 PM
1,4-Dichlorober	nzene	1.9	,	0.15	ppbV	1	6/18/2008 2:43:00 PM
1,4-Dioxane		ND	U	0.30	ppbV	1	6/18/2008 2:43:00 PM
2,2,4-trimethylp	pentane	2.7		1.5	ppbV	10	6/19/2008 6:03:00 AM
4-ethyltoluene		0.85		0.15	ppbV	1	6/18/2008 2:43:00 PM
Acetone		420		490 J	ppbV	1620	6/19/2008 10:51:00 PM
Allyl chloride		ND	\mathcal{O}	0.15	ppbV	1	6/18/2008 2:43:00 PM
Benzene		4.9		1.5	ppbV	10	6/19/2008 6:03:00 AM
Benzyl chloride	2	ND	U_{-}	0.15	ppbV	1	6/18/2008 2:43:00 PM
Bromodichioror	methane	ND	1	0.15	ppbV	1	6/18/2008 2:43:00 PM
Bromoform		ND		0.15	ppbV	1	6/18/2008 2:43:00 PM
Bromomethane	?	ND	V	0.15	ppbV	1	6/18/2008 2:43:00 PM
Carbon disulfid	e	4.0		1.5	ppbV	10	6/19/2008 6:03:00 AM
Carbon tetrach	loride	1.4		0.15	Vdqq	1	6/18/2008 2:43:00 PM
Chlorobenzene	2	ND	\mathcal{O}	0.15	ppbV	1	6/18/2008 2:43:00 PM
Chloroethane		ND	\cup	0.15	ppbV	1	6/18/2008 2:43:00 PM
Chloroform		47		1.5	ppbV	10	6/19/2008 6:03:00 AM
Chloromethane	2	ND	\cup	0.15	ppbV	1	6/18/2008 2:43:00 PM
cis-1,2-Dichloro	pethene	760		240	ppbV	1620	6/19/2008 10:51:00 PM
cis-1,3-Dichloro	opropene	ND	\mathcal{O}	0.15	ppbV	1	6/18/2008 2:43:00 PM
Cyclohexane		1.6		0.15	ppbV	1	6/18/2008 2:43:00 PM
Dibromochloro	methane	ND	\mathbf{V}	0.15	ppbV	1	6/18/2008 2:43:00 PM

Qualifiers: B Analyte detected in the associated Method Blank

П

E Value above quantitation range

ND Not Detected at the Reporting Limit

J Analyte detected at or below quantitation limits

JN Non-routine analyte, Quantitation estimated,

Holding times for preparation or analysis exceeded

S — Spike Recovery outside accepted recovery limits

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7/22/08

Date: 20-Jun-08

CLIENT:	Earth Tech			С	lient Sample II	D: SG37-0	61308
Lab Order:	C0806018				Tag Numbe	r: 359, 14	4
Project:	AFB 59 (BAE)				Collection Dat	e: 6/13/20	08
Lab ID:	C0806018-005A				Matri	x: AIR	
Analyses		Result	Limit	Qual	Units	DF	Date Analyzed
IUG/M3 BY ME	ETHOD TO15		TO	-15			Analyst: RJF
Ethyl acetate		ND \vee	0.25		ppbV	1	6/18/2008 2:43:00 PM
Ethylbenzene		2.7	1.5		ppbV	10	6/19/2008 6:03:00 AM
Freon 11		0.47 💙	0.15		ppbV	1	6/18/2008 2:43:00 PM
Freon 113		7.8	1.5		Vđqq	10	6/19/2008 6:03:00 AM
Freon 114		ND V	0.15		ppbV	1	6/18/2008 2:43:00 PM
Freon 12		0.69 V	0.15		ppbV	1	6/18/2008 2:43:00 PM
Heptane		8.5	1.5		ppbV	10	6/19/2008 6:03:00 AM
Hexachloro-1,3	3-butadiene	ND U	0.15		vdqq	1	6/18/2008 2:43:00 PM
Hexane		17	1.5		ppbV	10	6/19/2008 6:03:00 AM
isopropyl alcoh	10]	ND V	0.15		ppbV	1	6/18/2008 2:43:00 PM
m&p-Xylene		7.5	3.0		ppbV	10	6/19/2008 6:03:00 AN
Methyl Butyl Ke	etone	ND V	0.30		ppbV	1	6/18/2008 2:43:00 PN
Methyl Ethyl Ke	etone	13	3.0		ppbV	10	6/19/2008 6:03:00 AM
Methyl Isobutyl	Ketone	ND U	0.30		Vdqq	1	6/18/2008 2:43:00 PM
Methyl tert-buty	yl ether	ND U	0.15		ppbV	1	6/18/2008 2:43:00 PM
Methylene chlo	oride	0.44 V	0.15		ppbV	1	6/18/2008 2:43:00 PM
o-Xylene		4.4	1.5		ppbV	10	6/19/2008 6:03:00 AM
Propylene		ND U	0.15		ppbV	1	6/18/2008 2:43:00 PM
Styrene		ND · J	0.15		ppbV	1	6/18/2008 2:43:00 PM
Tetrachloroethy	ylene	5.5	1.5		ppbV	10	6/19/2008 6:03:00 AN
Tetrahydrofura	n	ND U	0.15		ppbV	1	6/18/2008 2:43:00 PN
Toluene		7.0	1.5		ppbV	10	6/19/2008 6:03:00 AM
trans-1,2-Dichle	oroethene	20	1.5		ppbV	10	6/19/2008 6:03:00 AN
trans-1,3-Dichl	oropropene	ND V	0.15		ppbV	1	6/18/2008 2:43:00 PN
Trichloroethene	e	7700	2200		Vdqq	14580	6/20/2008 6:50:00 AM
Vinyl acetate		1.4	0.15		ppbV	1	6/18/2008 2:43:00 PM
Vinyl Bromide		ND U	0.15		ppbV	1	6/18/2008 2:43:00 PM
Vinyl chloride		ND U	0.15		ppbV	1	6/18/2008 2:43:00 PM
Surr: Bromo	fluorobenzene	158	70-130	s	%REC	1	6/18/2008 2:43:00 PN
Surr: Bromo	fluorobenzene	131	70-130	S	%REC	10	6/19/2008 6:03:00 AN
Surr: Bromo	fluorobenzene	84.0	70-130		%REC	14580	6/20/2008 6:50:00 AM
Surr: Bromo	fluorobenzene	92.0	70-130		%REC	1620	6/19/2008 10:51:00 P
Surr: Bromo	fluorobenzene	84.0	70-130		%REC	6240	6/20/2008 12:37:00 A

Qualifiers:

- B Analyte detected in the associated Method Blank
- 11 Holding times for preparation or analysis exceeded
- E Value above quantitation range
 - Analyte detected at or below quantitation limits J
 - ND Not Detected at the Reporting Limit
- JN Non-routine analyte. Quantitation estimated. Spike Recovery outside accepted recovery limits S

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1122/08

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Centek Laboratories, LLC

Date: 20-Jun-08

CLIENT:	Earth Tech				C	lient Sample ID:		
Lab Order:	C0806018					Tag Number:	238, 2	62
Project:	AFB 59 (BAE)					Collection Date:	6/13/2	800
Lab ID: C0806018-006A					AIR			
Analyses		Result		Limit ()ual	Units	DF	Date Analyzed
FIELD PARAM	ETERS			FLC)			Analyst:
Vacuum Readir	ng "Hg	-3				"Hg		6/13/2008
HELIUM LEAK	TEST			GC				Analyst: RJI
Helium		ND		1.0		%	1	6/19/2008
1UG/M3 BY ME	ETHOD TO15			TO-1	5			Analyst: RJI
1,1,1-Trichloroe	ethane	2.0	<u>)</u>	0.15		ppb∨	1	6/18/2008 3:18:00 PM
1,1,2,2-Tetrach	loroethane	ND	\cup	0.15		ppbV	1	6/18/2008 3:18:00 PM
1,1,2-Trichloroe	ethane	ND	i	0.15		ppbV	1	6/18/2008 3:18:00 PM
1,1-Dichloroeth	ane	ND]	0.15		ppbV	1	6/18/2008 3:18:00 PM
1,1-Dichloroeth	ene	NĎ	1	0.15		ppbV	1	6/18/2008 3:18:00 PM
1,2,4-Trichlorot	penzene	ND	\checkmark	0.15		ppbV	1	6/18/2008 3:18:00 PM
1,2,4-Trimethyl	benzene	0.38	\mathcal{O}	0.15		ppbV	1	6/18/2008 3:18:00 PM
1,2-Dibromoeth	ane	ND	\cup	0.15		ppbV	1	6/18/2008 3:18:00 PM
1,2-Dichlorober	nzene	ND	i	0.15		ppbV	1	6/18/2008 3:18:00 PM
1,2-Dichloroeth	ane	ND		0.15		ppbV	1	6/18/2008 3:18:00 PM
1,2-Dichloropro	ppane	ND	V	0.15		ppbV	1	6/18/2008 3:18:00 PN
1,3,5-Trimethyl	benzene	0.45		0.15		ppbV	1	6/18/2008 3:18:00 PN
1,3-butadiene		ND	Ų.	0.15		ppbV	1	6/18/2008 3:18:00 PM
1,3-Dichlorober	nzene	ND	\cup	0.15		ppbV	1	6/18/2008 3:18:00 PN
1,4-Dichlorober	nzene	0.73	U	0.15		ppbV	1	6/18/2008 3:18:00 PM
1,4-Dioxane		ND	\cup	0.30		ppbV	1	6/18/2008 3:18:00 PM
2,2,4-trimethylp	pentane	ND	\mathcal{O}	0.15		ppbV	1	6/18/2008 3:18:00 PM
4-ethyltoluene		0.11		0.15	J	ppbV	1	6/18/2008 3:18:00 PM
Acetone		16	\mathcal{O}	3.0		ppbV	10	6/19/2008 7:15:00 AM
Allyl chloride		ND	V	0.15		ppbV	1	6/18/2008 3:18:00 PM
Benzene		2.1	5	0.15		ppbV	1	6/18/2008 3:18:00 PN
Benzyl chloride	;	ND	\cup	0.15		ppbV	1	6/18/2008 3:18:00 PM
Bromodichloro	methane	ND	ł	0.15		ppbV	1	6/18/2008 3:18:00 PM
Bromoform		ND		0.15		opbV	1	6/18/2008 3:18:00 PM
Bromomethane	\$	NÐ	\mathbf{V}	0.15		ppbV	1	6/18/2008 3:18:00 PM
Carbon disulfid	ie	0.44		0.15		ppbV	1	6/18/2008 3:18:00 PM
Carbon tetrach	loride	ND	U	0.15		ppbV	1	6/18/2008 3:18:00 PN
Chlorobenzene	2	ND	Ì	0.15		ppb∨	1	6/18/2008 3:18:00 PN
Chloroethane		ND		0.15		ppbV	1	6/18/2008 3:18:00 PN
Chloroform		ND		0.15		ppbV	1	6/18/2008 3:18:00 PM
Chloromethane	9	ND	V	0.15		ppbV	1	6/18/2008 3:18:00 PM
cis-1,2-Dichloro	pethene	0.25		0.15		ppbV	1	6/18/2008 3:18:00 PM
cis-1,3-Dichloro	opropene	ND	\overline{U}	0.15		ppbV	1	6/18/2008 3:18:00 PM
Cyclohexane		4.2		1.5		ppbV	10	6/19/2008 7:15:00 AN
Dibromochloro	methane	ND	V	0.15		ppbV	1	6/18/2008 3:18:00 PM

Qualifiers:

JN

B Analyte detected in the associated Method Blank
 H Holding times for preparation or analysis exceeded

E Value above quantitation range

ND Not Detected at the Reporting Limit

J Analyte detected at or below quantitation limits

Non-routine analyte. Quantitation estimated.

S — Spike Recovery outside accepted recovery limits

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Date: 20-Jun-08

CLIENT:	Earth Tech				(Client Sample 1D:	SG38-	061308
Lab Order:	C0806018					Tag Number:	238, 2	62
Project:	AFB 59 (BAE)					Collection Date:	6/13/2	008
Lab ID:	C0806018-006A					Matrix:	AIR	
Analyses		Result		Limit	Qual	Units	DF	Date Analyzed
1UG/M3 BY ME	ETHOD TO15			тс)-15			Analyst: RJF
Ethyl acetate		ND	U	0.25		ppbV	1	6/18/2008 3:18:00 PM
Ethylbenzene		0.55		0.15		ppbV	1	6/18/2008 3:18:00 PM
Freon 11		0.26	\mathcal{O}	0.15		ppbV	1	6/18/2008 3:18:00 PM
Freon 113		ND	\mathcal{O}	0.15		ppbV	1	6/18/2008 3:18:00 PM
Freon 114		ND	\cup	0.15		ppbV	1	6/18/2008 3:18:00 PM
Freon 12		0.45	U	0.15		ppbV	1	6/18/2008 3:18:00 PM
Heptane		3.0	J	1.5		ppbV	10	6/19/2008 7:15:00 AM
Hexachloro-1,3	-butadiene	ND	\cup	0.15		ppbV	1	6/18/2008 3:18:00 PM
Hexane		3.9	Ĵ	1.5		ppbV	10	6/19/2008 7:15:00 AM
Isopropyl alcoh	ol	ND	\cup	0.15		ppbV	1	6/18/2008 3:18:00 PM
m&p-Xylene		1.4		0.30		ppbV	1	6/18/2008 3:18:00 PM
Methyl Butyl Ke	etone	ND	\vee	0.30		ppbV	1	6/18/2008 3:18:00 PM
Methyl Ethyl Ke	etone	2.1	\cup	0.30		ppbV	1	6/18/2008 3:18:00 PM
Methyl Isobutyl	Ketone	3.0	5	3.0		ppbV	10	6/19/2008 7:15:00 AM
Methyl tert-buty	l ether	ND	U	0.15		ppbV	1	6/18/2008 3:18:00 PM
Methylene chlo	ride	0.32	\cup	0.15		ppbV	1	6/18/2008 3:18:00 PM
o-Xylene		0.43		0.15		ppbV	1	6/18/2008 3:18:00 PM
Propylene		ND	\cup	0.15		ppbV	1	6/18/2008 3:18:00 PM
Styrene		ND	\cup	0.15		ppbV	1	6/18/2008 3:18:00 PM
Tetrachloroethy	lene	1.8		0.15		ppbV	1	6/18/2008 3:18:00 PM
Tetrahydrofurai	n	ND	\cup	0.15		ppbV	1	6/18/2008 3:18:00 PM
Toluene		2.0		1.5		ppbV	10	6/19/2008 7:15:00 AM
trans-1,2-Dichle	oroethene	ND	\mathcal{V}	0.15		ppbV	1	6/18/2008 3:18:00 PM
trans-1,3-Dichle		ND	\cup	0.15		ppbV	1	6/18/2008 3:18:00 PM
Trichloroethene	· ·	2.1		0.15		ppbV	1	6/18/2008 3:18:00 PM
Vinyl acetate		ND	\cup	0.15		ppbV	1	6/18/2008 3:18:00 PM
Vinyl Bromide		ND	1	0.15		ppbV	1	6/18/2008 3:18:00 PM
Vinyl chloride		ND	Ĵ	0.15		ppbV	1	6/18/2008 3:18:00 PM
,	fluorobenzene	125	-	70-130		%REC	1	6/18/2008 3:18:00 PM

Qualifiers:

- B Analyte detected in the associated Method Blank
 - Holding times for preparation or analysis exceeded
- 11 JN Non-routine analyte. Quantitation estimated.
- Spike Recovery outside accepted recovery limits S
- E Value above quantitation range
- Analyte detected at or below quantitation limits J

ND Not Detected at the Reporting Limit

1/22/08

Date: 20-Jun-08

CLIENT: Lab Order:	Earth Tech C0806018			C	lient Sample ID: Tag Number:		
Project:	AFB 59 (BAE)				Collection Date:		
-					Matrix:		
Lab ID:	C0806018-008A						
Analyses		Result		Limit Qual	Units	DF	Date Analyzed
FIELD PARAMI				FLD			Analyst:
Vacuum Readir	ng "Hg	-3			"Нд		6/13/2008
HELIUM LEAK Helium	TEST	NĎ		GC 1.0	%	1	Analyst: RJP 6/19/2008
1UG/M3 BY ME	ETHOD TO15			TO-15			Analyst: RJP
1,1,1-Trichloroe		0.40	J	0.15	ppbV	1	6/18/2008 4:30:00 PM
1,1,2,2-Tetrach		ND	U	0.15	ppbV	1	6/18/2008 4:30:00 PM
1,1,2-Trichloroe		NÐ	1	0.15	ppbV	1	6/18/2008 4:30:00 PM
1,1-Dichloroeth		ND		0.15	ppbV	1	6/18/2008 4:30:00 PM
1,1-Dichloroeth		ND		0.15	ppbV	1	6/18/2008 4:30:00 PM
1,2,4-Trichlorob		ND	V	0.15	ppbV	1	6/18/2008 4:30:00 PM
1,2,4-Trimethyll		0.52	Ú.	0.15	ppbV	1	6/18/2008 4:30:00 PM
1,2-Dibromoeth		ND	Ù	0.15	ppbV	1	6/18/2008 4:30:00 PM
1,2-Dichlorober		ND	i	0.15	ppbV	1	6/18/2008 4:30:00 PM
1,2-Dichloroeth		ND		0.15	ppbV	1	6/18/2008 4:30:00 PM
1,2-Dichloropro		ND		0.15	ppbV	1	6/18/2008 4:30:00 PM
1,3,5-Trimethyli		0.50		0.15	ppbV	1	6/18/2008 4:30:00 PM
1,3-butadiene		ND	17	0.15	ppbV	1	6/18/2008 4:30:00 PM
1,3-Dichlorober	nzene	ND	V	0.15	ppbV	1	6/18/2008 4:30:00 PM
1,4-Dichlorober		1.2	\cup	0.15	ppbV	1	6/18/2008 4:30:00 PM
1,4-Dioxane		ND	\cup	0.30	ppbV	1	6/18/2008 4:30:00 PM
2,2,4-trimethylp	peniane	0.24		0.15	ppbV	1	6/18/2008 4:30:00 PM
4-ethyltoluene		0.18		0.15	ppbV	1	6/18/2008 4:30:00 PM
Acetone		20	\mathcal{O}	3.0	ppbV	10	6/19/2008 2:28:00 PM
Allyl chloride		ND	\cup	0.15	ppbV	1	6/18/2008 4:30:00 PM
Benzene		2.9	Ś	1.5	ppbV	10	6/19/2008 2:28:00 PM
Benzyl chloride	2	ND	Ŭ	0.15	ppbV	1	6/18/2008 4:30:00 PM
Bromodichloror		ND	i	0.15	ppbV	1	6/18/2008 4:30:00 PM
Bromoform		ND		0.15	ppbV	1	6/18/2008 4:30:00 PM
Bromomethane	2	ND	V	0.15	ppbV	1	6/18/2008 4:30:00 PM
Carbon disulfid		0.46	~	0.15	ppbV	1	6/18/2008 4:30:00 PM
Carbon tetrach		ND	\mathcal{O}	0.15	ppbV	1	6/18/2008 4:30:00 PM
Chlorobenzene		ND	4	0.15	ppbV	1	6/18/2008 4:30:00 PM
Chloroethane	-	ND	ļ	0.15	ppbV	1	6/18/2008 4:30:00 PM
Chloroform		ND		0.15	ppbV	1	6/18/2008 4:30:00 PM
Chloromethane	2	ND		0.15	ppbV	1	6/18/2008 4:30:00 PM
cis-1,2-Dichloro		0.19	Ť	0.15	ppbV	1	6/18/2008 4:30:00 PM
cis-1,3-Dichlord		ND	Ŭ	0.15	ppbV	1	6/18/2008 4:30:00 PM
	oproposio	11	v	1.5	ppb V ppbV	, 10	6/19/2008 2:28:00 PM
Cyclohexane	methane	ND	υ	0.15	ppbV ppbV	1	6/18/2008 4:30:00 PM

Qualifiers:

B Analyte detected in the associated Method Blank
 Holding times for preparation or analysis exceeded

E Value above quantitation range

J Analyte detected at or below quantitation limits

JN Non-routine analyte. Quantitation estimated.

S Spike Recovery outside accepted recovery limits

ND Not Detected at the Reporting Limit

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7/22/08

Date: 20-Jun-08

CLIENT:	Earth Tech				C	Client Sample ID:	SG38-	-061308DUP
Lab Order:	C0806018					Tag Number:	171, 2	.98
Project:	AFB 59 (BAE)					Collection Date:	6/13/2	2008
Lab ID:	C0806018-008A					Matrix:	AIR	
Analyses		Result				Units	DF	Date Analyzed
1UG/M3 BY ME	ETHOD TO15			тс	-15			Analyst: RJF
Ethyl acetate		ND	\vee	0.25		ppbV	1	6/18/2008 4:30:00 PM
Ethylbenzene		0.60		0.15		ppbV	1	6/18/2008 4:30:00 PM
Freon 11		0.27	\vee	0.15		ppbV	1	6/18/2008 4:30:00 PM
Freon 113		0.25	-	0.15		ppbV	1	6/18/2008 4:30:00 PM
Freon 114		ND	U	0.15		ppbV	1	6/18/2008 4:30:00 PM
Freon 12		0.47	\mathcal{V}	0.15		ppbV	1	6/18/2008 4:30:00 PM
Heptane		6.2	5	1.5		ppbV	10	6/19/2008 2:28:00 PM
Hexachloro-1,3	-butadiene	ND	\cup	0.15		ppbV	1	6/18/2008 4:30:00 PM
Hexane		8.1	J	1.5		ppbV	10	6/19/2008 2:28:00 PM
Isopropyl alcoh	ol	ND	Ũ	0.15		ppbV	1	6/18/2008 4:30:00 PM
m&p-Xylene		1.7		0.30		ppbV	1	6/18/2008 4:30:00 PM
Methyl Butyl Ke	etone	ND	U	0.30		ppbV	1	6/18/2008 4:30:00 PM
Methyl Ethyl Ke	etone	5.3	U	3.0		ppbV	10	6/19/2008 2:28:00 PM
Methyi isobutyi	Ketone	6.1	J	3.0		ppbV	10	6/19/2008 2:28:00 PM
Methyl tert-buty	/l ether	ND	U	0.15		ppb∨	1	6/18/2008 4:30:00 PM
Methylene chlo	ride	0.22	\cup	0.15		ppbV	1	6/18/2008 4:30:00 PM
o-Xylene		0.51		0.15		ppbV	1	6/18/2008 4:30:00 PM
Propylene		ND	U	0.15		ppbV	1	6/18/2008 4:30:00 PM
Styrene		0.35		0.15		ppbV	1	6/18/2008 4:30:00 PM
Tetrachloroethy	ylene	2.0		0.15		ppbV	1	6/18/2008 4:30:00 PM
Tetrahydrofura	n	ND	\mathcal{O}	0.15		Vdqq	1	6/18/2008 4:30:00 PM
Toluene		2.5		1.5		ppbV	10	6/19/2008 2:28:00 PM
trans-1,2-Dichl	oroelhene	ND	U	0.15		ppbV	1	6/18/2008 4:30:00 PM
trans-1,3-Dichl	oropropene	ND	\cup	0.15		ppbV	1	6/18/2008 4:30:00 PM
Trichloroethene	e	1.0		0.15		ppbV	1	6/18/2008 4:30:00 PM
Vinyl acetate		ND	\cup	0.15		ppbV	1	6/18/2008 4:30:00 PM
Vinyl Bromide		ND	1	0.15		ppbV	1	6/18/2008 4:30:00 PM
Vinyl chloride		ND	\mathcal{V}	0.15		ppbV	1	6/18/2008 4:30:00 PM
Surr: Bromo	fluorobenzene	94.0		70-130		%REC	1	6/18/2008 4:30:00 PM

Qualifiers:

- B Analyte detected in the associated Method Blank
- 11 Holding times for preparation or analysis exceeded
 - Non-routine analyte. Quantitation estimated. JN
 - S Spike Recovery outside accepted recovery limits
- E Value above quantitation range
-) Analyte detected at or below quantitation limits
- ND Not Detected at the Reporting Limit

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7/22/08

Date: 20-Jun-08

CLIENT:	Earth Tech			Client Sample ID:			
.ab Order:	C0806018			Tag Number:	202, 29)7	
Project:	AFB 59 (BAE)			Collection Date:	6/13/20	008	
Lab ID:	C0806018-007A			AIR	AIR		
Analyses		Result	Limit Qua	al Units	DF	Date Analyzed	
IELD PARAM	ETERS		FLD			Analyst:	
Vacuum Readir	ng "Hg	-2		"Hg		6/13/2008	
HELIUM LEAK	TEST		GC			Analyst: RJF	
Helium	1201	ND	1.0	%	1	6/19/2008	
UG/M3 BY ME	ETHOD TO15		TO-15			Analyst: RJF	
1,1,1-Trichloroe	ethane	0.35	0.15	ppbV	1	6/18/2008 3:54:00 PM	
1,1,2,2-Tetrach	loroethane	ND 🗸	0.15	ppbV	1	6/18/2008 3:54:00 PM	
1,1,2-Trichloroe	ethane	ND i	0.15	ppbV	1	6/18/2008 3:54:00 PM	
1,1-Dichloroeth	ane	ND	0.15	ppbV	1	6/18/2008 3:54:00 PM	
1,1-Dichloroeth	ene	ND	0.15	ppbV	1	6/18/2008 3:54:00 PM	
1,2,4-Trichlorot	penzene	ND 🕔	0.15	ppbV	1	6/18/2008 3:54:00 PM	
1,2,4-Trimethyl	benzene	4.2	1.5	ppbV	10	6/19/2008 1:16:00 PM	
1,2-Dibromoeth	nane	ND () 0.15	ppbV	1	6/18/2008 3:54:00 PM	
1,2-Dichlorober	nzene	ND	0.15	ppbV	1	6/18/2008 3:54:00 PM	
1,2-Dichloroeth	ane	ND	0.15	ppbV	1	6/18/2008 3:54:00 PM	
1,2-Dichloropro	pane	ND 👌	0.15	ppbV	1	6/18/2008 3:54:00 PM	
1,3,5-Trimethyl	benzene	2.9	1.5	ppbV	10	6/19/2008 1:16:00 PM	
1,3-butadiene		ND V	, 0110	ppbV	1	6/18/2008 3:54:00 PM	
1,3-Dichlorober	nzene	ND	0.15	ppb∨	1	6/18/2008 3:54:00 PM	
1,4-Dichlorobei	nzene	10	1.5	ppbV	10	6/19/2008 1:16:00 PM	
1,4-Dioxane		ND	0.30	ppbV	1	6/18/2008 3:54:00 PM	
2,2,4-trimethylp	pentane	ND' N	0.15	ppbV	1	6/18/2008 3:54:00 PM	
4-ethyltoluene		0.94	0.15	ppbV	1	6/18/2008 3:54:00 PM	
Acetone		980	320	ppbV	1080	6/19/2008 11:26:00 PI	
Allyl chloride		ND 🛝	J 0.15	ppbV	1	6/18/2008 3:54:00 PM	
Benzene		8.9	1.5	ppbV	10	6/19/2008 1:16:00 PM	
Benzyl chloride	9	ND () 0.15	ppbV	1	6/18/2008 3:54:00 PM	
Bromodichloro	methane	ND	0.15	ppbV	1	6/18/2008 3:54:00 PM	
Bromoform		ND	0.15	ppbV	1	6/18/2008 3:54:00 PM	
Bromomethane	j	ND	0.15	ppbV	1	6/18/2008 3:54:00 PM	
Carbon disulfid	le	6.0	1.5	ppbV	10	6/19/2008 1:16:00 PM	
Carbon tetrach	loride	ND \) 0.15	ppbV	1	6/18/2008 3:54:00 PM	
Chlorobenzene	9	ND	0.15	ppbV	1	6/18/2008 3:54:00 PM	
Chloroethane		ND	0.15	ppbV	1	6/18/2008 3:54:00 PM	
Chloroform		ND	0.15	ppbV	1	6/18/2008 3:54:00 PM	
Chloromethane	e	ND 🔻	2 0.15	ppbV	1	6/18/2008 3:54:00 PM	
cis-1,2-Dichlor	oethene	0.17	0.15	ppbV	1	6/18/2008 3:54:00 PM	
cis-1,3-Dichlor	opropene	ND V	J 0.15	ppbV	1	6/18/2008 3:54:00 PM	
Cyclohexane		ND	0.15	ppbV	1	6/18/2008 3:54:00 PM	
Dibromochloro	methane	ND	0.15	ppbV	1	6/18/2008 3:54:00 PM	

Qualifiers:

Н

JN

B Analyte detected in the associated Method Blank Holding times for preparation or analysis exceeded E Value above quantitation range

Analyte detected at or below quantitation limits J

ND Not Detected at the Reporting Limit

Non-routine analyte. Quantitation estimated. Spike Recovery outside accepted recovery limits S

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-122/02

Date: 20-Jun-08

CLIENT:	Earth Tech			Client	Sample ID:	SG39	-061308
Lab Order:	C0806018			т	ag Number:	202, 2	.97
Project:	AFB 59 (BAE)			Coll	ection Date:	6/13/2	2008
Lab ID:	C0806018-007A				Matrix:	AIR	
Analyses		Result	Limit	Qual Uni	ts	DF	Date Analyzed
1UG/M3 BY ME	ETHOD TO15	· · · · · · · · · · · · · · · · · · ·	тс)-15			Analyst: RJI
Ethyl acetate		ND V	0.25	ppb'	v	1	6/18/2008 3:54:00 PM
Ethylbenzene		1.4	0.15	ppb	V	1	6/18/2008 3:54:00 PM
Freon 11		0.31 - 1/*	0.15	ppb'	V	1	6/18/2008 3:54:00 PM
Freon 113		0.38	0.15	ppb	V	1	6/18/2008 3:54:00 PM
Freon 114		ND U	0.15	ppb	V	1	6/18/2008 3:54:00 PM
Freon 12		0.45 V	0.15	ppb	V	1	6/18/2008 3:54:00 PM
Heptane		40	6.0	ppb	V	40	6/19/2008 1:52:00 PM
Hexachloro-1,3	B-butadiene	ND U	0.15	ppb	V	1	6/18/2008 3:54:00 PM
Hexane		28	6.0	ppb	V	40	6/19/2008 1:52:00 PM
Isopropyl alcoh	ol	ND \lor	0.15	ppb	V	1	6/18/2008 3:54:00 PM
m&p-Xylene		3.6	0.30	ppb	V	1	6/18/2008 3:54:00 PM
Methyl Butyl Ke	elone	20	3.0	ppb	V	10	6/19/2008 1:16:00 PN
Methyl Ethyl Ke	elone	49	12	ppb	V	40	6/19/2008 1:52:00 PN
Methyl Isobutyl	Ketone	ND 🗸	0.30	ppb	V	1	6/18/2008 3:54:00 PN
Methyl tert-buty	/l ether	ND i	0.15	ppb	V	1	6/18/2008 3:54:00 PN
Methylene chlo	oride	ND 🗸	0.15	ppb	V	1	6/18/2008 3:54:00 PN
o-Xylene		1.5	0.15	ppb	V	1	6/18/2008 3:54:00 PM
Propylene		ND V	0.15	ppb	V	1	6/18/2008 3:54:00 PM
Styrene		ND V	0.15	ppb	v	1	6/18/2008 3:54:00 PM
Tetrachloroethy	ylene	1.9	0.15	ppb	V	1	6/18/2008 3:54:00 PN
Tetrahydrofura	•	ND V	0.15	dqq	V	1	6/18/2008 3:54:00 PM
Toluene		8.1	1.5	ppb	V	10	6/19/2008 1:16:00 PM
trans-1,2-Dichl	oroethene	ND V	0.15	ppb	V	1	6/18/2008 3:54:00 PN
trans-1,3-Dichl	oropropene	ND V	0.15	ppb	V	1	6/18/2008 3:54:00 PN
Trichloroethene	e	0.79 V	0.15	ppb	v	1	6/18/2008 3:54:00 PN
Vinyl acetate		ND V	0.15	ppb	V	1	6/18/2008 3:54:00 PN
Vinyl Bromide		ND	0.15	ppb	V	1	6/18/2008 3:54:00 PN
Vinyl chloride		ND 🗸	0.15	ppb	V	1	6/18/2008 3:54:00 PM
	fluorobenzene	114	70-130	%R	EC	1	6/18/2008 3:54:00 PN

Qualifiers:

- B Analyte detected in the associated Method Blank
- H Holding times for preparation or analysis exceeded
- Non-routine analyte. Quantitation estimated. JN
- S
 - Spike Recovery outside accepted recovery limits

E Value above quantitation range

Analyte detected at or below quantitation limits 3

ND Not Detected at the Reporting Limit

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7122/08

Date: 20-Jun-08

CLIENT: Lab Order: Project:	Earth Tech C0806018 AFB 59 (BAE)					lient Sampl Tag Nun Collection I	iber: Date:	290, 1 6/13/2	52 .	
Lab ID:	Matrix: AJR									
Analyses		Result		Limit	Qual	Units	<u> </u>	DF	Date Analyzed	
FIELD PARAM	ETERS			FL	.D				Analyst:	
Vacuum Readir	ng "Hg	-3				"Hg			6/13/2008	
1UG/M3 W/ 0.2	5UG/M3 CT-TCE-VC			то	-15				Analyst: RJP	
1,1,1-Trichloroe		ND	U	0.150		ppbV		1	6/18/2008 11:02:00 AM	
1,1,2,2-Tetrach	loroethane	ND	,	0.150		ppbV		1	6/18/2008 11:02:00 AM	
1,1,2-Trichloroe	ethane	ND		0.150		ppbV		1	6/18/2008 11:02:00 AM	
1,1-Dichloroeth	ane	ND	1	0.150		ppbV		1	6/18/2008 11:02:00 AM	
1,1-Dichloroeth	ene	ND		0.150		ppbV		1	6/18/2008 11:02:00 AM	
1,2,4-Trichlorob	benzene	ND	J.	0.150		ppbV		1	6/18/2008 11:02:00 AM	
1,2,4-Trimethyll	benzene	0.160		0.150		ppbV		1	6/18/2008 11:02:00 AM	
1,2-Dibromoeth	ane	ND	V	0.150		ppbV		1	6/18/2008 11:02:00 AM	
1,2-Dichlorober	nzene	ND	ı	0.150		ppbV		1	6/18/2008 11:02:00 AM	
1,2-Dichloroeth	ane	ND		0.150		ppbV		1	6/18/2008 11:02:00 AM	
1,2-Dichloropro	pane	ND		0.150		ppbV		1	6/18/2008 11:02:00 AM	
1,3,5-Trimethyll	benzene	ND		0.150		ppbV		1	6/18/2008 11:02:00 AM	
1,3-butadiene		ND		0.150		ppbV		1	6/18/2008 11:02:00 AM	
1,3-Dichlorober	nzene	ND	1	0.150		ppbV		1	6/18/2008 11:02:00 AM	
1,4-Dichlorober	nzene	0.240		0.150		ppbV		1	6/18/2008 11:02:00 AM	
1,4-Dioxane		ND	\mathcal{O}^{-}	0.300		ppbV		1	6/18/2008 11:02:00 AM	
2,2.4-trimethylp	pentane	ND	4	0.150		ppbV		1	6/18/2008 11:02:00 AM	
4-ethyltoluene		ND	J.	0.150		ppbV		1	6/18/2008 11:02:00 AM	
Acetone		31.0	Ŭ	6.00		ppbV		20	6/19/2008 1:21:00 AM	
Allyl chloride		ND	\mathcal{O}^{-}	0.150		ppbV		1	6/18/2008 11:02:00 AM	
Benzene		0.190		0.150		ppbV		1	6/18/2008 11:02:00 AM	
Benzyl chloride	9	ND	\cup	0.150		ppbV		1	6/18/2008 11:02:00 AM	
Bromodichloror	methane	ND	I	0.150		ppbV		1	6/18/2008 11:02:00 AM	
Bromoform		ND		0.150		ppbV		1	6/18/2008 11:02:00 AM	
Bromomethane	9	ND		0.150		ppbV		1	6/18/2008 11:02:00 AM	
Carbon disulfid	e	ND	ļ	0.150		ppbV		1	6/18/2008 11:02:00 AM	
Carbon tetrach	loride	ND	į	0.0400		ppbV		1	6/18/2008 11:02:00 AM	
Chlorobenzene	2	ND	-	0.150		ppbV		1	6/18/2008 11:02:00 AM	
Chloroethane		ND		0.150		ppbV		1	6/18/2008 11:02:00 AM	
Chloroform		ND	V	0.150		ppbV		1	6/18/2008 11:02:00 AM	
Chloromethane		0.480		0.150		ppbV		1	6/18/2008 11:02:00 AM	
cis-1.2-Dichloro	pethene	ND	Ú	0.150		ppbV		1	6/18/2008 11:02:00 AN	
cis-1,3-Dichloro	opropene	ND	i	0.150		ppbV		1	6/18/2008 11:02:00 AN	
Cyclohexane		ND		0.150		ppbV		1	6/18/2008 11:02:00 AN	
Dibromochloro	methane	ND	\sim	0.150		ppbV		1	6/18/2008 11:02:00 AN	
Ethyl acetate		0.150		0.250	J	ppbV		1	6/18/2008 11:02:00 AN	
Ethylbenzene		ND	\mathcal{V}^{-}	0.150		ppbV		1	6/18/2008 11:02:00 AM	

Qualifiers:

)-I

B Analyte detected in the associated Method Blank Holding times for preparation or analysis exceeded E Value above quantitation range

Analyte detected at or below quantitation limits j.

Non-routine analyte. Quantitation estimated. JN

Spike Recovery outside accepted recovery limits S

ND Not Detected at the Reporting Limit

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Date: 20-Jun-08

CLIENT:	Earth Tech	Client Sample ID:	AB-061308
Lab Order:	C0806018	Tag Number:	290, 152
Project:	AFB 59 (BAE)	Collection Date:	6/13/2008
Lab ID:	C0806018-009A	Matrix:	AIR

Analyses	Result		Limit	Qual	Units	DF	Đate Analyzed
1UG/M3 W/ 0.25UG/M3 CT-TCE-VC			то	-15			Analyst: RJP
Freon 11	0.280		0.150		ppb∨	1	6/18/2008 11:02:00 AM
Freon 113	ND	\cup	0.150		₽₽₽V	1	6/18/2008 11:02:00 AM
Freon 114	ND	\mathcal{O}	0.150		ppbV	1	6/18/2008 11:02:00 AM
Freon 12	0.590		0.150		ppbV	1	6/18/2008 11:02:00 AM
Heptane	NĎ	\cup	0.150		ppbV	1	6/18/2008 11:02:00 AM
Hexachloro-1,3-butadiene	ND	\cup	0.150		ppbV	1	6/18/2008 11:02:00 AM
Hexane	0.120		0.150	J	ppbV	1	6/18/2008 11:02:00 AM
Isopropyl alcohol	ND	U	0.150		ppbV	1	6/18/2008 11:02:00 AM
m&p-Xylene	0.170		0.300	J	ppb∨	1	6/18/2008 11:02:00 AM
Methyl Butyl Ketone	ND	V	0.300		ppbV	1	6/18/2008 11:02:00 AM
Methyl Ethyl Ketone	0.540		0.300		ppbV	1	6/18/2008 11:02:00 AM
Methyl Isobutyl Ketone	ND	υ	0.300		ppbV	1	6/18/2008 11:02:00 AM
Methyl tert-butyl ether	ND	U	0.150		ppbV	1	6/18/2008 11:02:00 AN
Methylene chloride	0.240		0.150		ppbV	1	6/18/2008 11:02:00 AN
o-Xylene	ND	J	0.150		Vđqq	1	6/18/2008 11:02:00 AN
Propylene	ND	i	0.150		Vđqq	1	6/18/2008 11:02:00 AN
Styrene	ND		0.150		ppbV	1	6/18/2008 11:02:00 AM
Tetrachloroethylene	NĎ		0.150		ppbV	1	6/18/2008 11:02:00 AM
Tetrahydrofuran	ND	\mathbf{V}	0.150		ppbV	1	6/18/2008 11:02:00 AM
Toluene	0.390		0.150		ppbV	1	6/18/2008 11:02:00 AN
trans-1,2-Dichloroethene	ND	\mathcal{O}	0.150		ppbV	1	6/18/2008 11:02:00 AM
trans-1,3-Dichloropropene	ND	\cup	0.150		ppbV	1	6/18/2008 11:02:00 AN
Trichloroethene	0.170		0.0400		ppbV	1	6/18/2008 11:02:00 AN
Vinyl acetate	ND	\cup	0.150		ppbV	1	6/18/2008 11:02:00 AM
Vinyl Bromide	ND	ì	0.150		ppbV	1	6/18/2008 11:02:00 AM
Vinyl chloride	ND	\mathbb{V}	0.0400		ppbV	1	6/18/2008 11:02:00 AN
Surr: Bromofluorobenzene	99.0		70-130		%REC	1	6/18/2008 11:02:00 AN

Qualifiers:

- В Analyte detected in the associated Method Blank
- Holding times for preparation or analysis exceeded
- JN Non-routine analyte. Quantitation estimated.
- S Spike Recovery outside accepted recovery limits

E Value above quantitation range

Analyte detected at or below quantitation limits Ļ

ND Not Detected at the Reporting Limit

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1.0 INTRODUCTION

This data quality review pertains to groundwater samples collected in June 2008 at Air Force Plant 59 (AFP-59). Parameters evaluated in groundwater samples included the total concentration of volatile organic constituent (VOC) and 1,4-dioxane. The samples were analyzed by Microbac Services, Marietta, Ohio.

Data quality review is an after-the-fact technical review of analytical data whereby the quality and usability of the data are determined based on a set of predefined criteria. These criteria depend upon the type of data involved and the purpose for which those data were collected. Data quality review assesses whether and to what extent specified criteria were met, and places restrictions on data use based on quality parameters. The data quality review process can range from a cursory review used to detect out-of-control situations to a detailed evaluation, depending on the analytical protocol, the associated quality control samples collected, and the intended data use.

Specific criteria for data quality review may include, but are not limited to: technical holding times, analysis of blanks, surrogate spike recovery, analysis of duplicates, and reported practical quantitation limits (PQLs). Where applicable, the recommendations of USEPA SW-846 Test Methods for Evaluating Solid Waste (Third Edition, December 1996) or USEPA Methods for Chemical Analysis of Water and Wastes (Revised March 1983) analytical method requirements, USEPA CLP National Functional Guidelines for Organic and Inorganic Data Review (February 1994, Functional Guidelines) data review guidance, and professional judgment.

Table 1 presents the data qualifiers applied during this review effort and their meanings.

Qualifier	Description
J	This is an estimated value.
UJ	The analyte was analyzed for but was not detected. The associated value is an estimate and may be inaccurate or imprecise.
R	The associated quality control indicates that the sample results are not useable.
U	The analyte was analyzed for, but not detected. The associated numerical value is at or below the MDL.

Table 1 Data Qualifiers

Table 2 provides a cross-reference list for field sample IDs and lab sample IDs.

Field Sample ID	Lab Sample ID	Field Sample ID	Lab Sample ID
59DW1WG1	L08060559-01	GS-95025WG-1	L08060559-10
TB061408	L08060559-02	GS-9505WG-1	L08060559-11
59SW1WG1	L08060559-03	59SW4WG1	L08060559-12
59SW3WG1	L08060559-04	59SW4WG9	L08060559-13
59DW3WG1	L08060559-05	G\$9506WG1	L08060559-14
59SW7WG1	L08060559-06	URS-2DWG1	L08060559-15
59SW7WG1-MS	L08060559-07	URS-2SWG1	L08060559-16
59SW7WG1-MSD	L08060559-08	URS-2SWG9	L08060559-17
GS-9502DWG-1	L08060559-09	EB061708	L08060559-18

 Table 2

 Field Sample ID/Lab Sample ID Cross Reference

During the data quality review process, laboratory qualified and unqualified data are verified against all available supporting documentation. Based on this review, qualifier codes may be added, deleted, or modified by the validator. Final results are therefore either qualified or unqualified. (Note: In those cases where the laboratory added a "U" flag indicated a non-detect result, and the validator agrees with this flag, then it remains intact, as noted on the corresponding Results Summary Form.) Changes to the data are reflected on the Results Summary Forms in Appendix A.

2.0 VOLATILE ORGANIC CONSTITUENTS

Volatile organic constituents were analyzed using EPA Test Method for Evaluating Solid Waste (SW-846) Method 8260B and SW8260B Selective Ion Monitoring (SIM). Samples were analyzed in batches:

<u>WG275043</u>: 59DW1WG1, GS-9502DWG-1, GS-95025WG-1, GS-9505WG-1, GS9506WG1, URS-2DWG1, URS-2SWG9, EB061708

<u>WG275878</u>: TB061408, 59SW1WG1, 59SW3WG1, 59DW3WG1, 59SW7WG1, 59SW7WG1-MS, 59SW7WG1-MSD, 59SW4WG1, 59SW4WG9

2.1 Holding Times

All samples were analyzed within prescribed hold times. No qualification is needed.

2.2 Calibration

<u>WG275043</u>: The Initial Calibration for this batch was performed on 06/24/2008. Standards were analyzed at 0.30, 0.4, 1, 2, 5, 20, 50, 100, and 200 ug/L in support of the Initial Calibration. The %RSD values were less than 30% for all target constituents. No qualification is needed.

The %D for the second source calibration was within $\pm 25\%$ for target constituents except for vinyl chloride, MTBE, and dichlorodifluoromethane. The validator qualifies UJ or J, the non detect and positive results, respectively, in the vinyl chloride, MTBE, and dichlorodifluoromethane results in the associated samples. The validator removes the "R" flag assigned by the laboratory to denote anomalies.

For the associated continuing calibration standard, none of the associated %D values were greater than 20% for target constituents. No qualification is needed.

<u>WG274878:</u> The Initial Calibration for this batch was performed on 06/29/2008. Standards were analyzed at 0.30, 0.4, 1, 2, 5, 20, 50, 100, and 200 ug/L in support of the Initial Calibration. The %RSD values were less than 30% for all target constituents. No qualification is needed.

The %D for the second source calibration was within $\pm 25\%$ for target constituents except for dichlorodifluoromethane. The validator qualifies UJ the dichlorodifluoromethane results in the associated samples. The validator removes the "R" flag assigned by the laboratory to denote anomalies.

For the associated continuing calibration standard, none of the associated %D values were greater than 20% for target constituents. No qualification is needed.

<u>WG275287</u>: The Initial Calibration for this batch was performed on 06/26/2008. Standards were analyzed at 0.30, 0.4, 1, 2, 5, 20, 50, and 100ug/L in support of the Initial Calibration. The %RSD values were less than 30% for all target constituents. No qualification is needed.

The %D for the second source calibration was within $\pm 25\%$ for target constituents.

For the associated continuing calibration standard, none of the associated %D values were greater than 20% for target constituents. No qualification is needed.

For all samples, it is noted that for those results which were less than the RL but greater than the MDL, the laboratory assigned an "F" flag, indicating an estimated value. Unless qualified otherwise, the validator removes the F flag and replaces it with the "J" qualifier, indicating an estimated value.

2.3 Laboratory Control Samples

<u>WG275043</u>: Laboratory control sample exhibited a 130% recovery for MTBE, which are above the LCS limits. Since MTBE was already qualified, no additional qualification is needed.

<u>WG274878:</u> Laboratory control sample exhibited a 130% recovery for MTBE, which are above the LCS limits and 59.3% recovery for 1,4-dioxane which are below the LCS limits. The validator qualifies J the MTBE positive results in the associated samples and qualifies R the 1,4-dioxane results in the associated samples.

<u>WG275287</u>: Constituent recoveries from the associated laboratory control samples were within control limits. No qualification is needed.

2.4 Blanks

<u>WG275043</u>: No constituents were detected in the associated method blank or equipment blank EB061708. No qualification needed.

<u>WG274878</u>: No constituents were detected in the associated method blank or trip blank TB061408. No qualification needed.

WG275287: No constituents were detected in the associated method blank or trip blank TB061408. No qualification needed.

2.5 Matrix Spike/Matrix Spike Duplicate

<u>WG275043</u>: Matrix Spike/Matrix Spike duplicate samples were not included in this sample delivery group. No qualification is needed.

<u>WG274878:</u> Sample 59SW7WG1 served as the MS/MSD. Recoveries of 1,4-dioxane was below the control limit in the MS and MSD samples but exceeded the %RPD. The validator has already qualified 1,4-dioxane result in 59SW7WG1 only. No additional qualification is necessary.

WG275287: Sample 59SW7WG1 served as the MS/MSD. Recoveries were within the control limits. No qualification is necessary.

2.6 Surrogate Recovery

<u>WG275043</u>: All surrogate recoveries were within control limits for all environmental and quality control samples. No qualification is needed.

<u>WG274878:</u> All surrogate recoveries were within control limits for all environmental and quality control samples. No qualification is needed.

<u>WG275287</u>: All surrogate recoveries were within control limits for all environmental and quality control samples. No qualification is needed.

2.7 Internal Standards

All internal standard area counts were within control limits for all samples. No qualification is needed based on the internal standard information provided.

2.8 Duplicates

Sample 59SW4 and URS-2D were analyzed in duplicate. One of two criteria was followed when evaluating field duplicates, depending on the amount detected. If the amount detected was greater than five times the reporting limit (RL), then the relative percent difference (RPD) should have been less than 25 percent. If the amount was less than five times the RL, then the difference between the duplicate and the sample concentrations should have been less than the RL. Agreement is excellent and no gualification is needed.

Analyte	Reporting Limit (RL)	59SW4WG1	59SW4WG9	Relative Percent Difference (RPD)
1,1,1-	1.0	2.98	2.87	3.8%
trichloroethane				
1,1-dichloroethane	1.0	1.51	1.48	2.0%
1,1-dichloroethene	1.0	0.751F	0.712F	5.3%
Cis-1,2- dichloroethene	1.0	4.35	4.13	5.2%
Trichloroethene	1.0	17.0	17.8	4.6
Tetrachloroethene	1.0	0.959F	0.965F	0.62%

Table 3: Duplicate Comparison (µg/L)

Analyte	Reporting Limit (RL)	URS-2DWG1	URS-2DWG9	Relative Percent Difference (RPD)
1,1,2,2- tetrachloroethane	1.0	2.20	2.25	2.2%
1,1-dichloroethane	1.0	0.569F	0.585F	2.8%
Chloroform	0.3	0.204F	0.197F	3.5%
Cis-1,2- dichloroethene	1.0	0.996F	0.966F	3.1%
Trichloroethene	1.0	2.19	2.22	1.4%

Analyte	Reporting Limit (RL)	598W4WG1	59SW4WG9	Relative Percent Difference (RPD)
1,4-dioxane	2.0	8.18	7.20	12.7 %

2.8 Summary

The data are acceptable with validator-assigned qualifiers.

Report Number: L08060559 Report Date :July 23, 2008

ample Number:L08060559-01 Client ID:59DW1WG1	PrePrep Hethod:NONE Prep Method:5030B		ument:HPMS14 Date:06/25/2001	3 17:54
Matrix:Water	Analytical Method:8260B		Date:06/24/200	
group Number: WG275043	Analyst:CMS		Date:06/25/200	
Collect Date:06/14/2008 12:20	Dilution:1		ID:14M06610	
Sample Tag:01	Units:ug/L			
10 - 10 - 10 - 10 - 10 - 10 - 10 - 10 -				
Analyte	CAS. Number Result	Qual	RL	MDL
1,1,1,2-Tetrachloroethane	; 630-20-6	U	0.500	0.250
1,1,1-Trichloroethane	: 71-55-6	U U	1.00	0.250
1,1,2,2-Tetrachloroethane	79-34-5	0	0.500	0.125
1,1,2-Trichloroethane	79-00-5	υ	1.00	0.125
1,1-Dichloroethane 1.1-Dichloroethene	75-34-3	υ	1.00	0.500
1,1-Dichloropropene	563-58-6	Ū	1.00	0.250
1,2,3-Trichlorobenzene	87-61-6	Ū	1.00	0.150
1,2,3-Trichloropropane	96-18-4	ŭ	1.00	0.500
1,2,4-Trichlorobenzene	120-82-1	Ψ	1.00	0.200
1,2,4-Trimethylbenzene	95-63-6	U	1.00	0.250
1,2-Dichloroethane	107-06-2	Ū	0.500	0.250
1,2-Dichlorobenzene	95-50-1	Ū	1.00	0.125
1,2-Dibromo-3-chloropropane	96-12-8	U	2.00	1.00
1,2-Dichloropropane	78-87-5	U	1.00	0.200
1,2-Dibromoethane	106-93-4	U	1.00	0.250
1,3,5-Trimethylbenzene	108-67-8	ប	1.00	0.250
1,3-Dichlorobenzene	541-73-1	ប	1.00	0.250
1,3-Dichloropropane	142-28-9	U	0.400	0.200
1,4-Dichlorobenzene	106-46-7	ប	0.500	0.125
1,4-Dioxane	123-91-1	U	100	50.0
1-Chlorohexane	544-10-5	U	1.00	0.125
2,2-Dichloropropane	594-20-7	U	1.00	0.250
2-Chlorotoluene	95-49-8	U	1.00	0.125
4-Chlorotoluene	106-43-4	U	1.00	0.250
Acetone	67-64-1	U U	10.0	2.50
Benzene	71-43-2	บ บ	0.400	0.125
Bromobenzene	108-86-1 74-97-5	U 1	1.00	0.200
Bromochloromethane Bromodichloromethane	75-27-4	0	0.500	0.250
Bromoform	75-25-2	υ	1.00	0.500
Bromomethane	74-83-9	Ū	3.00	0.500
Carbon tetrachloride	56-23-5	U	1.00	0.250
Chlorobenzene	108-90-7	σ	0.500	0.125
Chloroethane	75-00-3	σ	1.00	0.500
Chloroform	67-66-3	ប	0.300	0.125
Chloromethane	74-87-3	υ	1.00	0.250
cis-1,2-Dichloroethene	156-59-2	U	1.00	0.250
cis-1,3-Dichloropropene	10061-01-5	U	0.500	0.250
Dibromochloromethane	124-48-1	U	0.500	0.250
Dibromomethane	74-95-3	U	1.00	0,250
Dichlorodifluoromethane	75-71-8	-R- U	U 1.00	0.250
Ethylbenzene	100-41-4	U	1.00	0.250
Hexachlorobutadiene	87-68-3	U U	0.600	0.250
Isopropylbenzene	98-82-8	U U	1.00	0.250
Methylene chloride Methyl t-butyl ether (MTBE)	75-09-2 1634-04-4	0 .R~1	1.00	0.250
Methyl t-butyl ether (MTBE) MEK (2-Butanone)	1634-04-4 78-93-3	U	10.0	2.50
MER (2-Butanone) MIBK (methyl isobutyl ketone)	108-10-1	υ	10.0	2.50
n-Butylbenzene	108-10-1	υ	1.00	0.250
n-Propylbenzene	103-65-1	U U	1.00	0.125
m-,p-Xylene	136777-61-2	Ŭ	2.00	0.500
Naphthalene	91-20-3	บั	1.00	0.200
o-Xylene	95-47-6	Ū	1.00	0.250
p-Isopropyltoluene	99-87-6	U	1.00	0.250
sec-Butylbenzene	135-98-8	Ū	1.00	0.250
Styrene	100-42-5	U	1.00	0.125
Trichloroethene	79-01-6	U	1.00	0.250
tert-Butylbenzene	98-06-6	U	1.00	0.250

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1123/08

Report Number: L08060559 Report Date : July 23, 2008

Sample Number: L08060559-01	PrePrep Method:NONE	Instrument:HPMS1	4
Client ID:59DW1WG1	Prep Method:5030B	Prep Date:06/25	5/2008 17:54
Matrix:Water	Analytical Method:8260B	Cal Date:06/24	/2008 22:32
Workgroup Number: WG275043	Analvst:CMS	Run Date:06/25	5/2008 17:54
Collect Date:06/14/2008 12:20	Dilution:1	File ID:14M0661	0
Sample Tag:01	Units:ug/L		
haluta	CAS_Number Result	Oual RL	MDI.
Anaryce			
Tetrachloroethene	127-18-4	U 1.00	0.250

Toluene	108-88-3	-	U	1.00	0.250
trans-1,2-Dichloroethene	156-60-5		Ų	1.00	0.250
trans-1,3-Dichloropropene	10061-02-0	5	U	1.00	0.500
Trichlorofluoromethane	75-69-4		U	1.00	0.250
Vinyl chloride	75-01-4		<i>j</i> ∰]] 1.00	0.250
Surrogate	% Recovery	Lower	Upper	Qual	
Dibromofluoromethane	98.9	85	115		
1,2-Dichloroethane-d4	102	72	119		
Toluene-d8	103	81	120		
4-Bromofluorobenzene	102	76	119		

U Undetected; the analyte was analyzed for, but not detected. R Because of quality control deficiences for this analyte, this data may be rejected.

2 36 o£

Miletobal

Folky DC

Report Number: L08060559

Report Date :July 23, 2008

Sample Number:L08060559-05 Client ID:59DW3WG1	PrePrep Method:NO Prep Method:50		Instrume: Pren Dai	nt:HPMS8 te:06/24/2	009 15-00
Matrix:Water	Analytical Method:82				008 15:22
kgroup Number: WG274878	Analyst:PJ				008 15:22
Collect Date:06/15/2008 12:00	Dilution:1		File ID:		.000 13.22
Sample Tag:01	Units:ug	/ւ			
		,	•		
Ang Jub a	CAS. Number	Result	0.0.01	DI	MDI
Analyte 1,1,1,2-Tetrachloroethane	630-20-6	Result	Qual U	RL 0.500	MDL 0.250
1,1,1-Trichloroethane	71-55-6		บ	1.00	0.250
1,1,2,2-Tetrachloroethane	79-34-5		υ	0.500	0.125
1,1,2-Trichloroethane	79-00-5		U	1.00	0.250
1,1-Dichloroethane	75-34-3		U	1.00	0,125
1,1-Dichloroethene	75-35-4		U	1.00	0.500
1,1-Dichloropropene	563-58-6		U	1.00	0.250
1,2,3-Trichlorobenzene	87-61-6		a	1.00	0.150
1,2,3-Trichloropropane	96-18-4		U	1.00	0.500
1,2,4-Trichlorobenzene 1,2,4-Trimethylbenzene	120-82-1 95-63-6	· · · · · · · · · · · · · · · · · · ·	ប ប	1.00	0.200
1,2-Dichloroethane	107-06-2		U	0.500	0.250
1,2-Dichlorobenzene	95-50-1		U	1.00	0.125
1,2-Dibromo-3-chloropropane	96-12-8		Ū	2.00	1.00
1,2-Dichloropropane	78-87-5		U	1.00	0.200
1,2-Dibromoethane	105-93-4		U	1.00	0.250
1,3,5-Trimethylbenzene	108-67-8		υ	1.00	0.250
1,3-Dichlorobenzene	541-73-1		υ	1.00	0.250
1,3-Dichloropropane	142-28-9		U	0.400	0.200
1,4-Dichlorobenzene 1,4-Dioxane	106-46-7 123-91-1	· · . · · · · · · · · · · · · · · · · ·	U U	100	0.125 50.0
1-Chlorohexane	544-10-5	····	U K	1.00	0.125
2,2-Dichloropropane	594-20-7			1.00	0.250
2-Chlorotoluene	95-49-8		U	1.00	0.125
4-Chlorotoluene	106-43-4	• •••• • • • •••• ••• ••• •••	σ	1.00	0.250
Acetone	67-64-1		U	10.0	2.50
Benzene	71-43-2		U	0.400	0.125
Bromobenzene	108-86-1	·	U U	1.00	0.125
Bromochloromethane Bromodichloromethane	74-97-5 75-27-4		UUU	0.500	0.200
Bromoform	75-25-2		Ŭ	1.00	0.500
Bromomethane	74-83-9		υ	3.00	0.500
Carbon tetrachloride	56-23-5		υ	1.00	0.250
Chlorobenzene	108-90-7		U	0.500	0.125
Chloroethane	75-00-3		U	1.00	0.500
Chloroform	67-66-3		U	0.300	0.125
Chloromethane cis-1,2-Dichloroethene	74-87-3	72 1	U	1.00	0.250
cis-1, 2-Dichloropropene	10061-01-5	13.1	ម	0.500	0.250
Dibromochloromethane	124-48-1		Ű	0.500	0.250
Dibromomethane	74-95-3	• • • • • • • • • • • • • • • • • • •	υ·	1.00	0.250
Dichlorodifluoromethane	75-71-8		-R-U)	1.00	0.250
Ethylbenzene	100-41-4	:	U	1.00	0.250
Hexachlorobutadiene	87-68-3	:	U	0.600	0.250
Isopropylbenzene	98-82-8		ប ប	1.00	0.250
Methylene chloride Methyl t-butyl ether (MTBE)	75-09-2 1634-04-4		UUU	1.00	0.250
MEK (2-Butanone)	78-93-3	•••••••••••••••••••••••••••••••••••••••	υ	10.0	2.50
MIBK (methyl isobutyl ketone)	108-10-1		Ū	10.0	2.50
n-Butylbenzene	104-51-8	· · · · · · · · · · · · ·	U	1.00	0.250
n-Propylbenzene	103-65-1	· · · · · · · · · · · · · · · · · · ·	U	1.00	0.125
m-,p-Xylene	136777-61-2		U	2.00	0.500
Naphthalene	91-20-3	· ·	ប	1.00	0.200
o-Xylene	95-47-6		U	1.00	0.250
p-Isopropyltoluene	99-87-6		ช บ	1.00	0.250
sec-Butylbenzene Styrene	135-98-8 100-42-5		ប ប	1.00	0.125
Trichloroethene	79-01-6		U	1.00	0.125
tert-Butylbenzene	98-06-6		U	1.00	0.250

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7128108

Report Number: L08060559 Report Date : July 23, 2008

Sample Number:L08060559-05	PrePrep Method:NONE	Instrument: HPMS8
Client ID: 59DW3WG1	Prep Method:5030B	Prep Date:06/24/2008 15:22
Matrix:Water	Analytical Method:8260B	Cal Date:04/29/2008 15:06
Workgroup Number: WG274878	Analyst:FJB	Run Date:06/24/2008 15:22
Collect Date:06/15/2008 12:00	Dilution:1	File ID:8M345911
Sample Tag:01	Units:ug/L	

Analyte	CAS. Numbe	r	Result	Qual	RL	MDL
Tetrachloroethene	127-18-4			υ	1.00	0.250
Toluene	108-88-3			υ	1.00	0.250
trans-1,2-Dichloroethene	156-60-5			U	1.00	0.250
trans-1,3-Dichloropropene	10061-02-0	5		U	1.00	0.500
Trichlorofluoromethane	75-69-4			Ŭ	1.00	0.250
Vinyl chloride	75-01-4			U	1.00	0.250
Surrogate	% Recovery	Lower	Up	per C	ual	
Dibromofluoromethane	111	85	1	15		
1,2-Dichloroethane-d4	110	72	1	19		
Toluene-d8	93.5	81	1	20		
4-Bromofluorobenzene	95.7	76	1	19	:	

U Undetected; the analyte was analyzed for, but not detected.
R Because of quality control deficiences for this analyte, this data may be rejected.

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7128/08

Report Number: L08060559 Report Date : July 23, 2008

Sample Number:L08060559-05	PrePrep Method:NONE	Instrument:HPMS14
Client ID:59DW3WG1	Prep Method: 5030B	Prep Date:06/27/2008 12:37
Matrix:Water	Analytical Method:8260B	Cal Date:06/26/2008 20:24
Workgroup Number:WG275287	Analyst:CMS	Run Date:06/27/2008 12:37
Collect Date:06/15/2008 12:00	Dilution:1	File ID:14M06653
Sample Tag:01	Units:ug/L	

Analyte	CAS. Num	ber	Result	Qual	RL	MDL
1,4-Dioxane	123-91-	-1	14.3	l	2.00	1.00
Surrogate	% Recovery	Lower	Uppe	r Qual	1	
Dibromofluoromethane	118	54	138		į	
1,2-Dichloroethane-d4	117	51	135		}	

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1125/5°

Report Number: L08060559 Report Date :July 23, 2008

ample Number:L08060559-03 Client ID:59SWIWGL	PrePrep Method:N Prep Method:S	Instrument: HPMS8 Prep Date: 06/24/2008 14:15					
		Prep Method:5030B Prep Date:06/24/2008 14:17 nalytical Method:8260B Cal Date:04/29/2008 15:06					
Matrix:Water aroup Number:WG274878	Analytical Method:8 Analyst:B			Date:04/29/2 Date:06/24/2			
* *	•	and the second sec		D:8M345909	.005 14:1		
Collect Date: 06/14/2008 14:00	Dilution:1 Units:v		riie 1	1. 0H343303			
Sample Tag:01	0n) t S : <u>r</u>	ig7ь		•			
Analyte	CAS. Number	Result	Qual	RL	MDL		
1,1,1,2-Tetrachloroethane	630-20-6	3	υ	0.500	0.250		
L, 1, 1-Trichloroethane	71-55-6	······································	U	1.00	0.25		
L,1,2,2-Tetrachloroethane	79-34-5		U	0.500	0.12		
1,1,2-Trichloroethane	79-00-5		U	1.00	0.256		
L,1-Dichloroethane	75-34-3		υ.	1.00	0.12		
L,1-Dichloroethene	75-35-4		ប	1.00	0.500		
L, 1-Dichloropropene	563-58-6	n gana an Aran Ina Aran Ina Aran Ina Ar	υ	1.00	0.25		
1,2,3-Trichlorobenzene	87-61-6	2	υ	1.00	0.15		
L,2,3-Trichloropropane	96-18-4		υ	1.00	0.50		
1,2,4-Trichlorobenzene	120-82-1		U	1.00	0.20		
1,2,4-Trimethylbenzene	95-63-6	· ···· ···	U :	1.00	0.25		
1,2-Dichloroethane	107-06-2		U	0.500	0.25		
.2-Dichlorobenzene	95-50-1		U ;	1.00	0.12		
1,2-Dibromo-3-chloropropane	96-12-8	1	U	2.00	1.00		
1,2-Dichloropropane	78-87-5		U	1.00	0.20		
1,2-Dibromoethane	106-93-4	1	U	1.00	0.25		
1,3,5-Trimethylbenzene	108-67-8		U	1.00	0.25		
1,3-Dichlorobenzene	541-73-1		υ	1.00	0.25		
1,3-Dichloropropane	142-28-9		U	0.400	0.20		
1,4-Dichlorobenzene	106-46-7		U :	0.500	0.12		
l,4-Dioxane	123-91-1		_U. R	100	50.0		
1-Chlorohexane	544-10-5		υ·	1.00	0.12		
2,2-Dichloropropane	594-20-7		U	1.00	0.25		
2-Chlorotoluene	95-49-8		U	1.00	0.12		
4-Chlorotoluene	106-43-4	· · · · · · · · ·	U	1.00	0.25		
Acetone	67-64-1		U	10.0	2.50		
Benzene	71-43-2		U	0.400	0.12		
Bromobenzene	108-86-1	2	ប	1.00	0.12		
Bromochloromethane	74-97-5	. Lanca and second	U	1.00	0.20		
Bromodichloromethane	75-27-4		U	0.500	0.25		
Bromoform	75-25-2		U	1.00	0.50		
Bromomethane	74-83-9		<u> </u>	3.00	0.50		
Carbon tetrachloride	56-23-5			1.00	0.25		
Chlorobenzene	108-90-7	····	บ บ	0.500	0.12		
Chloroethane	75-00-3		U	1.00	0.50		
Chloroform	67-66-3		υ υ	0.300	0.12		
Chloromethane	74-87-3		U U	1.00	0.25		
cis-1,2-Dichloroethene	156-59-2 10061-01-5	.	U U	0.500	0.25		
cis-1,3-Dichloropropene Dibromochloromethane	10001-01-5		U	0.500	0.25		
Dibromochioromethane Dibromomethane	124-48-L 74-95-3		U U	1.00	0.25		
Dichlorodifluoromethane	71-93-3	. : :		3 1.00	0.25		
Ethylbenzene	100-41-4	······		1.00	0.25		
Hexachlorobutadiene	87-68-3		U	0.600	0.25		
Isopropylbenzene	98-82-8	····	U	1.00	0.25		
Methylene chloride	75-09-2	······ ····	U	1.00	0.25		
Methyl t-butyl ether (MTBE)	1634-04-4		υ	5.00	0.50		
MEK (2-Butanone)	78-93-3		U	10.0	2.50		
MIBK (methyl isobutyl ketone)	108-10-1		U	10.0	2.50		
n-Butylbenzene	104-51-8		U	1.00	0.25		
n-Propylbenzene	103-65-1		U .	1.00	0.12		
m-,p-Xylene	136777-61-2		U	2.00	0.50		
Naphthalene	91-20-3		U	1.00	0.20		
o-Xylene	95-47-6		U	1.00	0.25		
p-Isopropyltoluene	99-87-6	1	U	1.00	0.25		
sec-Butylbenzene	135-98-8		υ	1.00	0.25		
Styrene	100-42-5		U	1.00	0.12		
Trichloroethene	79-01-6		U	1.00	0.25		
tert-Butylbenzene	98-06-6		U	1.00	0.25		

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Report Number: L08060559 Report Date :July 23, 2008

Sample Number: L08060559-03	PrePrep Method:NONE	Instrument:HPMS8
Client ID:59SW1WG1	Prep Method: 5030B	Prep Date:06/24/2008 14:17
Matrix:Water	Analytical Method:8260B	Cal Date:04/29/2008 15:06
Workgroup Number:WG274878	Analyst:FJB	Run Date:06/24/2008 14:17
Collect Date:06/14/2008 14:00	Dilution:1	File ID:8M345909
Sample Tag:01	Units:ug/L	
Analyte	CAS. Number Result	Qual RL MDL
Tetrachloroethene	127-18-4	U 1.00 0.250

Tetrachloroethene	127-18-4		U	1.00	0.250
Toluene	108-88-3		ប	1.00	0.250
trans-1,2-Dichloroethene	156-60-5		1 ប	1.00	0.250
trans-1,3-Dichloropropene	10061-02-6		U	1.00	0.500
Trichlorofluoromethane	75-69-4		U	1.00	0.250
Vinyl chloride	75-01-4		: ប	1.00	0.250
Surrogate	% Recovery	Lower	Upper	Qual	
Dibromofluoromethane	108	85	115		
1.2-Dichloroethane-d4	107	72	119	-	
Toluene-d8	94 7	81	120		
4-Bromofluorobenzene	93.8	76	119		

U Undetected; the analyte was analyzed for, but not detected.
R Because of quality control deficiences for this analyte, this data may be rejected.

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Report Number: L08060559 Report Date : July 23, 2008

Sample Number:L08060559-03	PrePrep Method:NONE	Instrument: HPMS14
Client ID: 59SW1WG1	Prep Method: 5030B	Prep Date:06/27/2008 11:26
Matrix:Water	Analytical Method:8260B	Cal Date:06/26/2008 20:24
Workgroup Number:WG275287	Analyst: CMS	Run Date:06/27/2008 11:26
Collect Date:06/14/2008 14:00	Dilution:1	File ID:14M06651
Sample Tag:01	Units:ug/L	

Analyte		CAS. Numb	er	Result	Qual	RL	MDL
1,4-Dioxane		123-91-:	1		บ	2.00	1.00
Surrogate	÷ %	Recovery	Lower	Uppe	r Ç	Qual	
Dibromofluoromethane		117	54	138			
1.2-Dichloroethane-d4		117	51	135			

U Undetected; the analyte was analyzed for, but not detected.

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1/2 or

Report Number: L08060559

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Report Date : July 23, 2008

Methylene chloride Methylene chloride Methyl t-butyl ether (MTBE) MEK (2-Butanone) MIBK (methyl isobutyl ketone)

n-Butylbenzene

n-Propylbenzene

Naphthalene

o-Xylene

Styrene

m-,p-Xylene

p-Isopropyltoluene

sec-Butylbenzene

Trichloroethene

tert-Butylbenzene

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Sample Number: 1.08060559-04	PrePrep Method:N Prep Method:5	eres a second		nent:HPMS8 Date:06/24/	2008 14-50
Client ID: 59SW3WG1	Analytical Method:8			Date:04/29/	
Matrix:Water	-	A REAL PROPERTY AND A REAL PROPERTY A REAL PROPERTY A REAL PROPERTY AND A REAL PROPERTY A REAL PROPERTY AND A REAL PROPERTY A		Date:06/24/	
group Number: WG274878	Analyst:F Dilution:1			D:8M345910	2008 14:50
Collect Date:06/15/2008 09:30			. FILE I	0.00343910	
Sample Tag:01	Units:u	id\r			
Analyte	CAS, Number	Result	Qual	RL	MDL
1,1,1,2-Tetrachloroethane	630-20-6	Rebuxe	U U	0.500	0.250
1, 1, 1. Trichloroethane	71-55-6	0.661	P- 1	1.00	0.250
1,1,2,2-Tetrachloroethane	79-34-5			0.500	0.125
1,1.2-Trichloroethane	79-00-5			1.00	0.250
1,1-Dichloroethane	75-34-3	0.403	-1	1.00	0.125
1,1-Dichloroethene	75-35-4		u u v	1.00	0.500
1,1-Dichloropropene	563-58-6		U J	1.00	0.250
1,2,3-Trichlorobenzene	87-61-6		ů.	1.00	0.150
1,2,3-Trichloropropane	96-18-4	·· ••····· · · · · · · · · · · · · · ·	Ű	1.00	0.500
1,2,4-Trichlorobenzene	120-82-1	··· · ···· ··· ··· ··· ··· ·· ··		1.00	0.200
1,2,4-Trimethylbenzene	95-63-6		U	1.00	0.250
1,2-Dichloroethane	107-06-2		υ	0.500	0.250
1,2-Dichlorobenzene	95-50-1		U U	1.00	0.125
1,2-Dibromo-3-chloropropane	96-12-8			2.00	1.00
1, 2-Dichloropropane	78-87-5		- U	1.00	0.200
1, 2-Dichioropropane 1, 2-Dibromoethane	106-93-4		u u	1.00	0.250
1,3,5-Trimethylbenzene	108-67-8		Ű	1.00	0.250
1, 3-Dichlorobenzene	541-73-1		u	1.00	0.250
1, 3-Dichloropropane	142-28-9		<u> </u>	0,400	0.200
1, 4-Dichlorobenzene	105-46-7		บ	0.500	0.125
1,4-Dioxane	123-91-1	· · · [· · · · · · · · · · · · · · · · · · ·	- K	100	50.0
1-Chlorohexane	544-10-5		- U	1.00	0.125
2.2-Dichloropropane	594-20-7		Ŭ	1.00	0.250
2-Chlorotoluene	95-49-8	1911 1 1 1 1 1 1 1 1 1 1 1	U	1.00	0.125
4-Chlorotoluene	106-43-4	· · · · · · · · · · · · · · · · · · ·		1.00	0.250
Acetone	67-64-1		U	10.0	2.50
Benzene	71-43-2	··· •	U .	0.400	0.125
Bromobenzene	108-86-1		υ	1.00	0.125
Bromochloromethane	74-97-5	· • • · · · · · · · · · · · · · · · · ·	U U	1.00	0.200
Bromodichloromethane	75-27-4			0.500	0.250
Bromoform	75-25-2			1.00	0.500
Bromomethane	74-83-9		ប	3.00	0.500
Carbon tetrachloride	56-23-5		- ŭ	1.00	0.250
Chlorobenzene	108-90-7		n n	0.500	0.125
Chloroethane	75-00-3	·····		1.00	0.500
Chloroform	67-66-3		ŭ	0.300	0.125
Chloromethane	74-87-3	· · · · · · · · · · · · · · · · · · ·	U .	1.00	0.250
cis-1,2-Dichloroethene	156-59-2	1.45	· · · • • • • • • • • • • • • • • • • •	1.00	0.250
cis-1,3-Dichloropropene	10061-01-5		U U	0.500	0.250
Dibromochloromethane	124-48-1		. U ,	0.500	0.250
Dibromomethane	74-95-3	·····	U U	1.00	0.250
Dichlorodifluoromethane	75-71-8	· [· · · · · · · · · · · · · · · · · ·		1.00	0.250
Ethylbenzene	100-41-4		U U	1.00	0.250
Hexachlorobutadiene	87-68-3		υ	0.600	0.250
Teopropy) benzene	98-82-8		U U	1.00	0.250

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98-82-8

75-09-2

1634-04-4

78-93-3

104-51-8

103-65-1

136777-61-2

91-20-3

95-47-6

99-87-6

135-98-8

100-42-5

79-01-6

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Report Number: L08060559 Report Date : July 23, 2008

Sample Number:L08060559-04	PrePrep Method:NONE	Instrument: HPMS8
Client ID:59SW3WG1	Prep Method:5030B	Prep Date:06/24/2008 14:50
Matrix:Water	Analytical Method:8260B	Cal Date:04/29/2008 15:06
Workgroup Number: WG274878	Analyst:FJB	Run Date:06/24/2008 14:50
Collect Date: 06/15/2008 09:30	Dilution:1	File ID:8M345910
Sample Tag:01	Units:ug/L	
: Annlyten	CAS Number Result	Qual BL MDL

Analyte	CAS. Number	Res	ult Qua	L .	RL	MDL
Tetrachloroethene	127-18-4	1	ប		1.00	0.250
Toluene	108-88-3		U		1.00	0.250
trans-1,2-Dichloroethene	156-60-5		U		1.00	0.250
trans-1,3-Dichloropropene	10061-02-6	1	υ		1.00	0.500
Trichlorofluoromethane	75-69-4	l	U		1.00	0.250
Vinyl chloride	75-01-4		σ		1.00	0.250
Surrogate	% Recovery	Lower .	Upper	Qual	:	
Dibromofluoromethane	. 109	85	115			
1,2-Dichloroethane-d4	108	72	119			
Toluene-d8	94.3	81	120			
4-Bromofluorobenzene	94.1	76	119	1		

U Undetected; the analyte was analyzed for, but not detected.

P The analyte was positively identified, but the quantitation was below the RL.

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Report Number: L08060559 Report Date :July 23, 2008

Sample Number:L08060559-04	PrePrep Method:NONE	Instrument: HPMS14
Client ID:59SW3WG1	Prep Method:5030B	Prep Date:05/27/2008 12:01
Matrix:Water	Analytical Method:8260B	Cal Date:06/26/2008 20:24
Workgroup Number:WG275287	Analyst:CMS	Run Date:06/27/2008 12:01
Collect Date:06/15/2008 09:30	Dilution:1	File ID:14M06652
Sample Tag:01	Units:ug/L	

Analyte	CAS. Nur	nber	Result Qua	il RL	MDL
1,4-Dioxane	123-91	-1	U U	2.00	1.00
Surrogate	% Recovery	Lower	Upper	Qual	
Dibromofluoromethane	107	54	138		
1,2-Dichloroethane-d4	112	51	135		

U Undetected; the analyte was analyzed for, but not detected.

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1128/08

Report Number: L08060559

Report Date :July 23, 2008

ample Number:L08060559-12 Client ID:59SW4WG1	PrePrep Method:N Prep Method:S		Instrument:HPMS8 Prep Date:06/24/2008 18:24				
Matrix:Water	Analytical Method:8		Cal Date:04/29/2008 15:06				
group Number:WG274878	Analyst:F			Date:06/24/	and and and the second second second		
Collect Date:06/16/2008 10:50	Dilution:1		File ID:8M345916				
Sample Tag:01	Units:u						
Sample rag.vi		84.5 <u>5</u>					
Analyte	CAS. Number	Result	Qual	RL	MDL		
1,1,1,2-Tetrachloroethane	630-20-6		U	0.500	0.250		
1,1,1-Trichloroethane	71-55-6	2.98		1.00	0.250		
1,1,2,2-Tetrachloroethane	79-34-5		ប	0.500	0.125		
1,1,2-Trichloroethane	79-00-5	4	U	1.00	0.250		
1,1-Dichloroethane	75-34-3	1.51		1.00	0.125		
1,1-Dichloroethene	75-35-4	0.751	U	1.00	0.300		
1,1-Dichloropropene	563-58-6		σ	1.00	0.150		
1,2,3-Trichlorobenzene	87-61-6 96-18-4		U U	1.00	0.500		
1,2,3-Trichloropropane	120-82-1	·	ប	1.00	0.200		
1,2,4-Trichlorobenzene	95-63-6		UUU	1.00	0.250		
1,2,4-Trimethylbenzene 1,2-Dichloroethane	107-06-2		U U	0.500	0.250		
1,2-Dichloroethane 1,2-Dichlorobenzene	95-50-1		υ	1.00	0.125		
1,2-Dichiorobenzene 1,2-Dibromo-3-chloropropane	96-12-8		υ	2.00	1.00		
1,2-Dichloropropane	78-87-5		Ū	1.00	0.200		
1,2-Dibromoethane	106-93-4		υ	1.00	0.250		
1,3,5-Trimethylbenzene	108-67-8		U	1.00	0.250		
1,3-Dichlorobenzene	541-73-1		ប	1.00	0.250		
1,3-Dichloropropane	142-28-9		U	0.400	0.200		
1,4-Dichlorobenzene	106-46-7		ប	0.500	0.125		
1,4-Dioxane	123-91-1		-4- R	100	50.0		
1-Chlorohexane	544-10-5		U	1.00	0.125		
2,2-Dichloropropane	594-20-7		ប	1.00	0.250		
2-Chlorotoluene	95-49-8		U	1.00	0.125		
4-Chlorotoluene	106-43-4		U	1.00	0.250		
Acetone	67-64-1		σ	10.0	2.50		
Benzene	71-43-2		U	0.400	0.125		
Bromobenzene	108-86-1		U	1.00	0.125		
Bromochloromethane	74-97-5		U	1.00	0.200		
Bromodichloromethane	75-27-4		ប ្	0.500	0.250		
Bromoform	75-25-2		U	1.00	0.500		
Bromomethane	74-83-9 56-23-5		U U	1.00	0.250		
Carbon tetrachloride Chlorobenzene	108-90-7		U	0.500	0.125		
Chloroethane	75-00-3		υ.	1.00	0.500		
Chloroform	67-66-3		v	0.300	0.125		
Chloromethane	74-87-3		Ū	1.00	0.250		
cis-1,2-Dichloroethene	156-59-2	4.35		1.00	0.250		
cis-1,3-Dichloropropene	10061-01-5	· · · · · · · · · · · · · · · · · · ·	U	0.500	0.250		
Dibromochloromethane	124-48-1		ប	0.500	0.250		
Dibromomethane	74-95-3		ប	1.00	0.250		
Dichlorodifluoromethane	75-71-8		- R- (1) 1.00	0.250		
Ethylbenzene	100-41-4		U	1.00	0.250		
Hexachlorobutadiene	87-68-3		υ :	0.600	0.250		
Isopropylbenzene	98-82-8	·	ប	1.00	0.250		
Methylene chloride	75-09-2	· · · · · · · · · · · · · · · · · · ·	V	1.00	0.250		
Methyl t-butyl ether (MTBE)	1634-04-4		U	5.00	0.500		
MEK (2-Butanone)	78-93-3		U	10.0	2.50		
MIBR (methyl isobutyl ketone)	108-10-1		U	10.0	2.50		
n-Butylbenzene	104-51-8		U	1.00	0.250		
n-Propylbenzene	103-65-1	<u> </u>	ប ប	1.00	0.125		
m-,p-Xylene	136777-61-2		ប ប	1.00	0.200		
Naphthalene	91-20-3		U U	1.00	0.200		
o-Xylene	99-87-6	····	บ บ	1.00	0.250		
p-Isopropyltoluene	135-98-8		บ บ	1.00	0.250		
sec-Butylbenzene Styrene	135-58-8		UUU	1.00	0.125		
Trichloroethene	79-01-6	17.0		1.00	0.250		
tert-Butylbenzene	98-06-6		U	1.00	0.250		

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Report Number: L08060559

Report Date : July 23, 2008

Sample Number: L08060559-12 Client ID: 595W4WG1 Matrix: Water	PrePrep Method:NONE Prep Method:5030B Analytical Method:8260B	Instrument:HPMS8 Prep Date:06/24/2008 18:24 Cal Date:04/29/2008 15:06
₩orkgroup Number:₩G274878 Collect Date:06/16/2008 10:50 Sample Tag:01	Analyst:FJB Dilution:1 Units:ug/L	File ID:8M345916
Analyte	CAS. Number Result	Qual RL MDL

Tetrachloroethene	127-18-4		0.959	E •)	1.00		0.250
Toluene	108-88-3			U		1.00		0.250
trans-1,2-Dichloroethene	156-60-5			U	·	1.00	:	0.250
trans-1,3-Dichloropropene	10061-02-6			υ		1.00		0.500
Trichlorofluoromethane	75-69-4	1		U	i	1.90		0.250
Vinyl chloride	75-01-4			U		1.00	! .	0.250
Surrogate	% Recovery	Lower	Upper		Qual	1		
Dibromofluoromethane	108	85	115					
1,2-Dichloroethane-d4	107	72	119					
Toluene-d8	95.2	81	120					
4-Bromofluorobenzene	94.3	76	119			j		

U Undetected; the analyte was analyzed for, but not detected.

B Because of quality control deficiences for this analyte, this data may be rejected.
 F The analyte was positively identified, but the quantitation was below the RL.

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7128/08

Report Number: L08060559 Report Date :July 23, 2008

Sample Number:L08060559-12	PrePrep Method:NONE	Instrument: HPMS14
Client ID:59SW4WG1	Prep Method: 5030B	Prep Date:06/27/2008 13:12
Matrix:Water	Analytical Method:8260B	Cal Date:06/26/2008 20:24
Workgroup Number: WG275287	Analyst: CMS	Run Date:06/27/2008 13:12
Collect Date:06/16/2008 10:50	Dilution:1	File ID:14M06654
Sample Tag:01	Units:ug/L	

Analyte	CAS. Numb	er Res		Qual	RL	MDL
1,4-Dioxane	123-91-3		18		.00	1.00
Surrogate	% Recovery	Lower	Upper	Qual		
Dibromofluoromethane	110	54	138			
1,2-Dichloroethane-d4	112	51	135			
Anonimiento contrata antico antico antico antico antico a contrata antico antico antico antico antico antico a						

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1/2×103

Report Number: L08060559

o-Xylene p-Isopropyltoluene

sec-Butylbenzene

Trichloroethene

tert-Butylbenzene

Report Date :July 23, 2008

Sample Number: L08060559-13 Client ID: 595W4WG9	PrePrep Method:NC Prep Method:50	Instrument:HPMS8 Prep Date:06/24/2008 18:57					
Matrix:Water	Analytical Method:82			Date:04/29/2			
group Number:WG274878	Analytical Method: 52 Analyst: Fi		Run Date:06/24/2008 18:57				
Collect Date:06/16/2008 10:50	Dilution:1			D:8M345917	.000 10.57		
Sample Tag:01	Units:uc	т/т.		0.000333317	·····		
Sample ragion	onico, ag	, <u>-</u>					
Analyte	CAS. Number	Result	Qual	RL	MDL		
1,1,1,2-Tetrachloroethane	630-20-6		U	0.500	0.250		
1,1,1-Trichloroethane	71-55-6	2.87		1.00	0.250		
1,1,2,2-Tetrachloroethane	79-34-5		υ	0.500	0.125		
1,1,2-Trichloroethane	79-00-5		U	1.00	0.250		
1,1-Dichloroethane	75-34-3	1.48		1.00	0.125		
1,1-Dichloroethene	75-35-4	0.712	F }	1.00	0.500		
1,1-Dichloropropene	563-58-6	-	υ	1.00	0.250		
1,2,3-Trichlorobenzene	87-61-6		υ.	1.00	0.150		
1,2,3-Trichloropropane	96-18-4	1	ប	1.00	0.500		
1,2,4-Trichlorobenzene	120-82-1		υ	1.00	0.200		
1,2,4-Trimethylbenzene	95-63-6		U	1.00	0.250		
1,2-Dichloroethane	107-05-2		U	0.500	0.250		
1,2-Dichlorobenzene	95-50-1		υ	1.00	0.125		
1,2-Dibromo-3-chloropropane	96-12-8		U	2.00	1.00		
1,2-Dichloropropane	78-87-5		U	1.00	0.200		
1,2-Dibromoethane	106-93-4		U	1.00	0.250		
1,3,5-Trimethylbenzene	108-67-8		IJ	1.00	0.250		
1,3-Dichlorobenzene	541-73-1	1	U	1.00	0.250		
1,3-Dichloropropane	142-28-9		ប	0.400	0.200		
1,4-Dichlorobenzene	106-46-7	1	U	0.500	0.125		
1,4-Dioxane	123-91-1		#12	100	50.0		
1-Chlorohexane	544-10-5		U	1.00	0.125		
2,2-Dichloropropane	594-20-7		U	1.00	0.250		
2-Chlorotoluene	95-49-8	······	U	1.00	0.125		
4-Chlorotoluene	106-43-4	• .• · · · · · · · · · · · · · · · ·	υ	1.00	0.250		
Acetone	67-64-1		U	10.0	2.50		
Benzene	71-43-2	1	U	0.400	0.125		
Bromobenzene	108-86-1	1	U	1.00	0.125		
Bromochloromethane	74-97-5		U	1.00	0.200		
Bromodichloromethane	75-27-4		U	0.500	0.250		
Bromoform	75-25-2		U	1.00	0.500		
Bromomethane	74-83-9		U	3.00	0.500		
Carbon tetrachloride	56-23-5		UU	1.00	0.250		
Chlorobenzene	108-90-7		ບ ນ	0.500	0.125		
Chloroethane	75-00-3	4	U U	1.00	0.500		
Chloroform	67-66-3		UU	0.300	0.125		
Chloromethane	74-87-3		<u> </u>	1.00	0.250		
cis-1,2-Dichloroethene	156-59-2	4.13	u l	0.500	0.250		
cis-1,3-Dichloropropene	10061-01-5		ប	0.500	0.250		
Dibromochloromethane	124-48-1 74-95-3		υ. υ	1.00	0.250		
Dibromomethane	74-95-3			100 years and a second strategy and a 198 years	0.250		
Dichlorodifluoromethane	and a second second by the second			1.00	0.250		
Ethylbenzene	100-41-4 87-68-3		U	0.600	0.250		
Hexachlorobutadiene	and a second second second second second second second second		U U	1.00	0.250		
Isopropylbenzene	98-82-8		U	1.00	0.250		
Methylene chloride	75-09-2 1634-04-4		U	5.00	0.500		
Methyl t-butyl ether (MTBE)	- set and set of the set of th		U U	10.0	2.50		
MEK (2-Butanone)	78-93-3		U U	10.0	2.50		
MIBK (methyl isobutyl ketone)	108-10-1 104-51-8		U U	1.00	0.250		
n-Butylbenzene	104-51-8		U	1.00	0.125		
n-Propylbenzene	136777-61-2		U	2.00	0.500		
m-,p-Xylene	91-20-3		U U	1.00	0.200		
Naphthalene	91-20-3		ប	1.00	0.250		
o-Xylene	95-47-6		- U 17	1.00	0.250		

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99-87-6

135-98-8

100-42-5

79-01-6

98-06-6

Microbac

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7128108 pc

Report Number: L08060559 Report Date : July 23, 2008

Sample Number:L08060559-13	PrePrep	Method:NONE	Instrument: HPMS8	
Client ID:59SW4WG9	Prep	Method:5030B	Prep Date:06/24/2008 18:57	
Matrix:Water	Analytical	Method:8260B	Cal Date:04/29/2008 15:06	
Workgroup Number:WG274878	A	alyst:FJB	Run Date:06/24/2008 18:57	
Collect Date:06/16/2008 10:50	Di	lution:1	File ID:8M345917	
Sample Tag:01		Units:ug/L	······································	
	CAC No.	mhow	Popult Awal DI MDI	

Analyte	CAS. Numbe	r : 8	lesult	Qual	÷	RL	MDL
Tetrachloroethene	127-18-4	-	0.965		J	L.00	0.250
Toluene	108-88-3			υ		L.00	0.250
trans-1,2-Dichloroethene	156-60-5			U	1	1.00	0.250
trans-1,3-Dichloropropene	10061-02-	5		υ		L.00	0.500
Trichlorofluoromethane	75-69-4			U		1.00	0.250
Vinyl chloride	75-01-4			U	1	1.00	0.250
Surrogate	% Recovery	Lower	Uppe	er	Qual		
Dibromofluoromethane	108	85	115	5			
1,2-Dichloroethane-d4	110	72	119)			
Toluene-d8	92.7	81	120)			
4-Bromofluorobenzene	. 94.2	76	119)			

U Undetected; the analyte was analyzed for, but not detected.
R Because of quality control deficiences for this analyte, this data may be rejected.
F The analyte was positively identified, but the quantitation was below the RL.

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7/28/08

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Report Number: L08060559 Report Date :July 23, 2008

PrePrep Method:NONE	Instrument: HPMS14
Prep Method:5030B	Prep Date:05/27/2008 13:48
Analytical Method: 8260B	Cal Date:05/26/2008 20:24
Analyst: CMS	Run Date:06/27/2008 13:48
Dilution:1	File ID:14M06655
Units:ug/L	
	Prep Method:5030B Analytical Method:8260B Analyst: CMS Dilution:1

Analyte	CAS. Numb	er Re	esult	Qual	RL	MDL
1,4-Dioxane	123-91-	1	7.20		2.00	1.00
Surrogate	* Pogovory	Lower	Upper		Dual .	
- 1	3 Recovery		720		Zuar	
Dibromofluoromethane	TT2	24	130			
1,2-Dichloroethane-d4	117	51	135	· · · · · · · · · · · · · · · · · · ·		

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1125/08

Report Number: L08060559 Report Date :July 23, 2008

Sample Number: L08060559-06 Client ID:59SW7WG1 Matrix:Water Matrix: Water Workgroup Number: WG274878 Collect Date:06/15/2008 14:10 Sample Tag:01

 PrePrep Method: NONE
 Instrument: HPMS8

 Prep Method: 5030B
 Prep Date: 06/24/2008 15:55

 Analytical Method: 8260B
 Cal Date: 04/29/2008 15:06

 Analyst: FVB
 Run Date: 06/24/2008 15:55

 Dilution: 1
 File ID: 8M345912

 Units: ug/L
 File ID: 8M345912

Analyte	CAS. Number	Result	Qual	RL	MDL
1.1.1.2-Tetrachloroethane	630-20-6	·····	υ	0.500	0.250
1,1,1-Trichloroethane	71-55-6	2.50		1.00	0.250
1,1,2,2-Tetrachloroethane	79-34-5		U	0.500	0.125
1,1,2-Trichloroethane	79-00-5	· · · · · · · · · · · · · · · · · · ·	U	1.00	0.250
1,1-Dichloroethane	75-34-3	1.59		1.00	0,125
1,1-Dichloroethene	75-35-4		U	1.00	0.500
1,1-Dichloropropene	i 563-58-6		U	1.00	0.250
1,2,3-Trichlorobenzene	87-61-6	1	U	1.00	0.150
1,2,3-Trichloropropane	96-18-4	al p	U	1.00	0.500
1,2,4-Trichlorobenzene	120-82-1	······	v	1.00	0.200
1,2,4-Trimethylbenzene	95-63-6		U	1.00	0.250
1,2-Dichloroethane	107-06-2	÷	υ	0.500	0.250
1,2-Dichlorobenzene	95-50-1		บ	1.00	0.125
1,2-Dibromo-3-chloropropane	96-12-8		υ	2.00	1.00
	78-87-5	· · · · · · · · · · · · · · · · · · ·	υ	1.00	0.200
1,2-Dichloropropane	106-93-4	• • • • • • • • • • • • • • • • • • •	υ	1.00	0.250
1,2-Dibromoethane	and a second second second of the second	ļ			0.250
1,3,5-Trimethylbenzene	108-67-8		U	1.00	0.250
1,3-Dichlorobenzene	541-73-1		U	1.00	
1,3-Dichloropropane	142-28-9	s 2	U	0.400	0.200
1,4-Dichlorobenzene	106-46-7	1	υ	0.500	0.125
1,4-Dioxane	123-91-1		M	100	50.0
1-Chlorohexane	544-10-5		σ	1.00	0.125
2,2-Dichloropropane	594-20-7	• •	U,	1.00	0.250
2-Chlorotoluene	95-49-8	с Ј	U	1.00	0.125
4-Chlorotoluene	106-43-4		U	1.00	0.250
Acetone	67-64-1		U	10.0	2.50
Benzene	71-43-2		U	0.400	0.125
Bromobenzene	108-86-1		U	1.00	0.125
Bromochloromethane	74-97-5	•	υ	1.00	0.200
Bromodichloromethane	75-27-4		U	0.500	0.250
Bromoform	75-25-2		τ :	1.00	0.500
Bromomethane	74-83-9	•	U	3.00	0.500
Carbon tetrachloride	56-23-5	••••••••••••••••••••••••	σ	1.00	0.250
Chlorobenzene	108-90-7	· · · · · · · · · · · · · · · · · · ·	U U	0.500	0.125
Chloroethane	75-00-3		U	1.00	0.500
Chloroform	67-66-3	÷	v	0.300	0.125
Chloromethane	74-87-3		Ū	1.00	0.250
cis-1,2-Dichloroethene	156-59-2	6.34		1.00	0.250
	10061-01-5	0.34	U	0.500	0.250
cis-1,3-Dichloropropene			U U	0.500	0.250
Dibromochloromethane	124-48-1	· · · · · · · · · · · · · · · · · · ·	- and a room for some	any second a second a second second second scheme & s	0.250
Dibromomethane	74-95-3	· ·	U	1.00	
Dichlorodifluoromethane	75-71-8		R-()_[1.00	0.250
Ethylbenzene	100-41-4		U :	1.00	0.250
Hexachlorobutadiene	87-68-3	· · •	U	0.600	0.250
Isopropylbenzene	98-82-8		υ	1.00	0.250
Methylene chloride	75-09-2	• 	U	1.00	0.250
Methyl t-butyl ether (MTBE)	1634-04-4		υ	5.00	0.500
MEK (2-Butanone)	78-93-3		U	10.0	2.50
MIBK (methyl isobutyl ketone)	108-10-1	·	υ	10.0	2.50
n-Butylbenzene	104-51-8		υ	1.00	0.250
n-Propylbenzene	103-65-1		U	1.00	0.125
m-,p-Xylene	136777-61-2	· · · · · · · · · · · · · · · · · · ·	U	2.00	0.500
Naphthalene	91-20-3	:	U	1.00	0.200
o-Xylene	95-47-6		υ	1.00	0.250
p~Isopropyltoluene	99-87-6		U	1.00	0.250
sec-Butylbenzene	135-98-8	,	ΰ	1.00	0.250
Styrene	100-42-5		U	1,00	0.125
Trichloroethene	79-01-6	2.94		1.00	0,250
tert-Butylbenzene	98-06-6	4.33	U	1.00	0.250

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Sample Number:L08060559-06	PrePrep Method:NONE	Instrument:HPMS8
Client ID: 59SW7WG1	Prep Method: 5030B	Prep Date:06/24/2008 15:55
Matrix:Water	Analytical Method:8260B	Cal Date:04/29/2008 15:06
Workgroup Number:WG274878	Analyst FJB	Run Date:06/24/2008 15:55
Collect Date:06/15/2008 14:10	Dilution:1	File ID:8M345912
Sample Tag:01	Units:ug/L	
	· · · · · · · · · · · · · · · · · · ·	
analyte	CAS Number : Result	Oual RL MDL

Analyte) (20.0. 190, 190, 190, 190, 190, 190, 190, 190,	24 : I		1 Sugr	1	1111	
Tetrachloroethene	127~18-4		0.843	- #	1 1	L.00	0.250
Toluene	108-88-3			U]]	L.00	0.250
trans-1,2-Dichloroethene	156-60-5			U	3	L.00	0.250
trans-1,3-Dichloropropene	10061-02-	6		U]]	1.00	0.500
Trichlorofluoromethane	75-69-4			υ] I	ι.00	0.250
Vinyl chloride	75-01-4	1		υ	1	1.00	0.250
Surrogate	% Recovery	Lower	Uppe:	r !	Qual		
Dibromofluoromethane	111	85	115	1			
1,2-Dichloroethane-d4	114	72	119	1			
Toluene-d8	92.9	81	. 120	1			
4-Bromofluorobenzene	95.1	76	: 119	1			

M Matrix effect; the concentration is an estimate due to matrix effect. U Undetected; the analyte was analyzed for, but not detected. R Because of quality control deficiences for this analyte, this data may be rejected. F The analyte was positively identified, but the quantitation was below the RL.

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7/28/03 9U

Sample Number:L08060559-06	PrePrep Method:	NONE	Instrument: HPMS14	
Client ID: 59SW7WG1	Prep Method:	5030B	Prep Date:06/27/2008	14:23
Matrix:Water	Analytical Method:	8260B	Cal Date:06/26/2008	20:24
Workgroup Number: WG275287	Analyst:	CMS	Run Date:06/27/2008	1,4:23
Collect Date:06/15/2008 14:10	Dilution:		File ID:14M06656	
Sample Tag:01	Units:	ug/L		
Analyte	CAS. Number	Result	Qual RL	MDL

Analyte	CAS. Number	Result	Qual	КL	MDL	
1.4-Dioxane	123-91-1	4.66		2.00	1.00	
Surrogate		wer Uppe	r Qual	1		
Dibromofluoromethane	112 5	4 138				
1,2-Dichloroethane-d4	114	135				

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Sample Number: L08060559-07PrePrep Method: NONEInstrument: HPMS8Client ID: 595W7WG1-MSPrcp Method: 5030BPrep Date: 06/24/2008 16:47Matrix: WaterAnalytical Method: 8260BCal Date: 04/29/2008 15:06Workgroup Number: WG274878Analyst: PJBRun Date: 06/24/2008 16:47Collect Date: 06/15/2008 14:10Dilution:1File ID: 8M345913Sample Tag: 01Units: ug/LFile ID: 8M345913AnalyteCAS. NumberResultQualRLMDL1, 1, 2-Tetrachloroethane630-20-622.00.5009/250

1,1,1,2-Tetrachloroethane	630-20-6	22.0	(0.500	0.250
1,1,1-Trichloroethane	71-55-6	24.7		1.00	0.250
1,1,2,2-Tetrachloroethane	79-34-5	17.8	(.500	0.125
1,1,2-Trichloroethane	79-00-5	19.8		1.00	0.250
,1-Dichloroethane	75-34-3	22.5		1.00	0.125
,1-Dichloroethene	75-35-4	18.4	· · · · · · · · · ·	1.00/	0.500
, 1-Dichloropropene	563-58-6	19.1		1.00	0.250
,2,3-Trichlorobenzene	87-61-6	19.7		1.00	0.150
	96-18-4	19.6		1.00	0.500
., 2, 3-Trichloropropane	120-82-1	19.8		1.00	0,200
1,2,4-Trichlorobenzene	95-63-6	21.5		1.00	0.250
,2,4-Trimethylbenzene	107-06-2	22.9	and a second second second second second	1.00 0.500	0.250
,2-Dichloroethane	and a second second base of the second s	20.4		1.00	0.125
L,2-Dichlorobenzene	95-50-1				1.00
.,2-Dibromo-3-chloropropane	96-12-8	19.3		2.00	
,2-Dichloropropane	78-87-5	21.0	en la sur	1.00	0.200
1,2-Dibromoethane	106-93-4	19.2		1.00	0.250
1,3,5-Trimethylbenzene	108-67-8	18.7	and more a second second of	1.00	0.250
L,3-Dichlorobenzene	541-73-1	20.1		1.00	0.250
,3-Dichloropropane	142-28-9	20.0		0.400	0.200
,4-Dichlorobenzene	106-46-7	18.8		0.500	0.125
L,4-Dioxane	123-91-1	93.2	F	100	50.0
l-Chlorohexane	544-10-5	17.7		1.00	0.125
2,2-Dichloropropane	594-20-7	18.2		1.00	0.250
2-Chlorotoluene	95-49-8	21.0		1.00	0.125
-Chlorotoluene	106-43-4	21.2		1.00	0.250
Acetone	67-64-1	19.1		10.0	2.50
Benzene	71-43-2	21.4	····	0.400	0.125
3romobenzene	108-86-1	19.7		1.00	0.125
Bromochloromethane	74-97-5	20.1		1.00	0.200
Bromodichloromethane	75-27-4	23.7		0.500	0.250
Bromodichioromechane	75-25-2	20.8	••••	1.00	0.500
······································	75-25-2	16.2		3.00	0.500
Bromomethane	56-23-5	21.1		1.00	0.250
Carbon tetrachloride	108-90-7			0.500	0.125
Chlorobenzene		20.2			0.125
Chloroethane	75-00-3	17.4		1.00	CONTRACTOR AND A CONTRACTOR AND
Chloroform	67-66-3	22.2		0.300	0.125
Chloromethane	74-87-3	1.8.4		1.00	0.250
cis-1,2-Dichloroethene	156-59-2	28.3		1.00	0.250
cis-1,3-Dichloropropene	10061-01-5	20.9		0.500	0.250
Dibromochloromethane	124-48-1	. 18.7		0.500	0.250
Dibromomethane	74-95-3	22.3		1.00	0.250
Dichlorodifluoromethane	75-71-8	19.5	R	1.00	0.250
Ethylbenzene	100-41-4	21.7		1.00	0.250
Hexachlorobutadiene	87-68-3	19.4		0.600	0.250
Isopropylbenzene	98-82-8	16.8		1.00	0.250
Methylene chloride	75-09-2	18.3		1.00	0.250
Methyl t-butyl ether (MTBE)	1634-04-4	19.5		5.00	0.500
MEK (2-Butanone)	78-93-3	19.3		10.0	2.50
MIRK (2-Butanone) MIBK (methyl isobutyl ketone)	108-10-1	15.7		10.0	2.50
n-Butylbenzene	104-51-8	18.4		1.00	0.250
n-Propylbenzene	103-65-1	18.1		1.00	0.125
see see an All All and the second	136777-61-2	43.7		2.00	0.500
m-,p-Xylene				1.00	0.200
Naphthalene	91-20-3	21.1			
o-Xylene	95-47-6	21.4		1.00	0.250
p-Isopropyltoluene	99-87-6	17.9		1.00	0.250
sec-Butylbenzene	135-98-8	18.0		1.00	0.250
Styrene	100-42-5	19.0		1.00	0.125
Trichloroethene /	79-01-6	21.1	1	1.00	0.250
tert-Butylbenzene	98-06-6	17.5		1.00	0.250

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Report Number: L08060559

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Report Date : July 23, 2008

Sample Number: L08060559-07	PrePrep Method:N	ONE	Instrument:HPMS8	· ·
Client ID: 59SW7WG1-MS	Prep Method:5	030B	Prep Date:06/24/2	2008 16:47
Matrix:Water	Analytical Method:8	260B	Cal Date:04/29/2	2008 15:06
orkgroup Number:WG274878	Analyst: F	JB	Run Date:06/24/2	2008 16:47
Collect Date:06/15/2008 14:10	Dilution:1		File ID:8M345913	
Sample Tag:01	Units:u	g/L		
Analyte	CAS. Number	Result	Qual RL	MDL
Tetrachloroethene	127-18-4	17.8	1,00	0.250
Toluene	108-88-3	20.8	1.00	0.250
trans-1,2-Dichloroethene	156-60-5	21.3	/1.00	0.250
trans-1,3-Dichloropropene	10061-02-6	18.4	1.00	0.500
Trichlorofluoromethane	75-69-4	19.7	1.00	0.250
Vinyl chloride	75-01-4	26.3	1.00	0.250
Surrogate	% Recovery Lo	wer Upp	er / Qual ;	
	100	C 33	c /	

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Idea10685108721087210894.94-Bromofluorobenzene91.891.876R Because of quality control deficiences for this analyticF The analyte was positively ideative 119 R Because of quality control deficiences for this analyte, this data may be rejected. F The analyte was positively identified, but the quantitation was below the RL.

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Sample Number:L08060559-07	PrePrep Method	NONE	Ins	strument: HPMS	14
Client ID:59SW7WG1-MS	Prep Method	:5030B	Pi	rep Date:06/2	7/2008 14:58
Matrix:Water	Analytical Method	:8260B	(Cal Date:06/2	6/2008 20:24
kgroup Number:WG275287	Analyst	CMS	E. S. F.	Run Date:06/2	
Collect Date:06/15/2008 14:10	Dilution	1:1		le ID:14M0665	57/
Comple Tech 01	Units	uq/L			7
Sample Tag:01	•				
	CAS. Number		t Qua		MDL
	CAS. Number 123-91-1		t Qua		MDL 1.00
halyte	CAS. Number 123-91-1	Resul	t Qua	l RL	
Analyte 1,4-Dioxane	CAS. Number 123-91-1	Resul 13.8	t Qua	1 RL 2.00	

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		6	of	9
7				

Sample Number:L08060559-08	PrePrep Method:NONE	Instrument:HPMS8
Client ID:59SW7WG1-MSD	Prep Method:5030B	Prep Date:06/24/2008 17:19
Matrix:Water	Analytical Method:8260B	Cal Date:04/29/2008 15:06
Workgroup Number:WG274878	Analyst:FJB	Run Date:06/24/2008 17:19
Collect Date:06/15/2008 14:10	Dilution:1	File ID:8M345914
Sample Tag:01	Units:ug/L	and the second sec

Analyte	CAS. Number	Result	Qual RL	MDL
1,1,2-Tetrachloroethane	630-20-6	21.6	0.500	0.250
1,1-Trichloroethane	71-55-6	24.3	1.00	0.250
1,2,2-Tetrachloroethane	79-34-5	18,7	0.500	0.125
1,2-Trichloroethane	79-00-5	20.5	1.00	0.250
1-Dichloroethane	75-34-3	22.9	1.00	0.125
1-Dichloroethene	75-35-4	18.7	1.00	0.500
1-Dichloropropene	563-58-6	18.8	1.00	0.250
2,3-Trichlorobenzene	87-61-6	20.9	1.00	0.150
2,3-Trichloropropane	96-18-4	19.6	1.00	0.500
2,4-Trichlorobenzene	120-82-1	20.6	1.00	0.200
2,4-Trimethylbenzene	95-63-6	21.5	1.00	0.250
2-Dichloroethane	107-06-2	23.1	and a property of the second	0.250
2-Dichlorobenzene	95-50-1	20.7	1.00	0.125
2-Dibromo-3-chloropropane 2-Dichloropropane	96-12-8 78-87-5	20.2	2.00	0.200
	106-93-4	20.1	1.00	0.200
2-Dibromoethane	108-67-8	18,5	1.00	0.250
3,5-Trimethylbenzene	541-73-1	19.9	1.00	0.250
3-Dichlorobenzene	142-28-9	20.3	CONTRACTOR AND A D A CONTRACTOR AND AND A CONTRACTOR AND AND A CONTRACTOR AND AND A CONTRACTOR AND A CONTRACTOR AND A CONTRACTOR AND A CONTRAC	0.200
3-Dichloropropane 4-Dichlorobenzene	106-46-7	19.1	0.400	0.125
4-Dioxane	123-91-1	152	100	50.0
-Chlorohexane	544-10-5	17.7	1.00	0.125
2-Dichloropropane	594-20-7	17.9	1.00	0.250
-Chlorotoluene	95-49-8	22.0	1.00	0.125
-Chlorotoluene	106-43-4	19.7	1.00	0.250
cetone	67-64-1	21.5	10.0	2.50
enzene	71-43-2	21.3	0.400	0.125
comobenzene	108-86-1	20.2	1.00	0.125
romochloromethane	74-97-5	21.0	1.00	0.200
romodichloromethane	75-27-4	23.8	0.500	0.250
romoform	75-25-2	21.3	1.00	0.500
romomethane	74-83-9	17.2	3.00	0.500
arbon tetrachloride	56-23-5	20.3	1.00	0.250
hlorobenzene	108-90-7	19.9	0.500	0.125
hloroethane	75-00-3	18.1	1.00	0.500
hloroform	67-66-3	22.2	0.300	0.125
hloromethane	74-87-3	17.1	1.00	0.250
is-1,2-Dichloroethene	156-59-2	29.1	1.00	0.250
is-1,3-Dichloropropene	10061-01-5	21.4	0.500	0.250
ibromochloromethane	124-48-1	18.7	0.500	0.250
ibromomethane	74-95-3	23.2	1.00	0.250
ichlorodifluoromethane	75-71-8	19.5	R 1.00	0.250
thylbenzene	100-41-4	20.8	1.00	0.250
exachlorobutadiene	87-68-3	18.7	0.600	0.250
sopropylbenzene	98-82-8	16.5	1.00	0.250
ethylene chloride	75-09-2	18.2	1.00	0.250
ethyl t-butyl ether (MTBE)	1634-04-4	21.8	5.00	0.500
EK (2-Butanone)	78-93-3	22.2	10.0	2.50
IBK (methyl isobutyl ketone)	108-10-1	18.4	10.0	2.50
-Butylbenzene	104-51-8	18.2	1.00	0.250
-Propylbenzene	103-65-1	17.9	1.00	0.125
-,p-Xylene	136777-61-2	43.0	2.00	0.500
aphthalene	91-20-3	22.5	1.00	0.200
-Xylene	95-47-6	21.0	1.00	0.250
-Isopropyltoluene	99-87-6	17.7	1.00	0.250
ec-Butylbenzene	135-98-8	17.9	1.00	0.250
tyrene	100-42-5	18.6	1.00	0.125
richloroethene ert-Butylbenzene	79-01-6 98-06-6	21.1 17.3	1.00	0.250
	15 of 36		1.00	
1				
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Report Number: L08060559 Report Date :July 23, 2008

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Sample Number: L08060559-08	PreFrep Method:NONE	Instrument:HPMS8	1
Client ID: 59SW7WG1-MSD	Prep Method:5030B	Prep Date:06/24/200	8 17:19
Matrix:Water	Analytical Method:8260B	Cal Date:04/29/200	8 15:06
Workgroup Number:WG274878	Analyst PJB	Run Date:06/24/200	8 17:19
Collect Date:06/15/2008 14:10	Dilution:1	File ID:8M345914	2
Sample Tag:01	Units:ug/L		
Analyte	CAS. Number Result	Qual RL	MDL

Tetrachloroethene	127-18-4	:	17.5	1.00	0.250
Toluene	108-88-3		20.4	1.00	0.250
trans-1,2-Dichloroethene	156-60-5		22.0	1.00	0.250
trans-1,3-Dichloropropene	10061-02-6		19.0	1.00	0.500
Trichlorofluoromethane	75-69-4		18.5	1.00	0.250
Vinyl chloride	75-01-4		24.3	1.00	0.250
Surrogate	% Recovery	Lower	Upper	Qual	
Dibromofluoromethane	106	85	115		
1,2-Dichloroethane-d4	105	72	119		
Toluene-d8	94.0	81	120		
4-Bromofluorobenzene	92.2	76	119		

R Because of quality control deficiences for this analyte, this data may be rejected.

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Sample Number:L08060559-08	PrePrep Method:NONE	Instrument: HPMS14
Client ID:59SW7WG1-MSD	Prep Method:5030B	Prep Date:06/27/2008 15:34
Matrix:Water	Analytical Method:8260B	Cal Date:06/26/2008 20:24
Workgroup Number:WG275287	Analyst:CMS	Run Date:06/27/2008 15:34
Collect Date: 06/15/2008 14:10	Dilution:1	File ID:14M06658
Sample Tag:01	Units:ug/L	

Analyte	CAS.	Number		Result	Qual	RL	MDL
1,4-Dioxane	123-	-91-1		14.6		2.00	1.00
Surrogate	% Recove	ry :	Lower	Upper	r Qua	11	
Dibromofluoromethane	108		54	138			
1,2-Dichloroethane-d4	111	;	51	135	- 17		

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Report Number: L08060559 Report Date :July 23, 2008

Isopropylbenzene Methylene chloride

MIBK (methyl isobutyl ketone)

MEK (2-Butanone)

n-Propylbenzene

Trichloroethene

tert-Butylbenzene

o-Xylene

Styrene

m-,p-Xylene

p-Isopropyltoluene

sec-Butylbenzene

Naphthalene

Methyl t-butyl ether (MTBE)

n-Butylbenzene

Sample Number: L08060559-10	PrePrep Method:NC	Instrument: HPMS14							
Client ID:GS-95025WG-1		Prep Method:50308 Analytical Method:82608			Prep Date:06/25/2008 18:59 Cal Date:06/24/2008 22:32				
Matrix:Water									
kgroup Number:WG275043	Analyst : Ch	15	Run Date:06/25/2008 18:59						
Collect Date:06/15/2008 19:30	Dilution:1		, rile i	D:14M06612	· · · · · · · · · · · · · · · · · · ·				
Sample Tag:01	Units:ug	1/L							
Analyte	CAS. Number	Result	Qual	RL	MDL				
1,1,1,2-Tetrachloroethane	630-20-6		U	0.500	0.250				
1,1,1-Trichloroethane	71-55-6	9.04		1.00	0.250				
1,1,2,2-Tetrachloroethane	79-34-5		U	0.500	0.125				
1,1,2-Trichloroethane	79-00-5	1	υ	1.00	0.250				
1,1-Dichloroethane	75-34-3	;	U	1.00	0.125				
1,1-Dichloroethene	75-35-4	}	ប	1.00	0.500				
1,1-Dichloropropene	563-58-6	1	บ :	1.00	0.250				
1,2,3-Trichlorobenzene	87-61-6		U	1.00	0.150				
1,2,3-Trichloropropane	96-18-4		U	1.00	0.500				
1,2,4-Trichlorobenzene	120-82-1	1	U	1.00	0.200				
1,2,4-Trimethylbenzene	95-63-6		U :	1.00	0.250				
1,2-Dichloroethane	107-06-2		υ.	0.500	0.250				
1,2-Dichlorobenzene	95-50-1		U	1.00	0.125				
1,2-Dibromo-3-chloropropane	96-12-8		U ·	2.00	1.00				
1,2-Dichloropropane	78-87-5		U	1.00	0.200				
1,2-Dibromoethane	106-93-4		U	1.00	0.250				
1,3,5-Trimethylbenzene	108-67-8	-	ប	1.00	0.250				
1,3-Dichlorobenzene	541-73-1	1	υ.	1.00	0.250				
1,3-Dichloropropane	142-28-9		ប	0.400	0.200				
1,4-Dichlorobenzene	106-46-7		U	0.500	0.125				
1,4-Dioxane	123-91-1		U	100	50.0				
1-Chlorohexane	544-10-5		U	1.00	0.125				
2,2-Dichloropropane	594-20-7	·	Ŭ	1.00	0.250				
2-Chlorotoluene	95-49-8	· · · · · · · · · · · · · · · · · · ·	υ	1.00	0.125				
4-Chlorotoluene	1.06-43-4	·	ប	1.00	0.250				
Acetone	67-64-1		U	10.0	2.50				
Benzene	71-43-2		U	0.400	0.125				
Bromobenzene	108-86-1		U	1.00	0,125				
Bromochloromethane	74-97-5	ļ	U	1.00	0.200				
Bromodichloromethane	75-27-4		υ	0.500	0.250				
Bromoform	75-25-2		U	1.00	0.500				
Bromomethane	74-83-9		U	3.00	0.500				
Carbon tetrachloride	56-23-5		U	1.00	0.250				
Chlorobenzene	108-90-7		U	0.500	0.125				
Chloroethane	75-00-3	: 	U	1.00	0.500				
Chloroform	67-66-3		U	0.300	0.125				
Chloromethane	74-87-3		U	1.00	0.250				
cis-1,2-Dichloroethene	156-59-2		U U	1.00	0.250				
cis-1,3-Dichloropropene	10061-01-5		1	0.500	en la cara a diservaria di cherciano.				
Dibromochloromethane	124-48-1	.	U	1.00	0.250				
Dibromomethane	74-95-3		υ		and a set to and second some its stations do				
Dichlorodifluoromethane	75-71-8		R U U) 1.00	0.250				
Ethylbenzene	100-41-4 87-68-3	· · · · · · · · · · · · · · · · · · ·	U I	0.600	0.250				
Hexachlorobutadiene			U U	1.00	0.250				
Isopropylbenzene	98-82-8	1	1 U)	1.00 2	V.23V				

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75-09-2

1634-04-4

78-93-3

108-10-1

104-51-8

103-65-1

136777-61-2

91-20-3

95-47-6

99-87-6

135-98-8

100-42-5

79-01-6

98-06-6

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Sample Number:L08060559-10	PrePrep Method:NONE	Instrument: HPMS14
Client ID: GS-95025WG-1	Prep Method:5030	B Prep Date:06/25/2008 18:59
Matrix:Water	Analytical Method:8260	B Cal Date:06/24/2008 22:32
Workgroup Number:WG275043	Analyst:CMS	Run Date:06/25/2008 18:59
Collect Date:06/15/2008 19:30	Dilution:1	File ID:14M06612
Sample Tag:01	Units:ug/L	
Analyte	CAS. Number	Result Qual RL MDL

marycc		÷ .		Y GOL		
Tetrachloroethene	127-18-4	l		U	1.00	0.250
Toluene	108-88-3			υ	1.00	0.250
trans-1,2-Dichloroethene	156-60-5			U	1.00	0.250
trans-1,3-Dichloropropene	10061-02-6	5		υ	1.00	0.500
Trichlorofluoromethane	75-69-4			U	1.00	0.250
Vinyl chloride	75-01-4			R-())	1.00	0.250
Surrogate	% Recovery	Lower	Uppe	r Qua	1	
Dibromofluoromethane	100	85	115			
1,2-Dichloroethane-d4	104	72	119			
Toluene-d8	102	81	120			
4-Bromofluorobenzene	103	76	119			

U Undetected; the analyte was analyzed for, but not detected. R Because of quality control deficiences for this analyte, this data may be rejected.

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Sample Number: L08060559-09 Client ID:GS-9502DWG-1 Workgroup Number: Water Collect Date: 06/15/2008 18:10 Sample Tag:01

 PrePrep Method:NONE
 Instrument:HPMS14

 Prep Method:5030B
 Prep Date:06/25/2008 18:27

 Analytical Method:8260B
 Cal Date:06/24/2008 22:32

 Analyst:CMS
 Run Date:06/25/2008 18:27

 Dilution:1
 File ID:14M06611

 Units:ug/L

Analyte	CAS. Number	Result	Qual	RL	MDT,
1,1,1,2-Tetrachloroethane	630-20-6		U	0.500	0.250
1,1,1-Trichloroethane	71-55-6		ប	1.00	0.250
1,1,2,2-Tetrachloroethane	79-34-5		σ	0.500	0.125
1,1,2-Trichloroethane	79-00-5		U	1.00	0.250
1,1-Dichloroethane	75-34-3		U	1.00	0.125
1,1-Dichloroethene	75-35-4		U	1.00	0.500
1,1-Dichloropropene	563-58-6	····· ··· ···· ···· ··· ··· ··· ··· ··	U	1.00	0.250
1,2,3-Trichlorobenzene	87-61-6		υ	1.00	0.150
1,2,3-Trichloropropane	96-18-4		ប	1.00	0.500
1,2,4-Trichlorobenzene	120-82-1		U	1.00	0.200
1,2,4-Trimethylbenzene	95-63-6		υ	1.00	0.250
1,2-Dichloroethane	107-06-2		U	0.500	0.250
1,2-Dichlorobenzene	95-50-1	• • • • • • • • • • • • • • • • • • • •	Ū	1.00	0.125
1,2-Dibromo-3-chloropropane	96-12-8		U	2.00	1.00
1,2-Dichloropropane	78-87-5	· · · · · · · · · · · · · · · · · · ·	U	1.00	0.200
	106-93-4		- U	1.00	0.250
1,2-Dibromoethane			υ υ.	1.00	
1,3,5-Trimethylbenzene	108-67-8		.		0.250
1,3-Dichlorobenzene	541-73-1		U	1.00	0.250
1,3-Dichloropropane	142-28-9		U	0.400	0.200
1,4-Dichlorobenzene	106-46-7		U	0.500	0.125
1,4-Dioxane	123-91-1		Ū	100	50.0
1-Chlorohexane	544-10-5		U	1.00	0.125
2,2-Dichloropropane	594-20-7		U	1.00	0.250
2-Chlorotoluene	95-49-8		ប	1.00	0.125
4-Chlorotoluene	106-43-4		U	1.00	0.250
Acetone	67-64-1		U	10.0	2.50
Benzene	71-43-2		U	0.400	0.125
Bromobenzene	108-86-1		U	1.00	0.125
Bromochloromethane	74 - 97 - 5		U	1.00	0.200
Bromodichloromethane	75-27-4		U	0.500	0.250
Bromoform	75-25-2		U	1.00	0.500
Bromomethane	74-83-9		υ	3.00	0.500
Carbon tetrachloride	56-23-5		U	1.00	0.250
Chlorobenzene	108-90-7		U	0.500	0.125
Chloroethane	75-00-3	·····	υ	1.00	0.500
Chloroform	67-66-3		U	0.300	0.125
Chloromethane	74-87-3	·	U	1.00	0.250
cis-1,2-Dichloroethene	156-59-2	·· ·· ·· ·· ·· ·· ·· ·· ·· ·· ·· ·· ··	ប	1.00	0.250
cis-1,3-Dichloropropene	10061-01-5	·	υ	0.500	0.250
Dibromochloromethane	124-48-1		U	0.500	0.250
Dibromomethane	74-95-3	· · · · · · · · · · · · · · · · · · ·	U	1.00	0.250
Dichlorodífluoromethane	75-71-8	······	-R-11	1.00	0.250
Ethylbenzene	100-41-4			1.00	0.250
Hexachlorobutadiene	87-68-3		Ū	0.600	0.250
		- I	υ	1.00	
Isopropylbenzene	98-82-8		U	1.00	0.250
Methylene chloride	75-09-2	n general and an and an	de morrere de	and the second	0.500
Methyl t-butyl ether (MTBE)	1634-04-4		<u></u>) 5.00	et contras. In a contrast to
MEK (2-Butanone)	78-93-3		U	10.0	2.50
MIBK (methyl isobutyl ketone)	108-10-1		U	10.0	2.50
n-Butylbenzene	104-51-8		U	1.00	0.250
n-Propylbenzene	103-65-1		U	1.00	0.125
m-,p-Xylene	136777-61-2		U ;	2.00	0.500
Naphthalene	91-20-3		U	1.00	0.200
o-Xylene	95-47-6		U	1.00	0.250
p-Isopropyltoluene	99-87-6		U	1.00	0.250
sec-Butylbenzene	135-98-8		U	1.00	0.250
Styrene	100-42-5		U	1.00	0.125
Trichloroethene	79-01-6		U	1.00	0.250
tert-Butylbenzene	98-06-6		υ	1.00	0.250

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Milliobic

7/28/05

Sample Number: L08060559-09	PrePrep Method:NONE	Instrument: HPMS14
Client ID:GS-9502DWG-1	Prep Method:5030B	 Prep Date:06/25/2008 18:27
Matrix:Water	Analytical Method:8260B	 Cal Date:06/24/2008 22:32
Workgroup Number:WG275043	Analyst:CMS	 Run Date:06/25/2008 18:27
Collect Date:06/15/2008 18:10	Dilution:1	 File ID:14M06611
Sample Tag:01	Units:ug/L	

Analyte	CAS. Numbe	τ .	Result	Qual	RL	MDL
Tetrachloroethene	127-18-4			U	1.00	0.250
Toluene	108-88-3			U	1.00	0.250
trans-1,2-Dichloroethene	156-60-5			U	1.00	0.250
trans-1,3-Dichloropropene	10061-02-0	6		ប	1.00	0.500
Trichlorofluoromethane	75-69-4			U	1.00	0.250
Vinyl chloride	75-01-4			-17-) 1.00	0.250
Surrogate	% Recovery	Lower	Upp	er	Qual	
Dibromofluoromethane	99.8	85	11	5		
1,2-Dichloroethane-d4	103	72	. 11	9		
Toluene-d8	103	81	12	0		
4-Bromofluorobenzene	103	76	11	9		

U Undetected; the analyte was analyzed for, but not detected. R Because of quality control deficiences for this analyte, this data may be rejected.

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Microl

712×103

Sample Number:L08060559-11	PrePrep Method:NONE	Instrument:HPMS14
Client ID:GS-9505WG-1	Prep Method: 5030B	Prep Date:06/25/2008 19:32
Matrix:Water	Analytical Method:8260B	Cal Date:06/24/2008 22:32
Workgroup Number:WG275043	Analyst: CMS	Run Date:06/25/2008 19:32
Collect Date: 06/15/2008 20:45	Dilution:1	File ID:14M06613
Sample Tag:01	Units:ug/L	

Analyte	CAS. Number	Result	Qual	RL	MDL
,1,1,2-Tetrachloroethane	630-20-6	-	U	0.500	0.250
1,1-Trichloroethane	71-55-6	4.49		1.00	0.250
1,2,2-Tetrachloroethane	79-34-5		U	0.500	0.125
1,2-Trichloroethane	79-00-5		U	1.00	0.250
1-Dichloroethane	75-34-3	0.389	4	1.00	0.125
1-Dichloroethene	75-35-4	1	U	1.00	0.500
1-Dichloropropene	563-58-6	· · · · · · · · · · · · · · · · · · ·	U	1.00	0.250
,2,3-Trichlorobenzene	87-61-6		υ	1.00	0.150
,2,3-Trichloropropane	96-18-4		U	1.00	0.500
,2,4-Trichlorobenzene	120-82-1		U i	1.00	0.200
,2,4-Trimethylbenzene	95-63-6		U :	1.00	0.250
,2-Dichloroethane	107-06-2		U I	0.500	0.250
,2-Dichlorobenzene	95-50-1	······································	Ū	1.00	0.125
,2-Dibromo-3-chloropropane	96-12-8		ប	2.00	1.00
,2-Dichloropropane	78-87-5		U	1.00	0.200
,2-Dibromoethane	106-93-4		υ	1.00	0.250
, 3, 5-Trimethylbenzene	108-67-8		Ŭ	1.00	0.250
.,3,5-171metnyidenzene .,3-Díchlorobenzene	541-73-1		U	1.00	0.250
· · · · · · · · · · · · · · · · · · ·	142-28-9		U U	0.400	0.200
,3-Dichloropropane ,4-Dichlorobenzene	142-28-9		υ	0.500	0.125
	123-91-1		υ	100	50.0
.,4-Dioxane			U	1.00	Firsterner and strend strends and strends
-Chlorohexane	544-10-5		eres hannel is a fire	1.00	0.125
2,2-Dichloropropane	591-20-7		U	and a second	0.250
-Chlorotoluene	95-49-8		U	1.00	0.125
-Chlorotoluene	106-43-4		U	1.00	0.250
lectone	67 - 64 - 1		U .	10.0	2.50
Senzene	71-43-2		U ;	0.400	0.125
romobenzene	108-86-1		U	1.00	0.125
bromochloromethane	74-97-5		U	1.00	0.200
romodichloromethane	75-27-4		υ	0.500	0.250
Bromoform	75-25-2	4	U	1.00	0.500
Bromomethane	74-83-9		U	3.00	0.500
Carbon tetrachloride	56-23-5	;	U	1.00	0.250
'hlorobenzene	108-90-7		U	0.500	0.125
Chloroethane	75-00-3	1	υ,	1.90	0.500
Chloroform	67-66-3	1	U,	0.300	0.125
Chloromethane	74-87-3		U	1.00	0.250
is-1,2-Dichloroethene	156-59-2		U	1.00	0.250
cis-1,3-Dichloropropene	10061-01-5		υ	0.500	0.250
Dibromochloromethane	124-48-1		ប	0.500	0.250
Dibromomethane	74-95-3		υ	1.00	0.250
Dichlorodifluoromethane	75-71-8		R () 1.00	0.250
Sthylbenzene	100-41-4		U	1.00	0.250
fexachlorobutadiene	87-68-3		U 1	0.600	0.250
Isopropylbenzene	98-82-8		ບ [1.00	0.250
Methylene chloride	75-09-2		υ	1.00	0.250
Methyl t-butyl ether (MTBE)	1634-04-4		·) \ 5.00	0.500
MEK (2-Butanone)	78-93-3		Ū	10.0	2.50
fIBK (methyl isobutyl ketone)	108-10-1		Ū	10.0	2.50
1-Butylbenzene	104-51-8		Ū	1.00	0.250
1-BityiDenzene	103-65-1	····	υ	1.00	0.125
n-,p-Xylene	136777-61-2		υ	2.00	0.500
w-, p-xylene Vaphthalene	91-20-3	···•	υ	1.00	0.200
vapnthalene o-Xylene	91-20-3	· · · · · · · · · · · · · · · · · · ·	UUU	1.00	0.250
			UUU	1.00	0.250
p-Isopropyltoluene	99-87-6				0.250
sec-Butylbenzene	135-98-8		U	1.00	
Styrene	100-42-5		ប	1.00	0.125
Frichloroethene tert-Butylbenzene	79-01-6 98-06-6	····-	U :	1.00	0.250

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1123/22

PrePrep Method:NONE	lostrument:HPMS14
Prep Method:5030B	Prep Date:06/25/2008 19:32
Analytical Method:8260B	Cal Date:06/24/2008 22:32
Analyst:CMS	Run Date:06/25/2008 19:32
Dilution:1	File ID:14M06513
Units:ug/L	
	Prep Method: 5030B Analytical Method: 8260B Analyst: CMS Dilution: 1 Units: ug/L

CAS. Numbe	er	Result	Qual		RL		MDL
127-18-4			U		1.00		0.250
108-88-3			U		1.00		0.250
156-60-5			U		1.00		0.250
10061-02-	6		U		1.00	1	0.500
75-69-4			U	1	1.00		0.250
75-01-4			TR \	33	1.00	į	0.250
% Recovery	Lower	Upp	er	Qual			
100	85	11	5		-		
104	72	11	9		-		
103	81	12	0		1		
102	76	11	9		:		
	127-18-4 108-88-3 156-60-5 10061-02- 75-69-4 75-01-4 % Recovery 100 104 103	108-88-3 156-60-5 10061-02-6 75-69-4 75-01-4 % Recovery 100 85 104 72 103	127-18-4 108-88-3 156-60-5 10061-02-6 75-69-4 75-01-4 % Recovery 100 85 11 103 12	127-18-4 U 108-88-3 U 156-60-5 U 10061-02-6 U 75-69-4 U 75-01-4 R 8 Recovery Lower Upper 100 85 115 104 72 119 103 81 120	127-18-4 U 108-88-3 U 156-60-5 U 10061-02-6 U 75-69-4 U 75-01-4 R 100 85 104 72 103 81	127-18-4 U 1.00 108-88-3 U 1.00 156-60-5 U 1.00 10061-02-6 U 1.00 75-69-4 U 1.00 75-01-4 TK 1.00 8 Recovery Lower Upper 100 85 115 104 72 119 103 81 120	127-18-4 U 1.00 108-88-3 U 1.00 156-60-5 U 1.00 10061-02-6 U 1.00 75-69-4 U 1.00 75-01-4 T 00 8 Recovery Lower Upper 104 72 119 103 81 120

U Undetected; the analyte was analyzed for, but not detected.

R Because of quality control deficiences for this analyte, this data may be rejected.

F The analyte was positively identified, but the quantitation was below the RL.

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~122132

Sample Number: L08060559-14PrePrep Method: NONEInstrument: HPMS14Client ID: GS9506WG1Prep Method: 5030BPrep Date: 06/25/2008 20:05Matrix: WaterAnalytical Method: 8260BCal Date: 06/24/2008 22:32Workgroup Number: WG275043Analyst: CMSRun Date: 06/25/2008 20:05Collect Date: 06/16/2008 13:40Dilution: 1File ID: 14M06614Sample Tag: 01Units: ug/L

Analyte	CAS. Number	Result Qua	1, RL	MDL
,1,1,2-Tetrachloroethane	630-20-6	U	0.500	0.250
,1,1-Trichloroethane	71-55-6	υ	1.00	0.250
1,2,2-Tetrachloroethane	79-34-5	U	0.500	0.125
,1,2-Trichloroethane	79-00-5	υ	1.00	0.250
,1-Dichloroethane	1 75-34-3	υ	1.00	0.125
,1-Dichloroethene	75-35-4	υ	1.00	0.500
,1-Dichloropropene	563-58-6	υ	1.00	0.250
,2,3-Trichlorobenzene	87-61-6	Ŭ	1.00	0.150
,2,3-Trichloropropane	96-18-4		1.00	0.500
,2,4-Trichlorobenzene	120-82-1	υ	1.00	0.200
,2,4-Trimethylbenzene	95-63-6	Ū	1.00	0.250
,2-Dichloroethane	107-96-2	บี้	0.500	0.250
,2-Dichlorobenzene	95-50-1	U	1.00	0.125
,2-Dibromo-3-chloropropane	96-12-8	υ	2.00	1.00
,2-Dichloropropane	78-87-5	U	1.00	0.200
,2-Dibromoethane	105-93-4	U U	1.00	0.250
	108-67-8	U U	1.00	0.250
.,3,5-Trimethylbenzene	541-73-1	U U	1.00	0.250
.,3-Dichlorobenzene	142-28-9	U U U	0.400	0.200
.,3-Dichloropropane	142-28-9		0.500	0.125
.,4-Dichlorobenzene		U U	100	50.0
,4-Dioxane	123-91-1	บ บ	1.00	0.125
-Chlorohexane	544-10-5	way,		0.125
2,2-Dichloropropane	594-20-7	U	1.00	2. An and the second
2-Chlorotoluene	95-49-8	U		0.125
I-Chlorotoluene	106-43-4	U	where is an experiment of each of the second	: 0.250
lcetone	: 67 - 64 - 1	v.	10.0	2.50
Benzene	71-43-2	U		0.125
Bromobenzene	108-86-1	U		0.125
Bromochloromethane	74-97-5	U		0.200
Bromodichloromethane	75-27-4	ឋ		0.250
Bromoform	75-25-2	Ŭ		0.500
Bromomethane	74-83-9	ע		0.500
Tarbon tetrachloride	56-23-5	U	and a set of the set o	0.250
Chlorobenzene	108-90-7	ប		0.125
Chloroethane	75-00-3	U		0.500
Chloroform	67-66-3	ប		0.125
Chloromethane	74-87-3	ប	1.00	0.250
cis-1,2-Dichloroethene	156-59-2	ΰ	1.00	0.250
cis-1,3-Dichloropropene	10061-01-5	υ	0.500	0.250
Dibromochloromethane	124-48-1	υ	0.500	0.250
Dibromomethane	74-95-3	U	1.00	0.250
Dichlorodifluoromethane	75-71-8	R	- () 1.00	0.250
Ethylbenzene	100-41-4	U	1.00	0.250
Mexachlorobutadiene	87-68-3	U	0.600	0.250
Isopropylbenzene	98-82-8	U		0.250
Methylene chloride	75-09-2	U		0.250
Methyl t-butyl ether (MTBE)	1634-04-4	~- R	a substances and a second s	0.500
MEK (2-Butanone)	78-93-3	1		2.50
(IBK (methyl isobutyl ketone)	108-10-1	Ŭ		2.50
n-Butylbenzene	104-51-8			0.250
n-Propylbenzene	103-65-1			0.125
	136777-61-2		and a second	0.500
n-,p-Xylene	91-20-3			0.200
Naphthalene	91-20-3			0.250
o-Xylene	99-87-6		and the first of the second	0.250
p-Isopropyltoluene		1		0.250
sec-Butylbenzene	135-98-8			
Styrene	100-42-5	Ţ		0.125
Trichloroethene	79-01-6			
tert-Butylbenzene	98-06-6	ι	1.00	0.250

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1128108

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PrePrep Method:NONE	Instrument: HPMS14
Prep Method: 5030B	Prep Date:06/25/2008 20:05
Analytical Method:8260B	Cal Date:06/24/2008 22:32
Analyst:CMS	Run Date:06/25/2008 20:05
Dilution:1	File ID:14M06614
Units:ug/L	
	Prep Method: 5030B Analytical Method: 8260B Analyst: CMS Dilution: 1 Units: ug/L

Analyte	CAS. Numbe	r R	esult	Qual		RL	MDL
Tetrachloroethene	127-18-4		27.2			1.00	0.250
Toluene	108-88-3	:		U	1	1.00	0.250
trans-1,2-Dichloroethene	156-60-5			U		1.00	0.250
trans-1,3-Dichloropropene	10061-02-6	5		U	1	1.00	0.500
Trichlorofluoromethane	75-69-4			U		1.00	0.250
Vinyl chloride	75-01-4			-R- \	<i>(</i>)	1.00	0.250
Surrogate	% Recovery	Lower	Upp)er	Qual		
Dibromofluoromethane	101	85	: 11	.5			
1,2-Dichloroethane-d4	107	72	11	9			
Toluene-d8	102	81	12	0			
4-Bromofluorobenzene	103	76	11	.9		-	

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U Undetected; the analyte was analyzed for, but not detected. R Because of quality control deficiences for this analyte, this data may be rejected.

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7128/08 OL

Report Number: L08060559

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Report Date : July 23, 2008

Sample Number:L08060559-15	PrePrep Method:N	ONE	Instrument:HPMS14				
Client ID:URS-2DWG1	Prep Method:50)30B	Prep Date:06/25/2008 20:37				
Matrix:Water	Analytical Method:8	2608	Cal	Date:06/24/3	2008 22:32		
kgroup Number:WG275043	Analyst:C	1S	Run	Date:06/25/	2008 20:37		
Collect Date:06/16/2008 16:05	Dilution:1		File 1	ID:14M06615			
Sample Tag:01	Units:ug						
			···· į · · · · · · · · · · · · · · · ·				
λnalyte	CAS. Number	Result	Qual	RL	MDL		
1,1,1,2-Tetrachloroethane	630-20-6		U	0.500	0.250		
1,1,1-Trichloroethane	71-55-6	1 	U	1.00	0.250		
1,1,2,2-Tetrachloroethane	79-34-5		U	0.500	0.125		
1,1,2-Trichloroethane	79-00-5	79-00-5		1.00	0.250		
1,1-Dichloroethane	75-34-3	0.239		1.00	0.125		
1,1-Dichloroethene	75-35-4	1	U	1.00	0.500		
1,1-Dichloropropene	563-58-6		υ	1.00	0.250		
1,2,3-Trichlorobenzene	87-61-6	1	U	1.00	0.150		
1,2,3-Trichloropropane	96-18-4	· · · · · · · · · · · · · · · · · · ·	U	1.00	0.500		
1,2,4-Trichlorobenzene	120-82-1	:	υ	1.00	0.200		
1,2,4-Trimethylbenzene	95-63-6	1	U	1.00	0.250		
1,2-Dichloroethane	107-06-2		υ	0.500	0.250		
1,2-Dichlorobenzene	95-50-1	}	υ	1.00	0.125		
1,2-Dibromo-3-chloropropane	96-12-8		υ	2.00	1.00		
1,2-Dichloropropane	78-87-5		U	1.00	0.200		
1,2-Dibromoethane	106-93-4	and a set of the set o		1.00	0.250		

95-63-6				0.250
107-06-2		υ	· · · · · · · · · · · · · · · · · · ·	0.250
95-50-1		υί	1.00	0.125
96-12-8		U	2.00	1.00
78-87-5		U	1.00	0.200
106-93-4		U	1.00	0.250
108-67-8		Ŭ	1.00	0.250
541-73-1		U	1.00	0.250
142-28-9		U	0.400	0.200
106-46-7	,	υ	0.500	0.125
123-91-1		U	100	50.0
544-10-5		U	1.00	0.125
594-20-7		υ	1.00	0.250
95-49-8		U	1.00	0.125
106-43-4	and an oral of the second second	U	1.00	0.250
67-64-1		U.	10.0	2.50
71-43-2		U	0.400	0.125
108-86-1		U	1.00	0.125
74-97-5	······	U	1.00	0.200
75-27-4		U	0.500	0.250
75-25-2		U	1.00	0.500
74-83-9		ប	3.00	0.500
56-23-5		υ	1.00	0.250
108-90-7	;;	U	0.500	0.125
75-00-3		ប	1.00	0.500
67-66-3		U	0.300	0.125
74-87-3	······································	U	1.00	0.250
156-59-2	71.9	1	1.00	0.250
10061-01-5		υ	0.500	0.250
124-48-1	1	Ŭ	0.500	0.250
74-95-3		U	1.00	0.250
75-71-8			1.00	0.250
100-41-4		U	1.00	0.250
87-68-3	· ···· · ···· ··· ··· · · · · ·	U	0.600	0.250
98-82-8	•	U	1.00	0.250
75-09-2		U	1.00	0.250
1634-04-4		-R 1	5.00	0.500
78-93-3		ប	10.0	2.50
108-10-1		U	10.0	2.50
		U	1.00	0.250
103-65-1		U	1.00	0.125
136777-61-2		U	2.00	0.500
91-20-3	1	U	1.00	0.200
95-47-6		U	1.00	0.250
99-87-6		ប	1.00	0.250
135-98-8	A consideration of the constant	U	1.00	0.250
100-42-5	1	U		0.125
		U	1.00	0.250
	$\begin{array}{c} 95-50-1\\ 96-12-8\\ 78-87-5\\ 106-93-4\\ 108-67-8\\ 541-73-1\\ 142-28-9\\ 106-46-7\\ 123-91-1\\ 544-10-5\\ 594-20-7\\ 95-49-8\\ 106-43-4\\ 67-64-1\\ 71-43-2\\ 108-86-1\\ 74-97-5\\ 75-27-4\\ 75-25-2\\ 75-27-4\\ 75-25-2\\ 74-83-9\\ 56-23-5\\ 108-90-7\\ 75-00-3\\ 67-66-3\\ 74-87-3\\ 155-59-2\\ 1006-10-5\\ 124-48-1\\ 74-95-3\\ 75-71-8\\ 100-15\\ 124-48-1\\ 74-95-3\\ 75-02-2\\ 1634-04-4\\ 78-93-3\\ 108-10-1\\ 104-51-8\\ 99-87-6\\ 135-98-8\\ \end{array}$	$\begin{array}{c ccccccccccccccccccccccccccccccccccc$	$\begin{array}{c c c c c c c c c c c c c c c c c c c $	$\begin{array}{ c c c c c c c c c c c c c c c c c c c$

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717×188

Report Number: L08060559

Report Date : July 23, 2008

Sample Number:L08060559-15	PrePrep Method:NONE	Instrument: HPMS14
Client ID: URS-2DWG1	Prep Method:5030B	Prep Date:06/25/2008 20:37
Matrix:Water	Analytical Method:8260B	Cal Date:06/24/2008 22:32
Workgroup Number: WG275043	Analyst:CMS	Run Date:06/25/2008 20:37
Collect Date:06/16/2008 16:05	Dilution:1	File ID:14M06615
Sample Tag:01	Units:ug/L	
analute	CaS Number Result	ດາາລີ ຊີໄ. ທີ່ການ

Analyte	CAS. Number	R	lesult	Qual]	RL	MDL
Tetrachloroethene	127-18-4	· · · · · ·		ΰ		.00	0.250
Toluene	108-88-3			U	[-00	0.250
trans-1,2-Dichloroethene	156-60-5			U			0.250
trans-1,3-Dichloropropene	10061-02-6			U	1	.00	0.500
Trichlorofluoromethane	75-69-4			U		00	0.250
Vinyl chloride	75-01-4		0.354	-R	<u> </u>	1.00	0.250
Surrogate	% Recovery	Lower	Uppe	r j	Qual		
Dibromofluoromethane	101	85	115				
1,2-Dichloroethane-d4	105	72	. 119				
Toluene-d8	103	81	120				
4-Bromofluorobenzene	103	76	119		1		

U Undetected; the analyte was analyzed for, but not detected.
R Because of quality control deficiences for this analyte, this data may be rejected.
F The analyte was positively identified, but the quantitation was below the RL.

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Vildio

1178158

Report Number: L08060559 Report Date :July 23, 2008

Styrene

sec-Butylbenzene

tert-Butylbenzene

Trichloroethene

p-Isopropyltoluene

Sample Number: L08060559-16	PrePrep Method:NONE			Lostrument: HPMS14					
Client ID: URS-2SWG1	Prep Method: 5030B			Prep Date:06/25/2008 21:10					
Matrix:Water	Analytical Method:8260B			Date:06/24,	2008 22:32				
orkgroup Number:WG275043	Analyst:CNS			Date:06/25	2008 21:10				
Collect Date:06/16/2008 18:10	Dilution:1			D:14M06616					
Sample Tag:01	Units:u	a/L							
		<u></u>	• •						
	····				······				
Analyte	CAS. Number	Result	Qual	RL	MDL				
1,1,1,2-Tetrachloroethane	630-20-6	<u>.</u>	U	0.500	0.250				
1,1,1-Trichloroethane	71-55-6	2.20		1.00	0.250				
1,1,2,2-Tetrachloroethane	79-34-5		U	0.500	0.125				
1,1,2-Trichloroethane	79-00-5		U	1.00	0.250				
1,1-Dichloroethane	75-34-3	0.569)	1.00	0.125				
1,1-Dichloroethene	75-35-4		υΨ	1.00	0.500				
1,1-Dichloropropene	563-58-6		U	1.00	0.250				
1,2,3-Trichlorobenzene	87-61-6	-	U	1.00	0.150				
1,2,3-Trichloropropane	96-18-4		U	1.00	0.500				
1,2,4-Trichlorobenzene	120-82-1		U	1.00	0.200				
1,2,4-Trimethylbenzene	95-63-6		ប	1.00	0.250				
1,2-Dichloroethane	107-06-2		υ	0.500	0.250				
1,2-Dichlorobenzene	95-50-1		σ	1.00	0.125				
1,2-Dibromo-3-chloropropane	96-12-8		U	2.00	1.00				
1,2-Dichloropropane	78-87-5		U	1.00	0.200				
1,2-Dibromoethane	106-93-4	· · · · · · · · · · · · · · ·	σ	1.00	0.250				
1,3,5-Trimethylbenzene	108-67-8	1	U	1.00	0.250				
1.3-Dichlorobenzene	541-73-1		U	1.00	0.250				
1,3-Dichloropropane	142-28-9		v	0.400	0.200				
1,4-Dichlorobenzene	106-46-7		Ū	0.500	0.125				
1,4-Dioxane	123-91-1	••••••••••••	Ŭ	100	50.0				
1-Chlorohexane	544-10-5		U	1.00	0.125				
2,2-Dichloropropane	594-20-7		u u	1.00	0.250				
2-Chlorotoluene	95-49-8		U .	1.00	0.125				
4-Chlorotoluene	106-43-4		U	1.00	0.250				
Provide and the second se	67-64-1		บ	10.0	2.50				
Acetone	71-43-2		U U	0.400	0.125				
Benzene	108-86-1		υ.	1.00	0.125				
Bromobenzene	74-97-5		υ	1.00	0.200				
Bromochloromethane Bromodichloromethane	74-57-5		U U	0.500	0.200				
Bromodichioromethane	75-25-2		Ū	1.00	0.500				
2	74-83-9		U	3.00	0.500				
Bromomethane			- U						
Carbon tetrachloride	56-23-5			1.00	0.250				
Chlorobenzene	108-90-7		U	0.500	0.125				
Chloroethane	75-00-3		ប	1.00	0.500				
Chloroform	67-66-3	0.204		0.300	0.125				
Chloromethane	74-87-3		U	1.00	0.250				
cis-1,2-Dichloroethene	156-59-2	0.995	Ŧ.) 1.00	0.250				
cis-1,3-Dichloropropene	10061-01-5		ប	0.500	0.250				
Dibromochloromethane	124-48-1		ប	0.500	0.250				
Dibromomethane	74-95-3		U	1.00	0.250				
Dichlorodifluoromethane	75-71-8		-* U	J 1.00	0.250				
Ethylbenzene	100-41-4		U	1.00	0.250				
Hexachlorobutadiene	87-68-3		U	0.600	0.250				
Isopropylbenzene	98-82-8		U	1.00	0.250				
Methylene chloride	75-09-2		U	1.00	0.250				
Methyl t-butyl ether (MTBE)	1634-04-4		R U) 5.00	0.500				
MEK (2-Butanone)	78-93-3		U	10.0	2.50				
MIBK (methyl isobutyl ketone)	108-10-1		U	10.0	2.50				
n-Butylbenzene	104-51-8		U	1.00	0.250				
n-Propylbenzene	103-65-1		U	1.00	0.125				
m-,p-Xylene	136777-61-2		υ	2.00	0.500				
Naphthalene	91-20-3		U	1.00	0.200				
o-Xylene	95-47-6		U	1.00	0.250				
n-Tronwonviltoluono	00-07-5			1 00	0.250				

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99-87-6

135-98-8

100-42-5

79-01-6

98-06-6

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Micropac Laboratories inc.

Report Number: L08060559 Report Date : July 23, 2008

rep Method: 5030B	Prep Date:06/25/2008 21:10
cal Method:8260B	Cal Date:06/24/2008 22:32
Analyst:CMS	Run Date:06/25/2008 21:10
Dilution:1	File ID:14M06616
Units:ug/L	
_	Analyst:CMS Dilution:1

CAS. Number	:	Result	Qual	RL	MDL
127-18-4			υ	1.00	0.250
108-88-3			U	1.00	0.250
156-60-5			U	1.00	0.250
10061-02-6			ប	1.00	0.500
75-69-4			υ	1.00	0.250
75-01-4			-R- 1	/j 1.00	0.250
% Recovery	Lower	ប្រ	per	Qual	
102	85	1	15		
108	72	1	19	1	
102	81	1	20		
103	76	1	19		
	127-18-4 108-88-3 156-60-5 10061-02-6 75-69-4 75-01-4 % Recovery 102 108 102	108-88-3 156-60-5 10061-02-6 75-69-4 75-01-4 % Recovery 102 108 108 102 81	127-18-4 108-88-3 156-60-5 10061-02-6 75-69-4 75-01-4 % Recovery 102 108 72 102 102 102	127-18-4 U 108-88-3 U 156-60-5 U 10061-02-6 U 75-69-4 U 75-01-4 -R-V 102 85 108 72 102 81	127-18-4 U 1.00 108-88-3 U 1.00 156-60-5 U 1.00 10061-02-6 U 1.00 75-69-4 U 1.00 75-01-4 -R-V 1.00 102 85 115 108 72 119 102 81 120

R Because of quality control deficiences for this analyte, this data may be rejected.
 U Undetected; the analyte was analyzed for, but not detected.
 F The analyte was positively identified, but the quantitation was below the RL.

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Microbac

7128/08 Pr

Sample Number: L08060559-17PrePrep Method: NONEInstrument: HPMS14Client ID: URS-28WG9Prep Method: 5030BPrep Date: 06/25/2008 21:42Matrix: WaterAnalytical Method: 8260BCal Date: 06/24/2008 22:32Workgroup Number: WG275043Analyst: CMSRun Date: 06/25/2008 21:42Collect Date: 06/16/2008 18:10Dilution: 1Pile ID: 14M06617Sample Tag: 01Units: ug/LVerter

Analyte	CAS. Number	Result	Qual	RL.	MDL
1,1,1,2-Tetrachloroethane	630-20-6	}	U	0.500	0.250
,1,1-Trichloroethane	71-55-6	2.25		1.00	0.250
., 1, 2, 2-Tetrachloroethane	79-34-5	1	U	0.500	0.125
,1,2-Trichloroethane	79-00-5		υ	1.00	0.250
,1-Dichloroethane	75-34-3	0.585	- R)	1.00	0.125
,1-Dichloroethene	75-35-4	et e ser acceletationer e	υΨ	1.00	0.500
,1-Dichloropropene	563-58-6		υ	1.00	0.250
,2,3-Trichlorobenzene	87-61-6		Ū	1.00	0.150
,2,3-Trichloropropane	96-18-4	· · · · · · · · · · · · · · · · · · ·	U	1.00	0.500
,2,4-Trichlorobenzene	120-82-1		U	1.00	0.200
1,2,4-Trimethylbenzene	95-63-6	. <u>}</u>		1.00	0.250
L,2-Dichloroethane	107-06-2		υ	0.500	0.250
1,2-Dichlorobenzene	95-50-1	÷	U	1.00	0.125
	95-30-1	4	U	2.00	1.00
,2-Dibromo-3-chloropropane			UUU		
1,2-Dichloropropane	78-87-5			1.00	0.200
,2-Dibromoethane	106-93-4		U	1.00	0.250
.,3,5-Trimethylbenzene	108-67-8		U	1.00	0.250
,3-Dichlorobenzene	541-73-1	<u>.</u>	U	1.00	0.250
,3-Dichloropropane	142-28-9		υ	0.400	0.200
,4-Dichlorobenzene	106-46-7		ប	0.500	0.125
.,4-Dioxane	123-91-1	•	U	100	50.0
-Chlorohexane	544-10-5	1	U	1.00	0.125
2,2-Dichloropropane	594-20-7	1	U	1.00	0.250
-Chlorotoluene	95-49-8	1	U I	1.00	0.125
i-Chlorotoluene	106-43-4		U	1.00	0.250
Acetone	67-64-1		U	10.0	2.50
Benzene	71-43-2		U	0.400	0.125
Bromobenzene	108-86-1	;	U	1.00	0.125
Bromochloromethane	74-97-5		U	1.00	0.200
Bromodichloromethane	75-27-4		U	0.500	0.250
Bromoform	75-25-2		U	1.00	0.500
Bromomethane	74-83-9		U	3.00	0.500
Carbon tetrachloride	56-23-5		Ŭ	1.00	0.250
Chlorobenzene	108-90-7		U	0.500	0.125
Chloroethane	75-00-3			1.00	0.500
Chloroform	67-66-3	0.197	-P 1	0.300	0.125
Chloromethane	74-87-3	0.197		1.00	0.123
	A CONTRACTOR OF A CONTRACTOR O		enter a construction of the second	the state of the s	
cis-1,2-Dichloroethene	156-59-2	0.966	- P	1.00	0.250
zis-1,3-Dichloropropene	10061-01-5		U	0.500	0.250
Dibromochloromethane	124-48-1	·	U	0.500	0.250
Dibromomethane	74-95-3	and the second	U	1.00	0.250
Dichlorodifluoromethane	75-71-8) 1.00	0.250
Sthylbenzene	100-41-4		ប	1.00	0.250
fexachlorobutadiene	87-68-3	1	U	0.600	0.250
[sopropylbenzene	98-82-8	1	U	1.00	0.250
fethylene chloride	75-09-2		υ	1.00	0.250
Methyl t-butyl ether (MTBE)	1634-04-4		-R- Ų) 5.00	0.500
MEK (2-Butanone)	78-93-3		U	10.0	2.50
(IBK (methyl isobutyl ketone)	108-10-1		ប ·	10.0	2.50
n-Butylbenzene	104-51-8		U	1.00	0.250
1-Propylbenzene	103-65-1	1	υ	1.00	0.125
c-,p-Xylene	136777-61-2		U	2.00	0.500
Vaphthalene	91-20-3		U	1.00	0.200
-Xylene	95-47-6		Ŭ	1.00	0.250
>-Isopropyltoluene	99-87-6		Ŭ	1.00	0.250
sec-Butylbenzene	135-98-8		<u> </u>	1.00	0.250
Styrene	100-42-5		ບ : ບ	1.00	0.125
Styrene Frichloroethene	79-01-6	2.22		1.00	0.125
tert-Butylbenzene	98-06-6	4.44	U	1.00	0.250

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7128/08

Report Number: L08060559

Report Date : July 23, 2008

Sample Number:L08060559-17	PrePrep	Method:NONE		Instrume	nt:HPMS14		
Client ID: URS-2SWG9	Prep	Method: 50301		Prep Da	te:06/25/200	8 21:42	
Matrix:Water	Analytical	Method:82601	8	Cal Da	te:06/24/200	8 22:32	
Workgroup Number:WG275043	,	Analyst:CMS		Run Da	te:06/25/200	8 21:42	
Collect Date:06/16/2008 18:10	D	ilution:1			14M06617		
Sample Tag:01		Units:ug/L					
Analvte	CAS. Nu	mber	Result	Qual	RL	MDL	

. Analyte	; CAS. NUEDEI		esuit	Quar		RL	MDL	
Tetrachloroethene	127-18-4	:		U		1.00	0.250	
Toluene	108-88-3			U		1.00	0.250	į
trans-1,2-Dichloroethene	156-60-5			U		1.00	0.250	
trans-1,3-Dichloropropene	10061-02-6			U		1.00	0.500	
Trichlorofluoromethane	75-69-4			U		2.00	0.250	
Vinyl chloride	75-01-4			-R- \	11	1.00	0.250	
Surrogate	% Recovery	Lower	Uppe	r	Qual			
Dibromofluoromethane	101	85	115	1		1		
1,2-Dichloroethane-d4	108	72	119					
Toluene-d8	102	81	120					
4-Bromofluorobenzene	104	76	119			j		

U Undetected; the analyte was analyzed for, but not detected. R Because of quality control deficiences for this analyte, this data may be rejected. F The analyte was positively identified, but the quantitation was below the RL.

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micropac Laporatories inc.

Report Number: L08060559 Report Date :July 23, 2008

 Sample Number: L08060559-18
 PrePrep Method: NONE
 Instrument: HPMS14

 Client ID: EB061708
 Prep Method: 5030B
 Prep Date: 06/25/2008 16:49

 Matrix: Water
 Analytical Method: 8260B
 Cal Date: 06/24/2008 22:32

 Workgroup Number: W6275043
 Analyst: CMS
 Run Date: 06/25/2008 15:49

 Collect Date: 06/17/2008 15:05
 Dilution: 1
 File ID: 14M06608

 Sample Tag: 01
 Units: ug/L
 File ID: 14M06608

Analyte	CAS. Number	Result	Qual	RL	MDL
,1,1,2-Tetrachloroethane	630-20-6		υ	0.500	0.250
,1,1-Trichloroethane	71-55-6		U	1.00	0.250
,1,2,2-Tetrachloroethane	79-34-5		U	0.500	0.125
,1,2-Trichloroethane	79-00-5		U	1.00	0.250
,1-Dichloroethane	75-34-3		σ	1.00	0.125
,1-Dichloroethene	75-35-4		U	1.00	0.500
L,1-Dichloropropene	563-58-6		U	1.00	0.250
L,2,3-Trichlorobenzene	87-61-6		υ	1.00	0.150
L,2,3-Trichloropropane	96-18-4	į	U	1.00	0.500
L,2,4-Trichlorobenzene	120-82-1		υ	1.00	0.200
1,2,4-Trimethylbenzene	95-63-6		U	1.00	0.250
L,2-Dichloroethane	107-06-2		U	0.500	0.250
.2-Dichlorobenzene	95-50-1		U	1.00	0.125
1,2-Dibromo-3-chloropropane	96-12-8		U	2.00	1.00
1,2-Dichloropropane	78-87-5		U .	1.00	0.200
1,2-Dibromoethane	106-93-4	•••••	U	1.00	0.250
1,3,5-Trimethylbenzene	108-67-8		U	1.00	0.250
1,3-Dichlorobenzene	541-73-1		υ	1.00	0.250
1,3-Dichloropropane	142-28-9		U U U	9.400	0.200
1, 4-Dichlorobenzene	105-46-7		υ	0.500	0.125
1,4-Dichiorobenzene	123-91-1		υ	100	50.0
l-Chlorohexane	544-10-5		υ υ	1.00	0.125
2,2-Dichloropropane	594-20-7		UUU	1.00	0.125
2.2-Dichioropropane 2-Chlorotoluene	95-49-8		U U	1.00	0.250
	and a design of the second				
4-Chlorotoluene	106-43-4			1.00	0.250
Acetone	67-64-1		<u> </u>	10.0	2.50
Benzene	71-43-2		ម	0.400	0.125
Bromobenzene	108-86-1		<u>v</u>	1.00	0.125
Bromochloromethane	74-97-5		υ	1.00	0.200
Bromodichloromethane	75-27-4		U	0.500	0.250
Bromoform	75-25-2		្រា	1.00	0.500
Bromomethane	74-83-9		U	3.00	0.500
Carbon tetrachloride	56-23-5		ប	1.00	0.250
Chlorobenzene	108-90-7		U	0.500	0.125
Chloroethane	75-00-3		U	1.00	0.500
Chloroform	67-66-3		ប	0.300	0.125
Chloromethane	74 - 87 - 3		U	1.00	0.250
cis-1,2-Dichloroethene	156-59-2		U	1.00	0.250
cis-1,3-Dichloropropene	10061-01-5		U	0.500	0.250
Dibromochloromethane	124-48-1		υ	0.500	0.250
Dibromomethane	74-95-3	1	U	1.00	0.250
Dichlorodifluoromethane	75-71-8	1	-R- U) 1.00	0.250
Ethylbenzene	100-41-4		U :	1.00	0.250
Hexachlorobutadiene	87-68-3		U	0.600	0.250
Isopropylbenzene	98-82-8		U	1.00	0.250
Methylene chloride	75-09-2		U	1.00	0.250
Methyl t-butyl ether (MTBE)	1634-04-4	· · · · · · · · · · · · · · · · · · ·	R- ()		0.500
MEK (2-Butanone)	78-93-3		υ	10.0	2.50
MIBK (methyl isobutyl ketone)	108-10-1		U	10.0	2.50
n-Butylbenzene	104-51-8		U	1.00	0.250
n-Propylbenzene	103-65-1	1	<u> </u>	1.00	0.125
a-,p-Xylene	136777-61-2		υ	2.00	0.500
Naphthalene	91-20-3		υ.	1.00	0.200
o-Xylene	95-47-6			1.00	0.200
	99-87-6		U	1.00	0.250
p-Isopropyltoluene sec-Butylbenzene	135-98-8		U	1.00	0.250
	135-98-8		UU		0.250
Styrene				1.00	
Trichloroethene	79-01-6	1	U	1.00	0.250

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7128/18

Sample Number: L08060559-18	PrePrep Method:NONE	Instrument: HPMS14
Client ID:EB061708	Prep Method:5030B	Prep Date:06/25/2008 16:49
Matrix:Water	Analytical Method:8260B	Cal Date:06/24/2008 22:32
Workgroup Number:WG275043	Analyst:CMS	Run Date:06/25/2008 16:49
Collect Date:06/17/2008 15:05	Dilution:1	File ID:14M06608
Sample Tag:01	Units:ug/L	

Analyte	CAS. Numbe	CAS. Number Result		Qual	R	L	MDL
Tetrachloroethene	127-18-4			U	1.	00	0.250
Toluene	108-88-3			U	1.	00	0,250
trans-1,2-Dichloroethene	156-60-5			υ	1.	00	0.250
trans-1,3-Dichloropropene	10061-02-	6		U	1.	00	0.500
Trichlorofluoromethane	75-69-4	75-69-4		U	1.	00	0.250
Vinyl chloride	75-01-4			-R	$\left(\right) $ 1.	00	0.250
Surrogate	% Recovery	Lower	Upp	er	Qual		
Dibromofluoromethane	99.0	85	11	5	,		
1,2-Dichloroethane-d4	102	72	11	9			
Toluene-d8	103	81	12	0			
4-Bromofluorobenzene	102	76	11	9			

U Undetected; the analyte was analyzed for, but not detected.

R Because of quality control deficiences for this analyte, this data may be rejected.

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Report Number: L08060559

Report Date :July 23, 2008

Sample Number:L08060559-02	PrePrep Method:NONE	Instrument: HPMS8
Client ID:TB061408	Prep Method: 5030B	Prep Date:06/24/2008 12:40
Matrix:Water	Analytical Method:8260B	Cal Date:04/29/2008 15:06
Workgroup Number:WG274878	Analyst:FJB	Run Date:06/24/2008 12:40
Collect Date:06/14/2008 00:01	' Dilution:1	File 1D:8M345906
Sample Tag:01	Units:ug/L	
Sample Tag:01	Units:ug/L	

Analyte	CAS. Number	Result	Qual	RL	MDL
,1,1,2-Tetrachloroethane	630-20-6		U	0.500	0.250
,l,l-Trichloroethane	71-55-6		U	1.00	0.250
,1,2,2-Tetrachloroethane	79-34-5		U	0.500	0.125
,1,2-Trichloroethane	79-00-5	1	U	1.00	0.250
,1-Dichloroethane	75-34-3	1	υ	1.00	0.125
,l-Dichloroethene	75-35-4		U	1.00	0.500
,1-Dichloropropene	563-58-6	1	U	1.00	0.250
,2,3-Trichlorobenzene	87-61-6	· · · · · · · · · · · · · · · · · ·	U	1.00	0.150
,2,3-Trichloropropane	96-18-4		U U	1.00	0.500
,2,4-Trichlorobenzene	120-82-1		υ	1.00	0.200
,2,4-Trimethylbenzene	95-63-6		T T	1.00	6.250
,2-Dichloroethane	107-06-2		υ	0.500	0.250
,2-Dichlorobenzene	95-50-1	eg and second and second		1.00	0.125
The second s	96-12-8		u	2.00	1.00
,2-Dibromo-3-chloropropane	78-87-5		U	1.00	0.200
,2-Dichloropropane	· · · · · · · · · · · · · · · · · · ·		U U		
,2-Dibromoethane	106-93-4		and a farmer for	1.00	0.250
,3,5-Trimethylbenzene	108-67-8		U	1.00	0.250
, 3-Dichlorobenzene	541-73-1	· · · · · · · · · · · · · · · · · ·	U	1.00	0.250
,3-Dichloropropane	142-28-9		U	0.400	0.200
,4-Dichlorobenzene	106-46-7		U	0.500	0.125
,4-Dioxane	i 123-91-1		-B K	100 بر	50.0
-Chlorohexane	544-10-5		U	1.00	0.125
,2-Dichloropropane	594-20-7		U	1.00	0.250
-Chlorotoluene	95-49-8		U	1.00	0.125
-Chlorotoluene	106-43-4		U	1.00	0.250
cetone	67-64-1		U :	10.0	2.50
enzene	71-43-2		υ,	0.400	0.125
romobenzene	108-86-1	-	U I	1.00	0.125
romochloromethane	74-97-5		υ	1.00	0.200
romodichloromethane	75-27-4	•	U	0.500	0.250
Fromoform	75-25-2		υ :	1.00	0.500
Bromomethane	74-83-9	••• •• •• •• ••	U	3.00	0.500
Carbon tetrachloride	56-23-5		ប	1.00	0.250
hlorobenzene	108-90-7		U	0.500	0.125
Chloroethane	75-00-3		- l - ž	1.00	0.500
Thloroform	67-66-3	:	U	0.300	0.125
chloromethane	74-87-3	· • · · · · · · · · · · · · · · · · · ·	U U	1.00	0,250
sis-1,2-Dichloroethene	156-59-2		υ	1.00	0.250
is-1,3-Dichloropropene	10061-01-5		- U	0.500	0.250
bibromochloromethane	124-48-1		τ	0.500	0.250
)ibromochioromethane	74-95-3		υ	1.00	0.250
)ibromomethane Dichlorodifluoromethane	74-95-3			1.00	0.250
and a construction of the		· · · · · · · · · · · · · · · · · · ·	- <u>* U.</u>	1.00	0.250
Cthylbenzene	100-41-4	··•••·································	U U		
Mexachlorobutadiene	87-68-3		· · · · · · · · · · · · · · · · · · ·	0.600	0.250
Isopropylbenzene	98-82-8		U	1.00	0.250
fethylene chloride	75-09-2		U	1.00	0.250
fethyl t-butyl ether (MTBE)	1634-04-4		ប	5.00	0.500
IEK (2-Butanone)	78-93-3	; 	U	10.0	2.50
IIBK (methyl isobutyl ketone)	108-10-1		υ	10.0	2.50
-Butylbenzene	104-51-8		U	1.00	0.250
-Propylbenzene	103-65-1		υ,	1.00	0.125
1-,p-Xylene	136777-61-2		U	2.00	0.500
Taphthalene	91-20-3		U	1.00	0.200
p-Xylene	95-47-6		U	1.00	0.250
-Isopropyltoluene	99-87-6		υ	1.00	0.250
sec-Butylbenzene	135-98-8		U	1.00	0.250
Styrene	100-42-5		U	1.00	0.125
Frichloroethene	79-01-6	•	υ	1.00	0.250
tert-Butylbenzene	98-06-6		υ	1.00	0.250

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Report Number: L08060559

Report Date : July 23, 2008

Sample Number: L08060559-02	PrePrep Method:NONE	Instrument: HPMS8
Client ID:TB061408	Prep Method: 5030B	Prep Date:06/24/2008 12:40
Matrix:Water	Analytical Method:8260B	Cal Date:04/29/2008 15:06
Workgroup Number: WG274878	Analyst:FJB	Run Date:06/24/2008 12:40
Collect Date:06/14/2008 00:01	Dilution:1	File ID:8M345906
Sample Tag:01	Units:ug/f	

Analyte	CAS. Number	· I	Result	Qual	-	RL	MDL
Tetrachloroethene	127-18-4			U	-	1.00	0.250
Toluene	108-88-3			U		1.00	0.250
trans-1,2-Dichloroethene	156-60-5			U		1.00	0.250
trans-1,3-Dichloropropene	10061-02-6			υ		1.00	0.500
Trichlorofluoromethane	75-69-4			U		1.00	0.250
Vinvl chloride	75-01-4			U		1.00	0.250
Surrogate	% Recovery	Lower	Upper		Qual		
Dibromofluoromethane	106	85	115				
1,2-Dichloroethane-d4	101	72	119				
Toluene-d8	93.5	81	120				
4-Bromofluorobenzene	93.7	76	119			:	

•

U Undetected; the analyte was analyzed for, but not detected. R Because of quality control deficiences for this analyte, this data may be rejected.

1/2-1/04

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of 36

4

Sample Number: L08060559-02	PrePrep Method: NONE	Instrument: HPMS14
Client ID:TB061408	Prep Method: 5030B	Prep Date:06/27/2008 10:51
Matrix:Water	Analytical Method:8260B	Cal Date:06/26/2008 20:24
Workgroup Number:WG275287	Analyst: CMS	Run Date:06/27/2008 10:51
Collect Date:06/14/2008 00:01	Dilution:1	File ID:14M06650
Sample Tag:01	Units:ug/L	
• 7		• • •

1.4-Dioxane	CAS, Numb	1		Qual U	RL 2.00	MDL. 1.00
Surrogate	% Recovery	Lower	Upper	Qu	al	
Dibromofluoromethane	115	54	138	1		
1,2-Dichloroethane-d4	114	51	135			

U Undetected; the analyte was analyzed for, but not detected.

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1/2/08

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1.0 INTRODUCTION

This data quality review pertains to groundwater samples collected in November 2008 at Air Force Plant 59 (AFP-59). Parameters evaluated in groundwater samples included the total concentration of volatile organic constituent (VOC). The samples were analyzed by Microbac Services, Marietta, Ohio.

Data quality review is an after-the-fact technical review of analytical data whereby the quality and usability of the data are determined based on a set of predefined criteria. These criteria depend upon the type of data involved and the purpose for which those data were collected. Data quality review assesses whether and to what extent specified criteria were met, and places restrictions on data use based on quality parameters. The data quality review process can range from a cursory review used to detect out-of-control situations to a detailed evaluation, depending on the analytical protocol, the associated quality control samples collected, and the intended data use.

Specific criteria for data quality review may include, but are not limited to: technical holding times, analysis of blanks, surrogate spike recovery, analysis of duplicates, and reported practical quantitation limits (PQLs). Where applicable, the recommendations of USEPA SW-846 Test Methods for Evaluating Solid Waste (Third Edition, December 1996) or USEPA Methods for Chemical Analysis of Water and Wastes (Revised March 1983) analytical method requirements, USEPA CLP National Functional Guidelines for Organic and Inorganic Data Review (February 1994, Functional Guidelines) data review guidance, and professional judgment.

Table 1 presents the data qualifiers applied during this review effort and their meanings.

Qualifier	Description
J	This is an estimated value.
UJ	The analyte was analyzed for but was not detected. The associated value is an estimate and may be inaccurate or imprecise.
М	Matrix effect: The concentration is estimated due to a matrix effect.
R	One or more quality control criteria (for example, LCS recovery, surrogate spike recovery) failed.
U	The analyte was analyzed for, but not detected. The associated numerical value is at or below the MDL.

Table 1 Data Qualifiers

Table 2 provides a cross-reference list for field sample IDs and lab sample IDs.

Field Sample ID	Lab Sample ID	Field Sample ID	Lab Sample ID
59SW4WG1	L08110347-01	59SW3WG1	L08110347-05
59SW7WG1	L08110347-02	59DW3WG1	L08110347-06
59SW7WG1-MS	L08110347-03	59DW3WG9	L08110347-07
59SW7WG1-MSD	L08110347-04		

Table 2 Field Sample ID/Lab Sample ID Cross Reference

During the data quality review process, laboratory qualified and unqualified data are verified against all available supporting documentation. Based on this review, qualifier codes may be added, deleted, or modified by the validator. Final results are therefore either qualified or unqualified. (Note: In those cases where the laboratory added a "U" flag indicated a non-detect result, and the validator agrees with this flag, then it remains intact, as noted on the corresponding Results Summary Form.) Changes to the data are reflected on the Results Summary Forms in Appendix A.

2.0 VOLATILE ORGANIC CONSTITUENTS

Volatile organic constituents were analyzed using EPA Test Method for Evaluating Solid Waste (SW-846) Method 8260B. Samples were analyzed in one batch:

WG288769: 59SW4WG1, 59SW3WG1, 59DW3WG1, 59DW3WG9

WG288953: 59SW7WG1, 59SW7WG1-MS, 59SW7WG1-MSD

2.1 Holding Times

All samples were analyzed within prescribed hold times. No qualification is needed.

2.2 Calibration

<u>WG288769</u>: The Initial Calibration for this batch was performed on 11/22/2008. Standards were analyzed at 0.30, 0.4, 1, 2, 5, 20, 50, 100, and 200 ug/L in support of the Initial Calibration. The %RSD values were less than 30% for all target constituents. No qualification is needed.

The %D for the second source calibration was within $\pm 25\%$ for target constituents except for acetone. The validator qualifies UJ or J, the non detect and positive results, respectively, in the acetone results in the associated samples. The validator removes the "R" flag assigned by the laboratory to denote anomalies.

For the associated continuing calibration standard, none of the associated %D values were greater than 20% for target constituents except for dichlorodifluoromethane. The validator qualifies UJ or J, the non detect and positive results, respectively, in the dichlorodifluoromethane results in the associated samples. The validator removes the "R" flag assigned by the laboratory to denote anomalies.

WG288953: The Initial Calibration for this batch was performed on 11/24/2008. Standards were analyzed at 0.30, 0.4, 1, 2, 5, 20, 50, 100, and 200 ug/L in support of the Initial Calibration. The %RSD values were less than 30% for all target constituents. No qualification is needed.

The %D for the second source calibration was within $\pm 25\%$ for target constituents except for acetone. The validator qualifies UJ or J, the non detect and positive results, respectively, in the acetone results in the associated samples. The validator removes the "R" flag assigned by the laboratory to denote anomalies.

For the associated continuing calibration standard, none of the associated %D values were greater than 20% for target constituents except for dibromofluoromethane. The validator qualifies UJ or J, the non detect and positive results, respectively, in the dibromofluoromethane results in the associated samples.

For all samples, it is noted that for those results which were less than the RL but greater than the MDL, the laboratory assigned an "F" flag, indicating an estimated value. Unless qualified otherwise, the validator removes the F flag and replaces it with the "J" qualifier, indicating an estimated value.

2.3 Laboratory Control Samples

WG288769: Laboratory control sample exhibited a 155% recovery for acetone, which are above the LCS limits and a 37.5% for dichlorodifluoromethane which are below the LCS limits. The validator

qualifies J the acetone positive results in the associated samples and qualifies R the dichlorodifluoromethane results in the associated samples.

<u>WG288953</u>: Laboratory control sample exhibited a 159% recovery for acetone, which are above the LCS limits. The validator qualifies J the acetone positive results in the associated samples

2.4 Blanks

WG288769: No constituents were detected in the associated method blank. No qualification needed.

WG288953: No constituents were detected in the associated method blank. No qualification needed.

Note: A trip blank was not put into the cooler prior to shipping samples.

2.5 Matrix Spike/Matrix Spike Duplicate

<u>WG288769:</u> Matrix Spike/Matrix Spike duplicate samples were not included in this sample delivery group. No qualification is needed.

<u>WG288953</u>: Sample 59SW7WG1 served as the MS/MSD. Recoveries of 1,4-dioxane and acetone exceeded the %RPD and acetone, chloromethane, and vinyl chloride exceeded the upper control limits of MS/MSD samples. As a result, 1,4-dioxane, acetone, chloromethane, and vinyl chloride for sample 59SW7WG1 were qualified M. No further qualification was necessary.

2.6 Surrogate Recovery

<u>WG288769:</u> All surrogate recoveries were within control limits for all environmental and quality control samples. No qualification is needed.

<u>WG288953:</u> All surrogate recoveries were within control limits for all environmental and quality control samples. No qualification is needed.

2.7 Internal Standards

All internal standard area counts were within control limits for all samples. No qualification is needed based on the internal standard information provided.

2.8 Duplicates

Sample 59DW3were analyzed in duplicate. One of two criteria was followed when evaluating field duplicates, depending on the amount detected. If the amount detected was greater than five times the reporting limit (RL), then the relative percent difference (RPD) should have been less than 25 percent. If the amount was less than five times the RL, then the difference between the duplicate and the sample concentrations should have been less than the RL. Agreement is excellent and no qualification is needed.

Table 3: Duplicate Comparison (µg/L)

Analyte	Reporting Limit (RL)	5DW3WG1	59DW3WG9	Relative Percent Difference (RPD)
Cis-1,2-	1.0	65.2	ND	%

dichloroethene				
Trans-1,2-	1.0	1.09	1.18	7.9%
dichloroethene				

2.8 Summary

The data are acceptable with validator-assigned qualifiers. Trichlorofluoromethane in sample 01 may have attributed to carry-over contamination from a previous analysis.

LABORATORY REPORT

L08110347

12/09/08 11:12

Submitted By

Microbac Laboratories Inc. 158 Starlite Drive Marietta, OH 45750 (740) 373-4071

For

Account Name: Earth Tech, Inc 675 North Washington Street Suite 300 Alexandria, VA 22314 Attention: Devon Chicoine

Project Number: 2551.030 Project: Alex-Air Force Plant 59 Site: AFP 59

Sample Analysis Summary

Client ID	Lab ID	Method	Dilution	Date Received
595W4WG1	L08110347-01	8260B	1	13-NOV-08
59sw7wG1	L08110347-02	8260B	1	13-NOV-08
59SW7WG1-MS	L08110347-03	8260B	1	13-NOV-08
59SW7WG1-MSD	L08110347-04	8260B	1	13-NOV-08
59SW3WG1	L08110347-05	8260B	1	13-NOV-08
59DW3WG1	L08110347-06	8260B	1	13-NOV-08
59DW3WG9	L08110347-07	8260B	1	13-NOV-08

L1_A_PROD - Modified 03/05/2008 PDF File ID:1271360 Report generated: 12/09/2008 11:12 1 OF 1

Microbac

ç . Report Number: L08110347 Report Date :December 9, 2008

Sample Number:L08110347-01	PrePrep Mel		· ·		nt:HPMS		19.37	
Client ID:59SW4WG1		:hod:5030B	4		te:11/2			
Matrix:Water	Analytical Me	and the second second			te:11/1 te:11/2			
kgroup Number:WG288769		lyst:MES					10:37	
Collect Date:11/11/2008 14:16		ion:1 hits:ug/L	1	TTG IN	8M34995	4		
Sample Tag:01	UI UI	lics:ug/L						
		. . . <u>.</u>	- 1-	-				
Analyte	CAS. Numbe	r Res	ult 🛛	ual	RL 0.500		MDL 0.250	
1,1,1,2-Tetrachloroethane	630-20-6		513	v F J	1.00	-	0.250	-
1,1,1-Trichloroethane	71-55-6		513		0.500		0.125	
1,1,2,2-Tetrachloroethane 1,1,2-Trichloroethane	79-34-5			υ υ	1.00		0.250	
1,1,2-Trichloroethane 1,1-Dichloroethane	75-34-3		825		1.00		0.125	
1,1-Dichloroethene	75-35-4			T U	1.00		0.500	-
1,1-Dichloropropene	563-58-6	· · · · · · · · ·		Ū .	1.00		0.250	
1,2,3-Trichlorobenzene	87-61-6	1. 1. A. 1. 1. A. A. 1. 1. A. A. 1. 1. A.		U	1.00		0.150	
1,2,3-Trichloropropane	96-18-4			υ	1.00		0.500	
1,2,4-Trichlorobenzene	120-82-1			υ	1.00		0.200	
1,2,4-Trimethylbenzene	95-63-6			U.	1.00		0.250	
1,2-Dichloroethane	107-06-2			บ	0.500		0.250	
1,2-Dichlorobenzene	95-50-1		1	υ	1.00		0.125	
1,2-Dibromo-3-chloropropane	96-12-8			ບຼ	2.00		1.00	
1.2-Dichloropropane	78-87-5			U	1.00		0.200	
1,2-Dibromoethane	106-93-4			σ	1.00		0.250	
1,3,5-Trimethylbenzene	108-67-8			U T	1.00		0.250	
1.3-Dichlorobenzene	541-73-1			U H	1.00		0.250	
1,3-Dichloropropane	142-28-9		[บ บ	0.400 0.500		0.200	
1,4-Dichlorobenzene	106-46-7			υ υ	100		50.0	
1,4-Dioxane	123-91-1 544-10-5			UUU	1.00		0.125	
1-Chlorohexane	594-20-7		ţ	U U	1.00		0.250	
2,2-Dichloropropane 2-Chlorotoluene	95-49-8			ີບ	1.00		0.125	
4-Chlorotoluene	106-43-4			ບັ	1.00		0.250	
Acetone	67-64-1	•	I	-R- (2)	10.0		2.50	
Benzene	71-43-2	<i></i>	· •	0 0.0	0.400		0.125	
Bromobenzene	108-86-1			U	1.00		0.125	
Bromochloromethane	74-97-5			υ	1.00		0.200	
Bromodichloromethane	75-27-4			ប	0.500		0.250	
Bromoform	75-25-2			υ	1.00		0.500	
Bromomethane	74-83-9			U	3.00		0.500	
Carbon tetrachloride	56-23-5			U	1.00		0.250	
Chlorobenzene	108-90-7			U.	0.500		0.125	
Chloroethane	75-00-3			U.	1.00		0.500	
Chloroform	67-66-3			U .	0.300 1.00		0.125	
Chloromethane	74-87-3	· · · .	38	U.	1.00		0.250	
cis-1,2-Dichloroethene	10061-01-		°L	σ	0.500		0.250	
cis-1,3-Dichloropropene Dibromochloromethane	124-48-1			JU -	0.500		0.250	
Dibromochibromethane	74-95-3			U UU	1.00		0.250	
Dichlorodifluoromethane	75-71-8	· · · · · · · · · · · · · · · · · · ·		R	1.00		0.250	
Ethylbenzene	100-41-4		• •	υ	1.00		0.250	
Hexachlorobutadiene	87-68-3		1	υ	0.600		0.250	
Isopropylbenzene	98-82-8			υ	1.00		0.250	
Methylene chloride	75-09-2			ບໍ່	1.00		0.250	
Methyl t-butyl ether (MTBE)	1634-04-4	E		ບຸ	5.00		0.500	
MEK (2-Butanone)	78-93-3			ບໍ່	10.0		2.50	
MIBK (methyl isobutyl ketone)	108-10-1			U.	10.0		2,50	
n-Butylbenzene	104-51-8			U	1.00		0.250	
n-Propylbenzene	103-65-1			U .	1.00		0.125	
m-,p-Xylene	136777-61-	-2		U	2.00		0.500	
Naphthalene	91-20-3			U	1.00		0.200	
o-Xylene	95-47-6			U	1.00		0.250 0.250	
p-Isopropyltoluene	99-87-6		.	บ บ	1.00 1.00		0.250	
sec-Butylbenzene	135-98-8			บ บ	1.00		0.125	
Styrene	100-42-5 79-01-6		2.7	U.	1.00		0.125	
Trichloroethene tert-Butylbenzene	98-05-6	4 4	· · ·	U	1.00		0.250	

Microbac

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Report Number: L08110347 Report Date :December 9, 2008

Sample Number:L08110347-01	PrePrep Metho	d · NONE		Inst	rumer	t:HPMS8	, 1		
Client ID:59SW4WG1	-	Prep Method: 5030B			Prep Date:11/22/2008 18:37				
Matrix:Water	Analytical Method:8260B			Cal Date:11/14/2008 18:					
Workgroup Number: WG288769	Analys			Run Date: 11/22/2008 18:37					
Collect Date:11/11/2008 14:16	Dilutic			File ID:8M349951					
Sample Tag:01	Units:ug/L			1					
Analyte	CAS, Number	 F	esult	Qual	•••	RL	MDI.		
Tetrachloroethene	127-18-4	127-18-4 0.305			()	1.00	0.250		
Toluene	108-88-3	108-88-3		Ū		1.00	0.250		
trans-1,2-Dichloroethene	156-60-5	156-60-5		~~ F ~~~	· }	1.00	0.250		
trans-1,3-Dichloropropene	10061-02-6			υ		1.00	0.500		
Trichlorofluoromethane	75-69-4	75-69-4 0.651		-R-	- 1	1.00	0.250		
Vinyl chloride	75-01-4			ប		1.00	0.250		
Surrogate	% Recovery	Lower	Uppe	r i	Qual				
Dibromofluoromethane	100	85	115			-			
1,2-Dichloroethane-d4	106	72	119						
Toluene-d8	99.9	81	120						

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.

119

 1,2-Dichloroethane-d4
 106

 Toluene-d8
 99.9

 4-Bromofluorobenzene
 97.3

 U Undetected: the avaiute
 97.3
 U Undetected; the analyte was analyzed for, but not detected.

F Because of quality control deficiences for this analyte, this data may be rejected.
F The analyte was positively identified, but the quantitation was below the RL.

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12/12/08

ample Number: L08110347-6	02	PrePrep Metho		Instrument: HPMS8	
Client ID:595W7WG1		Prep Metho	d:5030B	Prep Date:11/25/200	
Matrix:Water		Analytical Metho	d:8260B	Cal Date:11/14/200	8 18:53
group Number:WG288953		Analys	t:MES	Run Date:11/25/200	8 00:26
Collect Date:11/11/2008	16:01	Dilutio	n:1	File ID:8M350017	
Sample Tag:01	•		s:ug/L		
	•		When any service		
			a		
Analyte		CAS. Number	Result	Qual RL	MDL
1,1,1,2-Tetrachloroethar	16	630-20-6		ບ 0.500	0.250
1.1.1-Trichloroethane		71~55-6	1.88	1.00	0.250
1,1,2,2-Tetrachloroethar	18	79-34-5		U 0.500	0.125
1,1,2-Trichloroethane		79-00-5		U 1.00	0.250
1,1-Dichloroethane		75-34-3	5.04	1.00	0.125
1,1-Dichloroethene		75-35-4		U 1.00	0.500
1,1-Dichloropropene		563-58-6		υ 1.00	0.250
1,2,3-Trichlorobenzene		87-61-6		U 1.00	0.150
1,2,3-Trichloropropane		96-18-4		U 1.00	0.500
1.2.4-Trichlorobenzene		120-82-1		U 1.00	0.200
1.2.4-Trimethylbenzene		95-63-6		U 1.00	0.250
1.2-Dichloroethane		107-06-2		U 0.500	0.250
1.2-Dichlorobenzene		95-50-1		U 1.00	0.125
1,2-Dibromo-3-chloroprop	ane	96-12-8		ບ 2.00	1.00
1,2-Dichloropropane		78-87-5	and a second second	U 1.00	0.200
1,2-Dibromoethane	1	106-93-4		U 1.00	0,250
1,3,5-Trimethylbenzene		108-67-8		U 1.00	0.250
1,3-Dichlorobenzene		541-73-1		v 1.00	0.250
1,3-Dichloropropane	A	142-28-9		u 0.400	0.200
1,4-Dichlorobenzene		106-46-7		u 0.500	0.125
	-	123-91-1	• · · · ·	M 100	50.0
1,4-Dioxane			1		0.125
1-Chlorohexane		544-10-5		U 1.00 U 1.00	0.125
2,2-Dichloropropane		594-20-7		U 1.00	0.125
2-Chlorotoluene		95-49-8			
4-Chlorotoluene		106-43-4		U 1.00	0.250 2.50
Acetone		67-64-1			
Benzene		71-43-2			0.125
Bromobenzene		108-86-1		U 1.00	0.125
Bromochloromethane		74-97-5		U 1.00	0.200
Bromodichloromethane		75-27-4		U 0.500	0.250
Bromoform		75-25-2	· - · · · · · · ·	U 1.00	0.500
Bromomethane		74-83-9		U 3.00	0.500
Carbon tetrachloride		56-23-5		U 1.00	0.250
Chlorobenzene		108-90-7		U 0.500	0.125
Chloroethane		75-00-3		υ 1.00	0.500
Chloroform		67-66-3		U 0.300	0.125
Chloromethane		74-87-3		M 1.00	0.250
cis-1,2-Dichloroethene		156-59-2	35.3	M 1.00	0.250
cis-1,3-Dichloropropene		10061-01-5	and a second second	U 0.500	0.250
Dibromochloromethane		124-48-1		_R_ () 0.500	0.250
Dibromomethane		74~95-3		υ 1.00	0.250
Dichlorodifluoromethane		75-71-8		U 1.00	0.250
Ethylbenzene		100-41-4		U 1.00	0.250
Hexachlorobutadiene		87-68-3		U 0.600	0.250
Isopropylbenzene		98-82-8		U 1.00	0.250
Methylene chloride		75-09-2		U 1.00	0.250
Methyl t-butyl ether (MT	BE)	1634-04-4		υ 5.00	0.500
MEK (2-Butanone)	• • •	78-93-3		U 10.0	2.50
MIBK (methyl isobutyl ke	tone)	108-10-1		U 10.0	2.50
n-Butylbenzene		104-51-8		υ 1.00	0.250
n-Propylbenzene		103-65-1		υ 1.00	0.125
m-,p-Xylene		136777-61-2	· ··· •	ບ 2.00	0.500
Naphthalene	• -	91-20-3		U 1.00	0.200
o-Xylene		95-47-6		U 1.00	0.250
p-Isopropyltoluene		99-87-6		U 1.00	0.250
sec-Butylbenzene	÷	135-98-8		U 1.00	0.250
Styrene		100-42-5	e de la companya de l	U 1.00	0.125
		N	8.15	1.00	0.250
Trichloroethene		79-01-6			

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WINDL

Sample Number: L08110347-02	PrePrep Method:NONE			Instr	ument:HPMS8	,
Client ID:59SW7WG1	Prep Meth	od:5030E	3	Prep	Date:11/25/	2008 00:26
Matrix:Water	Analytical Meth	od:8260E	3	Cal	Date:11/14/	2008 18:53
Vorkgroup Number:WG288953	Analy	st:MES		Run	Date:11/25/	2008 00:26
Collect Date:11/11/2008 16:01 Sample Tag:01	Dilution:1 Units:ug/L				1D:8M350017	
Analyte	CAS. Number		Result	Qual	RL	MDL
Tetrachloroethene	127-18-4		0.590	- P	1.00	0.250
Toluene	108-88-3			v 1	1.00	0.250
trans-1,2-Dichloroethene	156-60-5		0.302		1.00	0.250
trans-1,3-Dichloropropene	10061-02-6	· · · · · · · · ·		U	1.00	0.500
Trichlorofluoromethane	75-69-4			ົບ	1.00	0.250
Vinyl chloride	75-01-4		1.21	M	1.00	0.250
Surrogate	% Recovery	Lower	Upp		Qual	
					· · ·	
Dibromofluoromethane	96.6	85	· 11	5		

Dibromofluoromethane96.61,2-Dichloroethane-d498.4Toluene-d81074-Bromofluorobenzene101
 72
 119

 81
 120

 76
 119

M Matrix effect; the concentration is an estimate due to matrix effect.
U Undetected; the analyte was analyzed for, but not detected.
R Because of quality control deficiences for this analyte, this data may be rejected.
F The analyte was positively identified, but the quantitation was below the RL.

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12/12 D

Sample Number:L08110347-03 Client ID:59SWTWG1-MS Matrix:Water Workgroup Number:WG288953 Collect Date:11/11/2008 16:01 Sample Tag:01	PrePrep Metho Prep Metho Analytical Metho Analys Dilutio Unit:	d:5030B d:8260B t:MES	Instrument:HPMS Prep Date:11/2 Cal Date:11/1 Run Date:11/2 File ID:8M35002	5/2008 00:58 4/2008 18:53 5/2008 00:58
Analyte	CAS. Number	Result	Qual RL	MDL
1,1,1,2-Tetrachloroethane	630-20-6	22.1	0.500	0.250
1,1,1-Trichloroethane 1,1,2,2-Tetrachloroethane	71-55-6 79-34-5	20.7	0.500	0.125
1,1,2-Trichloroethane	79-00-5	21.8	1.00	0.250
1.1-Dichloroethane	75-34-3	23.6	1.00	0.125
1,1-Dichloroethene	75-35-4	19.7	1.00	0.500
1,1-Dichloropropene 1,2,3-Trichlorobenzene	563-58-6 87-61-6	19.5 15.9	1.00	0.150
1,2,3-Trichloropropane	96-18-4	19.0	1.00	0.500
1,2,4-Trichlorobenzene	120-82-1	16.0	1.00	0.200
1,2,4-Trimethylbenzene	95-63-6	19.7	1.00	0.250
1.2-Dichloroethane	107-06-2	21.3 19.3	0.500	0.250
1,2-Dichlorobenzene 1,2-Dibromo-3-chloropropane	95-50-1 96-12-8	21.4	2.00	1.00
1,2-Dichloropropane	78-87-5	22.5	1.00	0.200
1.2-Dibromoethane	106-93-4	20.8	1.00	0.250
1,3,5-Trimethylbenzene	108-67-8	18.7	1.00	0.250
1,3-Dichlorobenzene	541-73-1 142-28-9	18.9 22.1	1.00	0.250 0.200
1,3-Dichloropropane 1,4-Dichlorobenzene	106-46-7	19.2	0.500	0.125
1,4-Dioxane	123-91-1	149	100	50.0
1-Chlorohexane	544-10-5	19.8	1.00	0.125
2,2-Dichloropropane	594-20-7	18.4	1.00	0.250
2-Chlorotoluene 4-Chlorotoluene	95-49-8 106-43-4	20.5	1.00	0.250
Acetone	67-64-1	46.2	R 10.0	2.50
Benzene	71-43-2	20.7	0.400	0.125
Bromobenzene	108-86-1	18.7	1.00	0.125
Bromochloromethane	74-97-5	19.7	1.00	0.200 0.250
Bromodichloromethane Bromoform	75-27-4 75-25-2	21.3 15.3	0.500	0.500
Bromomethane	74-83-9	25.2	3.00	0.500
Carbon tetrachloride	56-23-5	19.8	1.00	0.250
Chlorobenzene	108-90-7	19.8	0.500	0.125
Chloroethane	75-00-3	24.3	1.00	0.500 0.125
Chloroform Chloromethane	67-66-3 74-87-3	46.7	1.00	0.250
cis-1,2-Dichloroethene	156-59-2	48.8	1.00	0.250
cis-1,3-Dichloropropene	10061-01-5	19.8	0.500	0.250
Dibromochloromethane	124-48-1	19.5	R 0.500	0.250
Dibromomethane Dichlorodifluoromethane	74-95-3	21.9 25.5	1.00	0.250
Ethylbenzene	100-41-4	20.3	1.00	0.250
Hexachlorobutadiene	87-68-3	16.6	0.600	0.250
Isopropylbenzene	98-82-8	18.4	1.00	0.250
Methylene chloride	75-09-2	18.9 19.1	1.00	0.250
Methyl t-butyl ether (MTBE) MEK (2-Butanone)	1634-04-4 78-93-3	23.7	10.0	2.50
MIBK (methyl isobutyl ketone)	108-10-1	20.2	10.0	2.50
n-Butylbenzene	104-51-8	20.3	1.00	0.250
n-Propylbenzene	103-65-1	20.4	1.00	0.125
m-,p-Xylene	136777-61-2 91-20-3	41.2	2.00	0.500
Naphthalene o-Xylene	91-20-3	21.1	1.00	0.250
p-Isopropyltoluene	99-87-6	18.5	1.00	0.250
sec-Butylbenzene	135-98-8	20.0	1.00	0.250
Styrene	100-42-5	20.6	1.00	0.125
Trichloroethene	79-01-6 98-06-6	24.4 18.8	1.00	0.250
tert-Butylbenzene		20.0		

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Sample Number: £08110347-03	PrePrep Method:NONE			Instr	ument:HPMS8	ļ.	
Client ID:59SW7WG1-MS	Prep Me	thod: 50301	3	Prep	Date:11/25	5/2008 00	0:58
Matrix:Water	Analytical Me	thod: 82601	3	Cal	Date:11/14	/2008 18	8:53
Workgroup Number:WG288953	Ana	lyst:MES		Rur	Date:11/25	5/2008 0	0:58
Collect Date:11/11/2008 16:01	Dilu	tion:1		File	ID:8M350018	3	
Sample Tag:01	U	nits:ug/L			,		
Analyte	CAS. Numbe	er	Result	Qual	RL	Ň	юL
Tetrachloroethene	127-18-4		18.5		1,00	0.	250
Toluene	108-88-3		21.0		1.00	ο.	250
trans-1,2-Dichloroethene	156-60-5		19.8		/ 1.00	0.	250
trans-1,3-Dichloropropene	10061-02-	6	18.9		1.00	Ó,	500
Trichlorofluoromethane	75-69-4		20.3		1.00	0.	250
Vinyl chloride	75-01-4		44.7		1.00	0.	250
Surrogate	% Recovery	Lower	Upp	er	Qual		
Dibromofluoromethane	92.7	85	11	5.			
1,2-Dichloroethane-d4	97.3	72	11	9			
Toluene-d8	105	81	12	3			
4-Bromofluorobenzene	93.7	76	11	9			

4-Bromofluorobenzene93.776119R Because of quality control deficiences for this analyte, this data may be rejected.

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Sample Number: 108110347-04	PrePrep Method:NONE	Instrument: HPMS8
Client ID: 59SW7WG1-MSD	Prep Method:5030B	Prep Date:11/25/2008 01:30
Matrix:Water	Analytical Method:8260B	Cal Date:11/14/2008 18:53
Workgroup Number: WG288953	Analyst : MES	Run Date:11/25/2008 01:30
Collect Date:11/11/2008 16:01	Dilution:1	File ID:8M350019
Sample Tag:01	Units ug/L	
Analyte	CAS. Number Result	Qual RL 7 MDL
1.1.1.2 Tetrachloroethane	630-20-6 22.9	0.500 0.250

1, 1, 2; - Trichloroethame 7: -55 - 6 20.7 0.66 0.250 1, 1, 2; - Trichloroethame 73-00-5 23.3 1.00 0.253 1, 1, 2: Trichloroethame 73-35.4 25.9 1.00 0.233 1, 1, 2: Trichloroethame 73-35.4 0.7 1.00 0.235 1, 2: Trichloroethame 73-35.4 0.7 1.00 0.235 1, 2: Trichloroethame 76-61.6 17.3 1.00 0.200 1, 2: A Trichloroethame 107-06-2 21.0 0.500 0.250 1, 2: Dichloroethame 107-06-2 21.0 0.500 0.250 1, 2: Dichloroethame 106-35.4 21.4 1.00 0.250 1, 2: Dichloroethame 106-37.4 21.4 1.00 0.250 1, 3: Dichloroethame 106-37.4 21.4 1.00 0.250 1, 3: Dichloropropane 106-37.4 21.4 1.00 0.250 1, 3: Dichloropropane 106-46-7 19.4 1.00 0.250 1, 3: Dichloropropane 106-46-7 19.4 1.00 0.250 1, 3: Dichloropropane	1,1,1,2-Tetrachloroethane	630-20-6	22.9	0.500	0.250
1,1,2,2-retrachlorosthame 79-34-5 22.8 0.500 0.125 1,1,2-rtichlorosthame 79-34-3 25.9 1.00 0.125 1,1-Dichlorosthame 75:34-3 25.9 1.00 0.125 1,1-Dichlorosthame 75:34-4 20.7 1.00 0.500 1,2-Strichlorosthame 86-86-6 19.8 1.00 0.250 1,2-Strichlorosthame 86-18-4 21.7 1.00 0.500 1,2-Strichlorosthame 107.96-2 21.8 1.00 0.500 1,2-Strichlorosthame 107.96-2 21.8 1.00 0.250 1,2-Strichlorosthame 107.96-2 21.8 1.00 0.250 1,2-Dichtorosthame 106-93-4 21.4 1.00 0.250 1,2-Dichtorostopane 95-10-1 19.4 1.00 0.250 1,2-Dichtorostopane 95-10-1 19.4 1.00 0.250 1,2-Dichtorostopane 95-10-1 19.4 1.00 0.250 1,2-Dichtorostopane 106-97-8 19.5 1.00 0.200 1,3-Stritanthylbemsene 106-97-7 19.4<		71-55-6	20.7	1.00	0.250
1.1.2-Trichlorosthane 79-00-5 22.3 1.00 0.250 1.3-Diskhorosthane 75:35-4 20.7 1.00 0.123 1.1-Diskhorosthane 75:35-4 20.7 1.00 0.250 1.2-Diskhorosthane 87-61-6 17.3 1.00 0.250 1.2.3-Trichlorosthane 87-61-6 17.3 1.00 0.250 1.2.3-Trichlorosthane 107-66-7 21.8 0.500 0.250 1.2.4-Trichlorosthane 107-66-7 21.8 0.500 0.250 1.2-Diskhorosthane 107-66-7 21.8 0.500 0.250 1.2-Diskhorosthane 106-63-4 21.4 1.00 0.250 1.2-Disknosthane 108-67-5 22.8 1.00 0.250 1.3-Diskhorosthane 108-67-7 19.5 1.00 0.250 1.3-Diskhorosthane 108-67-7 19.5 1.00 0.250 1.3-Diskhorosthane 108-67-7 19.5 1.00 0.250 1.3-Diskhorosthane 142-28-9 2.7 0.400 <th></th> <th></th> <th>22.8</th> <th>0:500</th> <th>0.125</th>			22.8	0:500	0.125
1.1-Dichlorosthame 75:35:4 25.9 1.00 0.125 1.1-Dichlorosthame 75:35:4 20.7 1.00 0.500 1.2.3-Trichlorosthame 56:36:6 19.8 1.00 0.250 1.2.3-Trichlorosthame 56:36:6 19.8 1.00 0.250 1.2.3-Trichlorosthame 120:42:1 17.3 1.00 0.250 1.2-bichlorosthame 39:63:6 20.1 1.00 0.250 1.2-bichlorosthame 39:63:6 20.1 1.00 0.250 1.2-bichlorosthame 39:63:6 21.3 1.00 0.250 1.2-bichlorosthame 106:67:6 19.5 1.00 0.250 1.3-bichlorosthame 106:67:6 19.5 1.00 0.250 1.3-bichlorostograme 54:47:3-1 19.2 1.00 0.250 1.3-bichlorostograme 54:47:3-1 19.5 1.00 0.250 1.3-bichlorostograme 54:41:0-5 19.5 1.00 0.250 1.3-bichlorostograme 54:41:0-5 19.5 1.00 0.255 1.3-bichlorostograme 54:41:0-5 19.5 <th></th> <th></th> <th></th> <th></th> <th>0.250</th>					0.250
1.1-Dischloroptopene 75:35.4 20.7 1.00 0.500 1.2-Dischloroptopene 56:36.6 19.8 1.00 0.250 1.2.3-Trichloroptopene 57:61:6 17.3 1.00 0.250 1.2.3-Trichloroptopene 120:62:1 1.7 1.00 0.250 1.2.4-Trichloroptopene 120:62:1 17.3 1.00 0.250 1.2-Dischloroptopene 107:66:2 21.9 0.500 0.250 1.2-Dischloroptopene 35:50:1 19.4 1.00 0.250 1.2-Dischloroptopene 36:12:8 21.3 2.00 0.200 1.2-Dischloroptopene 108:57:4 21.4 1.00 0.200 1.3-Dischloroptopene 108:67:4 12.5 1.00 0.250 1.3-Dichloroptopene 106:43:4 20:7					0.125
1.1.1.01chloropropene 563-56.4 19.8 1.00 0.250 1.2.3 Trichloropopene 97-61-6 17.3 1.00 0.150 1.2.3 Trichloropopene 97-61-6 17.3 1.00 0.250 1.2.4 TrinchyDersene 120-62-1 17.3 1.00 0.250 1.2.2 Trichloropensene 107-66-2 21.8 0.500 0.255 1.2-Dichloropensene 95-50-1 19.4 1.00 0.255 1.2-Dichloropensene 78-87-5 22.8 1.00 0.255 1.3-Dichloropense 106-93-4 21.4 1.00 0.255 1.3-Dichloropensene 106-83-4 21.4 1.00 0.255 1.3-Dichloropensene 106-83-7 19.5 1.00 0.255 1.3-Dichloropensene 108-67-7 19.4 0.500 0.125 1.4-Dichloropensene 108-64-7 19.4 0.500 0.125 1.4-Dichloropensene 108-64-7 19.4 0.500 0.125 1.4-Dichloropensene 122-41-1 195		and the second of the second s			0.500
1.2.3 - Trichlorobarsene 97-61:6 17.3 1.00 0.150 1.2.3 - Trichlorobarsene 120-62-1 17.3 1.00 0.200 1.2.4 - Trichlorobarsene 120-62-1 17.3 1.00 0.250 1.2.4 - Trichlorobarsene 107-06-2 21.8 0.500 0.255 1.2-bichlorobarsene 35-50-1 19.4 1.00 0.255 1.2-bichlorobarsene 36-12-8 21.3 2.00 1.00 1.2-bichloropropane 78-87-5 2.2 2.0 1.00 0.250 1.3-bichloropropane 108-67-8 19.5 1.00 0.250 1.3-bichloropropane 124-22-1 0.00 0.250 1.3-5 1.3-bichloropropane 124-23-1 19.5 1.00 0.250 1.3-bichloropropane 124-24-2 19.5 1.00 0.250 1.4-bichlorobarsene 123-91-1 19.5 1.00 0.250 2.2-chorotoluene 59-49-6 19.3 1.00 0.250 2.2-chorotoluene 57-64-1 33.2 R 10.0 0.250 2-chlorotoluene		(1) (1) (1) (1) (1) (1) (1) (1) (1) (1)		1	
1.2.3 Trichloropropene 56.18-4 21.7 1.00 0.500 1.2.4 Trinethylbenzene 120.62.1 17.3 1.00 0.250 1.2.bithlorosethane 107.66-2 21.8 0.500 0.255 1.2.bithlorosethane 95-50.1 19.4 1.00 0.255 1.2.bithlorosethane 95-12.4 21.3 2.00 1.00 1.2.bithlorosethane 106-63.4 21.4 1.00 0.2250 1.3.bithlorosethane 106-63.4 21.4 1.00 0.2250 1.3.bithlorosethane 106-63.4 21.4 1.00 0.2250 1.3.bithlorosethane 124.28.9 1.9.2 1.000 0.2250 1.3.bithlorosethane 124.428.9 22.7 0.400 0.2250 1.4.bithorosethane 124.428.9 1.9.5 1.00 0.125 1.4.bithorosethane 124.428.9 1.9.5 1.00 0.255 1.4.bithorosethane 124.420.7 19.5 1.00 0.255 2.Chiorotoluene 104.43.4 20.3 1.00 0.255 2.Chiorotoluene 74.97.5 20.9<		a second s			
1.2.4.Trichlorobersone 12.0.4.Trichlorobersone 17.3 1.0.0 0.200 1.2.4.Trichlorobersone 95-63.6 20.1 1.0.0 0.250 1.2.Jubicorobersone 95-63.6 21.8 0.500 0.255 1.2.Jubicorobersone 95-63.6 21.8 1.00 0.255 1.2.Jubicorobersone 95-63.6 21.8 2.00 1.00 1.2.Jubicorobersone 96-87.6 22.8 1.00 0.250 1.3.S.Trinchylbersone 106-67.8 19.5 1.00 0.250 1.3.S.Trinchylbersone 106-67.8 19.5 1.00 0.250 1.3.S.Trinchylbersone 106-646.7 19.4 0.500 0.125 1.4.Dichlorobersone 106-46.7 19.5 1.00 0.250 1.4.Dichlorobersone 594-20.7 19.0 1.00 0.250 2.7.Dichloropropane 594-20.7 19.0 1.00 0.250 2.7.Dichloropropane 594-20.7 19.0 1.00 0.250 2.7.Dichloropropane 106-43.4 20.3 1.00 0.250 2.1.Dichloropropane 106-43.4					
1.2.4.Trimethylbersene 75-63-6 20.1 1.00 0.250 1.2.9bichlorsethane 107-66-2 21.8 0.500 0.255 1.2.9bichlorsethane 95-50-1 19.4 1.00 0.255 1.2.9bichlorsethane 96-12-8 21.3 2.00 1.00 0.255 1.2.9bichlorsethane 106-53-4 21.4 1.00 0.255 1.3.9bichlorsethane 106-67-8 19.5 1.00 0.255 1.3.9bichlorsethane 106-67-1 19.4 0.00 0.255 1.3.9bichlorsethane 123-91-1 195 1.00 0.255 1.4-bicknesene 123-91-1 195 1.00 0.125 2bichlorsepane 594-20-7 19.5 1.00 0.255 2bichlorsepane 594-9.6 19.3 1.00 0.255 2chiorstoluene 67-64+1 33.2 R 10.00 0.255 2chiorstoluene 74-97-5 20.9 1.00 0.255 2chiorstoluene 75-27-4 20.9 </th <th></th> <th></th> <th></th> <th></th> <th></th>					
1 2-Dischlorochane 10 ² -6-2 21.8 0.500 0.250 1,2-Dikronobenzene 96-50-1 21.8 1.00 0.125 1,2-Dikronos3-chloropropane 96-12-8 22.3 2.00 1.00 1,2-Dikronos4-chloropropane 106-93-4 21.4 1.00 0.250 1,2-Dichloropropane 106-63-4 21.4 1.00 0.250 1,3-Dichlorobenzene 106-64-7 19.5 1.00 0.250 1,3-Dichlorobenzene 106-46-7 19.4 0.500 0.125 1,4-Dicklorobenzene 106-46-7 19.4 0.500 0.125 1,4-Dickloropenzene 544-10-5 19.5 100 50.0 1-Chlorohexane 544-10-5 19.5 100 0.250 2-Chlorotoluene 554-9-8 19.3 1.00 0.250 2-Chlorotoluene 105-43-4 20.3 1.00 0.250 2-Chlorotoluene 71-43-2 20.4 0.400 0.125 Bromobenzene 71-52-7 20.9 1.00					
1.2-Dichlorobersene 95-50-1 1.4.1 1.40 0.125 1.2-Dichloropropane 78-87-5 21.8 1.00 0.250 1.2-Dichloropropane 106-63-4 21.4 1.00 0.250 1.3-J.5thinorobensene 106-67-8 19.5 1.00 0.250 1.3-J.5thinorobensene 102-67-8 19.5 1.00 0.250 1.3-J.5thinorobensene 102-67-8 19.4 0.500 0.125 1.4-Jichlorobensene 102-67-7 19.4 0.500 0.125 1.4-Jichlorobensene 102-67-8 19.4 0.500 0.125 1.4-Jichlorobensene 102-64-7 19.0 0.00 0.250 2Alcrotoluene 594-20-7 19.0 1.00 0.250 2Chlorotoluene 594-20-7 19.0 1.00 0.250 2Chlorotoluene 75-49 20.3 1.00 0.250 2Chlorotoluene 75-27-2 20.4 0.400 0.125 Bromochloromethane 75-27-2 20.5 0.500					
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Cis-1, 2-Dichloroethene156-59-251.31.000.250cis-1, 3-Dichloropropene10061-01-520.00.5000.250Dibromochloromethane124-48-120.6R0.5000.250Dibromomethane74-95-322.41.000.250Dichlorodifluoromethane75-71-829.31.000.250Ethylbenzene100-41-420.01.000.250Hexachlorobutadiene87-68-316.80.6000.250Methylene chloride75-79-220.51.000.250Methylene chloride75-09-220.51.000.250Methylene chloride78-93-326.210.02.50MiBK (methyl isobutyl ketone)108-10-121.510.02.50n-Propylbenzene103-65-120.71.000.250m-Propylbenzene103-65-120.21.000.250m-Propylbenzene103-65-120.21.000.250n-Propylbenzene136777-61-239.32.000.500n-Propylbenzene95-47-620.31.000.250n-Propyltoluene99-87-618.71.000.250sec-sutylbenzené135-98-819.51.000.250sec-sutylbenzené135-98-819.51.000.250sec-sutylbenzené135-98-819.51.000.250sec-sutylbenzené135-98-819.51.000.250sec-sutylbenzené135-98-819.51.000.250 </th <th></th> <th></th> <th></th> <th></th> <th></th>					
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Dibromochloromethane 124-48-1 20.6 R 0.500 0.250 Dibromochloromethane 74-95-3 22.4 1.00 0.250 Dibromothlane 75-71-8 29.3 1.00 0.250 Ethylbenzene 100-41-4 20.0 1.00 0.250 Hexachlorobutadiene 87-68-3 16.8 0.600 0.250 Isopropylbenzene 98-82-8 18.1 1.00 0.250 Methylene chloride 75-09-2 20.5 1.00 0.250 Methylene chloride 78-93-3 26.2 10.0 2.50 Methyl t-butyl ether (MTBE) 1634-04-4 21.6 5.00 0.500 MEK (2-Butanone) 108-10-1 21.5 10.0 2.50 n-Butylbenzene 104-51-8 20.7 1.00 0.250 n-Propylbenzene 103-65-1 20.2 1.00 0.250 m-propylbenzene 13677-61-2 39.3 2.00 0.500 n-Propylbenzene 91-20-3 18.7 1.00	cis-1,2-Dichloroethene	the second se			
Dibromomethane 74-95-3 22.4 1.00 0.250 Dichlorodifluoromethane 75-71-8 29.3 1.00 0.250 Ethylbenzene 100-41-4 20.0 1.00 0.250 Hexachlorobutadiene 87-68-3 16.8 0.600 0.250 Isopropylbenzene 98-82-8 18.1 1.00 0.250 Methylene chloride 75-09-2 20.5 1.00 0.250 Methyl t-butyl ether (MTBE) 1634-04-4 21.6 5.00 0.500 MEK (2-Butanone) 78-93-3 26.2 10.0 2.50 MIBK (methyl isobutyl ketone) 108-10-1 21.5 10.0 2.50 n-Butylbenzene 103-65-1 20.2 1.00 0.250 n-Propylbenzene 103-65-1 20.2 1.00 0.250 m-propylbenzene 91-20-3 18.7 1.00 0.250 m-propylbenzene 91-20-3 18.7 1.00 0.250 m-propylbenzene 91-20-3 18.7 1.00 0.250<				1	
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Ethylbenzene100-41-420.01.000.250Hexachlorobutadiene87-68-316.80.6000.250Isoproylbenzene98-82-818.11.000.250Methylene chloride75-09-220.51.000.250Methylene chloride75-09-220.51.000.250Methylene chloride78-93-326.210.02.50MEK (2-Butanone)108-10-121.510.02.50NIBK (methyl isobutyl ketone)108-10-121.510.02.50n-Butylbenzene103-65-120.21.000.125m-, p-Xylene136777-61-239.32.000.500Naphthalene91-20-318.71.000.250p-Isopropyltoluene99-87-618.71.000.250sec-Butylbenzené135-98-819.51.000.250rrichloroethene79-01-625.81.000.250					
Hexachlorobutadiene87-68-316.80.6000.250Isopropylbenzene98-82-818.11.000.250Methylene chloride75-09-220.51.000.250Methyl t-butyl ether (MTBE)1634-04-421.65.000.500MEK (2-Butanone)78-93-326.210.02.50MIBK (methyl isobutyl ketone)108-10-121.510.02.50n-Butylbenzene104-51-820.71.000.250n-Propylbenzene103-65-120.21.000.125m-, p-Xylene136777-61-239.32.000.500Naphthalene91-20-318.71.000.250o-Xylene95-47-620.31.000.250sec-Butylbenzene135-98-819.51.000.250rystene135-98-819.51.000.250rystene135-98-819.51.000.250rtichloroethene79-01-625.81.000.250	Dichlorodifluoromethane				
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n-Butylbenzene104-51-820.71.000.250n-Propylbenzene103-65-120.21.000.125m.pXylene136777-61-239.32.000.500Naphthalene91-20-318.71.000.220o-Xylene95-47-620.31.000.250p-Isopropyltoluene99-87-618.71.000.250sc-Butylbenzené135-98-819.51.000.250Styrene100-42-520.11.000.250Trichloroethene79-01-625.81.000.250					
n-Propylbenzene 103-65-1 20.2 1.00 0.125 m-,p-Xylene 136777-61-2 39.3 2.00 0.500 Naphthalene 91-20-3 18.7 1.00 0.220 o-Xylene 95-47-6 20.3 1.00 0.250 p-Isopropyltoluene 99-87-6 18.7 1.00 0.250 sec-Butylbenzené 135-98-8 19.5 1.00 0.250 Styrene 100-42-5 20.1 1.00 0.125 Trichloroethene 79-01-6 25.8 1.00 0.250					
m.,pXylene 136777-61-2 39.3 2.00 0.500 Naphthalene 91-20-3 18.7 1.00 0.200 o-Xylene 95-47-6 20.3 1.00 0.250 p-Isopropyltoluene 99-87-6 18.7 1.00 0.250 sec-Butylbenzené 135-98-8 19.5 1.00 0.250 Styrene 100-42-5 20.1 1.00 0.125 Trichloroethene 79-01-6 25.8 1.00 0.250	n-Butylbenzene				
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o-Xylene95-47-620.31.000.250p-Isopropyltoluene99-87-618.71.000.250sec-Butylbenzene135-98-819.51.000.250Styrene100-42-520.11.000.125Trichloroethene79-01-625.81.000.250	m-,p-Xylene				
p-Isopropyltoluene 99-87-6 18.7 1.00 0.250 sec-Butylbenzené 135-98-8 19.5 1.00 0.250 Styrene 100-42-5 20.1 1.00 0.125 Trichloroethene 79-01-6 25.8 1.00 0.250					
Jospinistic 135-98-8 19.5 1.00 0.250 Styrene 100-42-5 20.1 1.00 0.125 Trichloroethene 79-01-6 25.8 1.00 0.250	o-Xylene				
Styrene 100-42-5 20.1 1.00 0.125 Trichloroethene 79-01-6 25.8 1.00 0.250	p-Isopropyltoluene	99-87-6			
Styrene 100-42-5 20.1 1.00 0.125 Trichloroethene 79-01-6 25.8 1.00 0.250	sec-Butylbenzene				
ilicatoroechene	Styrene	100-42-5		1	
tert-Butylbenzene 98-06-6 19.3 1.00 0.250	Trichloroethene	79-01-6	25.8		
	tert-Butylbenzene	98-06-6	19.3	1.00	0.250

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,

Sample Number: L08110347-04	PrePrep Meth	od:NONE		Instrumen	L:HPMS8	
Client ID:59SW7WG1-MSD	Prep Meth	od:5030B		Prep Dat	e:11/25	/2008 01:30
Matrix:Water	Analytical Meth	od:8260B		Cal Dat	e:11/14	/2008 18:53
rkgroup Number:WG288953	Analy	st:MES		Run Dat	e:11/25	/2008 01:30
Collect Date:11/11/2008 16:01	Diluti	on:1		File ID:8	M350019)
Sample Tag:01	Uni	ts:ug/L				
					7	
Analyte	CAS. Number	R	esult	Qual	RĹ	MDL
Tetrachloroethene	127-18-4		18.6		1.00	0.250
Toluene	108-88-3		20.9	111	1.00	0.250
trans-1,2-Dichloroethene	156-60-5		21.0	• • • • •	1.00	0.250
trans-1,3-Dichloropropene	10061-02-6		19.3		1.00	0.500
Trichlorofluoromethane	75-69-4	1	20.2		1.00	0.250
Vinyl chloride	75-01-4		40.9		1.00	0.250
Surrogate	% Recovery	Lower	Upper	Qual		
Dibromofluoromethane	95.5	85	115			
1,2-Dichloroethane-d4	97.6	72	119			
Toluene-d8	105	81	120			
4-Bromofluorobenzene	96.6	76	119			

R Because of quality control deficiences for this analyte, this data may be rejected.

Microbac

of 14

Sample Number:L08110347-05PrePrep Method:NONEInstrument:HPMS8Client ID:S9SW3WG1Prep Method:5030BPrep Date:11/22/2008 19:09Matrix:WaterAnalytical Method:8260BCal Date:11/14/2008 18:53Workgroup Number:WG288769Analyst:MESRun Date:11/12/2008 19:09Collect Date:11/11/2008 16:39Dilution:1File ID:88349952Sample Tag:01Units:ug/LVOR

1,1,1,2-Tetrachloroethane		630-20-6		υ ο	. 500	0.250
1,1,1-Trichloroethane		71-55-6	0.345		. 00	0.250
1,1,2,2-Tetrachloroethane		79-34-5			.500	0.125
1,1,2-Trichloroethane	1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 - 1999 -	79-00-5	. :	ีขั 1	.00	0.250
1,1-Dichloroethane		75-34-3	······································	់ប ់ រ	.00	0.125
1,1-Dichloroethene		75-35-4		ניט		0.500
1,1-Dichloropropene		563-58-6		נ ע	1.00	0.250
1,2,3-Trichlorobenzene		87-61-6		ີ 🗘 🗍 1		0.150
1,2,3-Trichloropropane		96-18-4	· · · · · · · · · · · ·	נ ט	L.00	0.500
1,2,4-Trichlorobenzene		120-82-1		ับ 1		0.200
1,2,4-Trimethylbenzene		95-63-6		U 1		0.250
1,2-Dichloroethane		107-06-2		ບ 0	.500	0.250
1,2-Dichlorobenzene		95-50-1			1.00	0.125
1,2-Dibromo-3-chloropropane		96-12-8			2.00	1.00
1,2-Dichloropropane		78-87-5			1.00	0.200
1,2-Dibromoethane		106-93-4			L.00	0.250
1,3,5-Trimethylbenzene		108-67-8	· · · · · · · · · · · · · · · · · · ·		L.00	0.250
1,3-Dichlorobenzene		541-73-1	.		L.00	0.250
1.3-Dichloropropane		142-28-9			.400	0.200
1,4-Dichlorobenzene		105-46-7			.500	0.125
1.4-Dioxane		123-91-1			100	50.0
l-Chlorohexane		544-10-5		1. 7	L.00	0.125
2.2-Dichloropropane		594-20-7		1 1 1	1.00	0.250
2-Chlorotoluene		95-49-8		1	1.00	0.125
4-Chlorotoluene		106-43-4		1. 7	1.00 10.0	0.250 2.50
Acetone		67-64-1			.400	0.125
Benzene		71-43-2			1.00	0.125
Bromobenzene		74-97-5			1.00	0.200
Bromochloromethane		75-27-4			.500	0.250
Bromodichloromethane Bromoform		75-25-2			1.00	0.500
Bromomethane	- · · · ·	74-83-9			3.00	0.500
Carbon tetrachloride		56-23-5	· · · · · · · ·		1.00	0.250
Chlorobenzene	1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1	108-90-7	· · · · · · · · · · · · · · · · · · ·		500	0.125
Chloroethane	··· ···	75-00-3			1.00	0.500
Chloroform	· ·	67-66-3			.300	0.125
Chloromethane		74-87-3		υ	1.00	0.250
cis-1,2-Dichloroethene	· ·	156-59-2		υ	1.00	0.250
cis-1,3-Dichloropropene		10061-01-5		บ 0	.500	0.250
Dibromochloromethane		124-48-1		-0-1110	.500	0.250
Dibromomethane		74-95-3		υ	1.00	0.250
Dichlorodifluoromethane		75-71-8		R	1.00	0.250
Ethylbenzene		100-41-4			1.00	0.250
Hexachlorobutadiene		87-68-3			.600	0.250
Isopropylbenzene		98-82-8		a sheep a contra	1.00	0.250
Methylene chloride		75-09-2			1.00	0.250
Methyl t-butyl ether (MTBE)		1634-04-4			5.00	0.500
MEK (2-Butanone)		78-93-3		1	10.0	2.50
MIBK (methyl isobutyl ketone)		108-10-1		d . T	10.0	2.50
n-Butylbenzene		104-51-8			1.00	0.250
n-Propylbenzene	1	103-65-1		A DECEMBER OF THE T	1.00	0.125
m-,p-Xylene		136777-61-2		· · · ·	2.00	0.500
Naphthalene		91-20-3		1 * .	1.00	0.250
o-Xylene		95-47-6			1.00	0.250
p-Isopropyltoluene		99-87-6 135-98-8	.		1.00	0.250
sec-Butylbenzene		100-42-5		1 T	1.00	0.125
Styrene		79-01-6	0.759		1.00	0.250
Trichloroethene tert-Butylbenzene	· ·	98-06-6	0.137		1.00	0.250
eere bacythemsene		20 00 0	· · · · · · · · · · · · · · · · · · ·	.1 7		

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12/12/08

Sample Number: L08110347-05	PrePrep Method:NONE			Instrument: HPMS8			
Client ID: 59SW3WG1	Prep Met	hod:5030B		Prep	Date:11/22	2008 19:09	
Matrix:Water	Analytical Met	hod:8260B		Cal	Date:11/14	2008 18:53	
Workgroup Number:WG288769	Anal	yst:MES		Run	Date:11/22	2008 19:09	
Collect Date:11/11/2008 16:39	Dilut	ion:1		File	ID:8M349952		
Sample Tag:01	Ur	its:ug/L					
			nii	0	77	MIDT	
Analyte	CAS. Numbe	r	Result	Qual	КГ	MDL	
Tetrachloroethene	127-18-4			U :	1.00	0.250	
Toluene	108-88-3			υ.	1.00	0.250	
trans-1,2-Dichloroethene	156-60-5			σ	1.00	0.250	
trans-1,3-Dichloropropene	10061-02-6	5		บ	1.00	0.500	
Trichlorofluoromethane	75-69-4			υ	1.00	0.250	
Vinyl chloride	75-01-4			U	1.00	0.250	
Surrogate	% Recovery	Lower	Upper	ç)ual		
Dibromofluoromethane	103	85	115				
1,2-Dichloroethane-d4	109	72	119				
Toluene-d8	99.7	81	120				
4-Bromofluorobenzene	95.6	76	119				
· · · · · · · · · · · · · · · · · · ·							

U Undetected; the analyte was analyzed for, but not detected. R Because of quality control deficiences for this analyte, this data may be rejected. F The analyte was positively identified, but the quantitation was below the RL.

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Microbac

12/12/38

Sample Number: L08110347-06	PrePrep Method:NONE			ment:HPMS8	· · · · · · · · · · · · · · · · · · ·
Client ID: 59DW3WG1	Prep Method:50	30B		Date:11/22/2	
Matrix:Water	Analytical Method:82	60B	Cal	Date:11/14/2	2008 18:53
Workgroup Number: WG288769	Analyst:ME	S	Run	Date:11/22/2	2008 19:41
Collect Date:11/11/2008 18:25	Dilution:1		File 3	D:8M349953	
Sample Tag:01	Units:ug	/L			
Analyte	CAS. Number	Result	Qual	RL	MDL
1,1,1,2-Tetrachloroethane	630-20-6		U	0.500	0.250
1,1,1-Trichloroethane	71-55-6		υ	1.00	0.250
1,1,2,2-Tetrachloroethane	79-34-5		U	0.500	0.125
1,1,2,2-Trichloroethane	79-00-5	· · · · · ·	u i	1.00	0.250
1,1-Dichloroethane	75-34-3		1 1	1.00	0.125
	75-35-4		- <u>-</u>	1.00	0.500
1,1-Dichloroethene	563-58-6		u u	1.00	0.250
1,1-Dichloropropene	87-61-6		-l - n	1.00	0.150
1,2,3-Trichlorobenzene	96-18-4		- ñ	1.00	0.500
1,2,3-Trichloropropane	120-82-1		1 II	1.00	0,200
1,2,4-Trichlorobenzene	and the second	• • • •		1.00	0.250
1,2,4-Trimethylbenzene	95-63-6			the second se	0.250
1,2-Dichloroethane	107-06-2	1. a	U	0.500	a service and the service of the ser
1,2-Dichlorobenzene	95-50-1	·	U	1.00	0.125
1 2 Dibromo 2 ablaropropana	96-12-8		i U	2.00	1.00

1,1-Dichloropropene	563-58-6	U 1.00	0.250
1,2,3-Trichlorobenzene	87-61-6	U 1.00	0.150
1,2,3-Trichloropropane	96-18-4	U 1.00	0.500
1,2,4-Trichlorobenzene	120-82-1	U 1.00	0.200
1,2,4-Trimethylbenzene	95-63-6	υ 1.00	0.250
1,2-Dichloroethane	107-06-2	ບ 0.500	0.250
1,2-Dichlorobenzene	95-50-1	U 1.00	0.125
1,2-Dibromo-3-chloropropane	96-12-8	υ 2.00	1.00
1,2-Dichloropropane	78-87-5	U 1.00	0.200
1,2-Dibromoethane	106-93-4	U 1.00	0.250
1,3,5-Trimethylbenzene	108-67-8	U 1.00	0.250
1,3-Dichlorobenzene	541-73-1	U 1.00	0.250
1,3-Dichloropropane	142-28-9	υ 0.400	0.200
1,4-Dichlorobenzene	106-46-7	ບ 0.500	0.125
1.4-Dioxane	123-91-1	U 100	50.0
1-Chlorohexane	544-10-5	ບ 1.00	0.125
2,2-Dichloropropane	594-20-7	υ 1.00	0.250
2-Chlorotoluene	95-49-8	U 1.00	0.125
4-Chlorotoluene	106-43-4	U 1.00	0.250
Acetone	67-64-1	R()) 10.0	2.50
Benzene	71-43-2	U V 0.400	0.125
Bromobenzene	108-86-1	U 1.00	0.125
Bromochloromethane	74-97-5	U 1.00	0.200
Bromodichloromethane	75-27-4	U 0.500	0.250
Bromoform	75-25-2	U 1.00	0.500
Bromomethane	74-83-9	U 3.00	0.500
Carbon tetrachloride	56-23-5	U 1.00	0.250
Chlorobenzene	108-90-7	ບ 0.500	0.125
Chloroethane	75-00-3	U 1.00	0.500
Chloroform	67-66-3	U 0.300	0.125
Chloromethane	74-87-3	U 1.90	0.250
	156~59-2 65.2	1.00	0.250
cis-1,2-Dichloroethene cis-1,3-Dichloropropene	10061-01-5	U 0.500	0.250
Dibromochloromethane	124-48-1	-U- () 0.500	0.250
Dibromochloromethane	74-95-3	0 1.00	0.250
Dichlorodifluoromethane	75-71-8	R 1.00	0.250
	100-41-4	u 1.00	0.250
Ethylbenzene	87-68-3	U 0.600	0.250
Hexachlorobutadiene	98-82-8	U 1.00	0.250
Isopropylbenzene	75-09-2	U 1.00	0.250
Methylene chloride	1634-04-4	υ 5.00	0.500
Methyl t-butyl ether (MTBE)	78-93-3	ບ 10.0	2.50
MEK (2-Butanone)	108-10-1	U 10.0	2.50
MIBK (methyl isobutyl ketone)	104-51-8	u 1.00	0.250
n-Butylbenzene	103-65-1	U 1.00	0.125
n-Propylbenzene	136777-61-2	U 2.00	0.500
m-,p-Xylene	91-20-3	U 1.00	0.200
Naphthalene	95-47-6	U 1.00	0.250
o-Xylene	99-87-6	U 1.00	0.250
p-Isopropyltoluene	135-98-8	U 1.00	0.250
sec-Butylbenzene	100-42-5	u 1.00	0.125
Styrene	79-01-6	U 1.00	0.250
Trichloroethene	98-06-6	U 1.00	0.250
tert-Butylbenzene	30-00-0	1	
	11 of 14		

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2

Sample Number:L08110347-06 Client ID:59DW3WG1 Matrix:Water Workgroup Number:WG288769 Collect Date:11/11/2008 18:25 Sample Tag:01	Diluti	od:5030E od:8260E st:MES		Prep Cal Run	ment:HPMS8 Date:11/22/ Date:11/14/ Date:11/22/ D:8M349953	
	·			Louis		2007
Analyte	CAS. Number		Result	Qual	κ <u>υ</u>	MDL
Tetrachloroethene	127-18-4			<u> </u>	1.00	0.250
Toluene	108-88-3			. <u>v</u>	1.00	0.250
trans-1,2-Dichloroethene	156-60-5		1.09		1.00	0.250
trans-1,3-Dichloropropene	10061-02-6			U	1.00	0.500
Trichlorofluoromethane	75-69-4	•••		U	1.00	0.250
Vinyl chloride	75-01-4			U	1.00	0.250
Surrogate	% Recovery	Lower	Uppe	er Qı	lal	
Dibromofluoromethane	101	85	115	5		
1,2-Dichloroethane-d4	108	72	119			
Toluene-d8	101	81	120) .		
4-Bromofluorobenzene	99.2	76	119	•		

U Undetected; the analyte was analyzed for, but not detected. R Because of quality control deficiences for this analyte, this data may be rejected.

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Microbac

12/12/05

Report Number: L08110347

Report Date : December 15, 2008

Sample Number:L08110347-07 Client ID:59DW3WG9 Matrix:Water Workgroup Number:WG288769 Collect Date:11/11/2008 18:25 Sample Tag:01 PrePrep Method:NONE Prep Method:5030B Analytical Method:8260B Analyst:MES Dilution:1 Units:ug/L Instrument:HPMS8 Prep Date:11/22/2008 20:13 Cal Date:11/14/2008 18:53 Run Date:11/22/2008 20:13 File ID:8M349954

a compatibility of the state of the	(1) (1) Marsham	Result	Qual	RL	MDL
Analyte	CAS. Number 630-20-6	Result	U	0.500	0.250
1,1,1,2-Tetrachloroethane	71-55-6		U	1.00	0.250
1,1,1-Trichloroethane	79-34-5		Ŭ	0.500	0.125
1,1,2,2-Tetrachloroethane	and the second		UU	1.00	0.250
1,1,2-Trichloroethane	79-00-5 75-34-3	0.410	- 17-	1.00	0.125
1,1-Dichloroethane	75-35-4	0.410	n n	1.00	0.500
1,1-Dichloroethene	563-58-6		ថ	1.00	0.250
1,1-Dichloropropene	87-61-6		បី	1.00	0.150
1,2,3-Trichlorobenzene	96-18-4		ີຫໍ່	1.00	0.500
1,2,3-Trichloropropane	120-82-1		Ū	1.00	0.200
1,2,4-Trichlorobenzene	95-63-6	· ·	ប	1.00	0.250
1,2,4-Trimethylbenzene	107-06-2		บั	0.500	0.250
1,2-Dichloroethane	95-50-1		u U	1.00	0.125
1,2-Dichlorobenzene	96-12-8		บ	2.00	1.00
1,2-Dibromo-3-chloropropane	78-87-5	•	ີຫ	1.00	0.200
1,2-Dichloropropane	106-93-4	(4) (4) (4) (4) (4) (4) (4) (4) (4) (4)	ΰ	1.00	0.250
1,2-Dibromoethane	108-53-4		ŭ	1.00	0.250
1,3,5-Trimethylbenzene	541-73-1		ŭ	1.00	0.250
1,3-Dichlorobenzene	142-28-9	· ·	ับ	0.400	0.200
1,3-Dichloropropane	106-46-7		์ บ	0.500	0.125
1,4-Dichlorobenzene	123-91-1	• •	ិប	100	50.0
1.4-Dioxane	544-10-5		ີ້ຫັ	1.00	0.125
1-Chlorohexane	594-20-7	• · · · · · ·	ੱਧ	1.00	0.250
2,2-Dichloropropane 2-Chlorotoluene	95-49-8	· · · · · ·	័ម	1.00	0.125
4-Chlorotoluene	106-43-4	• • • • •	υ	1.00	0.250
Acetone	67-64-1			10.0	2.50
Benzene	71-43-2	and the second second second	. . .	0.400	0.125
Bromobenzene	108-86-1	100 C	ំ ប៊ី	1.00	0.125
Bromochloromethane	74-97-5	· · · · · ·	ŭ	1.00	0.200
Bromodichloromethane	75-27-4		់ប់	0.500	0.250
Bromoform	75-25-2		Ū	1.00	0.500
Bromomethane	74-83-9		Ū	3.00	0.500
Carbon tetrachloride	56-23-5		Ū	1.00	0.250
Chlorobenzene	108-90-7		υ	0.500	0.125
Chloroethane	75-00-3	· · ·	ប	1.00	0.500
Chloroform	67-66-3		ប	0.300	0.125
Chloromethane	74-87-3		ΰ	1.00	0.250
cis-1,2-Dichloroethene	156-59-2	67.3		1.00	0.250
cis-1,3-Dichloropropene	10061-01-5		៍ប	0.500	0.250
Dibromochloromethane	124-48-1		U·	0.500	0.250
Dibromomethane	74-95-3		់ប	1.00	0.250
Dichlorodifluoromethane	75-71-8		R	1.00	0.250
Ethylbenzene	100-41-4		U	1.00	0.250
Hexachlorobutadiene	87-68-3		υ	0.600	0.250
Isopropylbenzene	98-82-8		υ	1.00	0.250
Methylene chloride	75-09-2		υ	1.00	0.250
Methyl t-butyl ether (MTBE)	1634-04-4		U	5.00	0.500
MEK (2-Butanone)	78-93-3		្មប	10.0	2.50
MIBK (methyl isobutyl ketone)	108-10-1		្ឋា	10.0	2.50
n-Butylbenzene	104-51-8		υ	1.00	0.250
n-Propylbenzene	103-65-1		. U	1.00	0.125
m-,p-Xylene	136777-61-2		U	2.00	0.500
Naphthalene	91-20-3	· · · · · · · · · · · · · · · · · · ·	υ	1.00	0.200
o-Xylene	95-47-6		U	1.00	0.250
p-Isopropyltoluene	99-87-6		U.	1.00	0.250
sec-Butylbenzene	135-98-8		U	1.00	0.250
Styrene	100-42-5		U	1.00	0.125
Trichloroethene	79-01-6		σ	1.00	0.250
tert-Butylbenzene	98-06-6		U	1.00	0.250

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12/12/08

MICLODAC HADOLALOLIES INC.

Report Number: L08110347

Report Date : December 15, 2008

Sample Number:L08110347-07 Client ID:59DW3WG9 Matrix:Water Workgroup Number:WG288769 Collect Date:11/11/2008 18:25	PrePrep Method:N Prep Method:5 Analytical Method:8 Analyst:M Dilution:1	030B 260B ES	Prep Cal Run	ment: HPMS8 Date: 11/22/2 Date: 11/14/2 Date: 11/22/2 D: 8M349954	2008 18:53
Sample Tag:01	Units: u				
Analyte	CAS. Number	Result	Qual	RL	MDL
Tetrachloroethene	127-18-4		ប	1.00	0.250
Toluene	108-88-3		U	1.00	0.250
trans-1,2-Dichloroethene	1.56-60-5	1.18		1.00	0.250
trans-1,3-Dichloropropene	10061-02-6		ΰ	1.00	0.500
Trichlorofluoromethane	75-69-4	0.285	F	1.00	0.250
Vinvl chloride	75-01-4		U	1.00	0.250

Vinyl chloride	75-01-4	U -		
Surrogate	% Recovery	Lower	Upper	Qual
Dibromofluoromethane	104	85	115	
1,2-Dichloroethane-d4	108	72	119	
Toluene-d8	102	81	120	
4-Bromofluorobenzene	103	76	119	

U Undetected; the analyte was analyzed for, but not detected.

R Because of quality control deficiences for this analyte, this data may be rejected.
 F The analyte was positively identified, but the quantitation was below the RL.

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Microbac

12/12/02

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1.0 INTRODUCTION

2.0 VOLATILE ORGANIC CONSTITUENTS

- 2.1 Holding Times
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- 2.6 Surrogate Recovery
- 2.7 Duplicates
- 2.8 Summary

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- 2 Field Sample ID/Lab Sample ID Cross Reference

APPENDICES

A Hand-Annotated Results Summary Forms

1.0 INTRODUCTION

This data quality review pertains to a soil sample collected in November 2008 at Air Force Plant 59 (AFP-59). Parameters evaluated in soil samples included the total concentration of volatile organic constituent (VOC). The samples were analyzed by Microbac Services, Marietta, Ohio.

Data quality review is an after-the-fact technical review of analytical data whereby the quality and usability of the data are determined based on a set of predefined criteria. These criteria depend upon the type of data involved and the purpose for which those data were collected. Data quality review assesses whether and to what extent specified criteria were met, and places restrictions on data use based on quality parameters. The data quality review process can range from a cursory review used to detect out-of-control situations to a detailed evaluation, depending on the analytical protocol, the associated quality control samples collected, and the intended data use.

Specific criteria for data quality review may include, but are not limited to: technical holding times, analysis of blanks, surrogate spike recovery, analysis of duplicates, and reported practical quantitation limits (PQLs). Where applicable, the recommendations of USEPA SW-846 *Test Methods for Evaluating Solid Waste* (Third Edition, December 1996) or USEPA *Methods for Chemical Analysis of Water and Wastes* (Revised March 1983) analytical method requirements, USEPA *CLP National Functional Guidelines for Organic and Inorganic Data Review* (February 1994, *Functional Guidelines*) data review guidance, and professional judgment.

Table 1 presents the data qualifiers applied during this review effort and their meanings.

Table 1 Data Qualifiers

Qualifier	Description
J	This is an estimated value.
UJ	The analyte was analyzed for but was not detected. The associated value is an estimate and may be inaccurate or imprecise.
U	The analyte was analyzed for, but not detected. The associated numerical value is at or below the MDL.

Table 2 provides a cross-reference list for field sample IDs and lab sample IDs.

Table 2 Field Sample ID/Lab Sample ID Cross Reference

Field Sample ID		Field Sample ID	Lab Sample ID
TRIP BLANK	L08110371-01	59SG36-3.0-SO1	L08110371-02

During the data quality review process, laboratory qualified and unqualified data are verified against all available supporting documentation. Based on this review, qualifier codes may be added, deleted, or modified by the validator. Final results are therefore either qualified or unqualified. (Note: In those cases where the laboratory added a "U" flag indicated a non-detect result, and the validator agrees with this flag, then it remains intact, as noted on the corresponding Results Summary Form.) Changes to the data are reflected on the Results Summary Forms in Appendix A.

2.0 VOLATILE ORGANIC CONSTITUENTS

Volatile organic constituents were analyzed using EPA Test Method for Evaluating Solid Waste (SW-846) Method 8260B. Samples were analyzed in batches:

WG288953: TRIP BLANK

WG288783: 59SG36-3.0-SO1

2.1 Holding Times

All samples were analyzed within prescribed hold times. No qualification is needed. Samples were collected via 5035 preserved by freezing. Sample prep proceeded as normally.

2.2 Calibration

WG288953: The Initial Calibration for this batch was performed on 11/24/2008. Standards were analyzed at 0.30, 0.4, 1, 2, 5, 20, 50, 100, and 200 ug/L in support of the Initial Calibration. The %RSD values were less than 30% for all target constituents. No qualification is needed.

The %D for the second source calibration was within $\pm 25\%$ for target constituents except for acetone. The validator qualifies UJ or J, the non detect and positive results, respectively, in the acetone results in the associated samples. The validator removes the "R" flag assigned by the laboratory to denote anomalies.

For the associated continuing calibration standard, dibromochloromethane associated %D values were greater than 20% for target constituents. The validator qualifies UJ or J, the non detect and positive results, respectively, in the acetone results in the associated samples. The validator removes the "R" flag assigned by the laboratory to denote anomalies.

<u>WG288783</u>: The Initial Calibration for this batch was performed on 11/22/2008. Standards were analyzed at 0.30, 0.4, 1, 2, 5, 20, 50, and 100 ug/kg in support of the Initial Calibration. The %RSD values were less than 30% for all target constituents. No qualification is needed.

The %D for the second source calibration was within $\pm 25\%$ for target constituents No qualification is needed.

For the associated continuing calibration standard, none of the associated %D values were greater than 20% for target constituents. No qualification is needed.

For all samples, it is noted that for those results which were less than the RL but greater than the MDL, the laboratory assigned an "F" flag, indicating an estimated value. Unless qualified otherwise, the validator removes the F flag and replaces it with the "J" qualifier, indicating an estimated value.

2.3 Laboratory Control Samples

<u>WG288953</u>: Laboratory control sample exhibited a 159% recovery for acetone, which are above the LCS limits. Since acetone was already qualified, no additional qualification is needed.

<u>WG288783</u>: Constituent recoveries from the associated laboratory control samples were within control limits. No qualification is needed.

2.4 Blanks

WG288953: No constituents were detected in the associated method blank or trip blank. No qualification needed.

WG288783: Acetone was detected at 6.37 ug/kg, naphthalene at 0.646 ug/kg, and 1,2,3trichlorobenzene at 0.563 ug/kg in associated method blank. The validator qualifies U any positive result less than or equal to 63.7 ug/kg in acetone; 3.23 ug/kg in naphthalene, and 2.815 ug/kg in 1,2,3trichlorobenzene.

2.5 Matrix Spike/Matrix Spike Duplicate

<u>WG288953</u>: Matrix Spike/Matrix Spike duplicate samples were not included in this sample delivery group. No qualification is needed.

<u>WG288783</u>: Matrix Spike/Matrix Spike duplicate samples were not included in this sample delivery group. No qualification is needed.

2.6 Surrogate Recovery

<u>WG288953</u>: All surrogate recoveries were within control limits for all environmental and quality control samples. No qualification is needed.

<u>WG288783:</u> All surrogate recoveries were within control limits for all environmental and quality control samples. No qualification is needed.

2.7 Internal Standards

All internal standard area counts were within control limits for all samples. No qualification is needed based on the internal standard information provided.

2.8 Duplicates

No sample was analyzed in duplicate. No qualification is needed.

2.8 Summary

The data are acceptable with validator-assigned qualifiers.

LABORATORY REPORT

L08110371

12/09/08 09:04

Submitted By

Microbac Laboratories Inc. 158 Starlite Drive Marietta, OH 45750 (740)373-4071

For

Account Name: Earth Tech, Inc 675 North Washington Street Suite 300 Alexandria, VA 22314 Attention: Devon Chicoine

Project Number: 2551.030 Project: Alex-Air Force Plant 59 Site: AFP59

	Sample Ana	lysis Su	mmary		
Client ID	L	ab ID	Method	Dilution	Date Received
TRIP BLANK	LÖ81	10371-01	8260B	1	14-NOV-08
59SG36-3.0-SO 1	L081	10371-02	8260B	1	14-NOV-08

L1_A_PROD - Modified 03/06/2008 PDF File ID:1270491 Report generated: 12/09/2008 09:04 1 OF 1



 PrcPrep Method:NONE
 Instrument:HPMS8

 Prep Method:5030B
 Prep Date:11/24/2008 23:55

 Analytical Method:8260B
 Cal Date:11/14/2008 18:53

 Analyst:MES
 Run Date:11/24/2008 23:55

 Dilution:1
 File ID:8M350016

 Units:ug/L
 File ID:8M350016
 Sample Number:L08110371-01 Client ID: TRIP BLANK Matrix: Water Workgroup Number: WG288953 Collect Date: 11/12/2008 00:01 Sample Tag:01
 Qual
 RL
 MDL

 U
 0.500
 0.250

 U
 1.00
 0.250

 U
 0.500
 0.125

 Analyte
 CAS. Number
 Result

 1,1,1,2-Tetrachloroethane
 630-20-6

 1,1,1-Trichloroethane
 71-55-6

 1,2,2-Tetrachloroethane
 79-34-5
 1,1,1-Trichloroethane 1,1,2,2-Tetrachloroethane 1,1,2-Trichloroethane 0.500 79-34-5 U 0.125 79-00-5 ΥT 1.00 0.250 ------
 1,1-Dichloroethane
 75-34-3

 1,1-Dichloroethene
 75-35-4

 1,1-Dichloropropene
 563-58-6
 1.00 0.125 ប 1.00 0.500 υ U U 0.250 1.00 1.....

1,1-Dichloropropene	0~02~00		U 1.00	
1,2,3-Trichlorobenzene	87-61-6			0.150
1,2,3-Trichloropropane	96-18-4		U 1.00	0.500
1,2,4-Trichlorobenzene	120-82-1	[U 1.00	0.200
1,2,4-Trimethylbenzene	95-63-6		υ 1.00	0.250
1,2-Dichloroethane	107-06-2		ປ 0.500	0.250
1,2-Dichlorobenzene	95-50-1		U 1.00	0.125
1.2-Dibromo-3-chloropropane	96-12-8		ບ 2.00	1.00
1.2-Dichloropropane	78-87-5		U 1.00	0.200
1,2-Dibromoethane	106-93-4		υ 1.00	0.250
1,3,5-Trimethylbenzene	108-67-8		U 1.00	0.250
1,3-Dichlorobenzene	541-73-1	<u> </u>	υ 1.00	0.250
1.3-Dichloropropane	142-28-9		U 0.400	0.200
1,4-Dichlorobenzene	106-46-7		U 0.500	0.125
1.4-Dioxane	123-91-1		U 100	50.0
1-Chlorohexane	544-10-5		υ 1.00	0.125
2.2-Dichloropropane	594-20-7		U 1.00	0.250
	95-49-8		υ 1.00	0.125
2-Chlorotoluene	106-43-4		u 1.00	0.250
4-Chlorotoluene	67-64-1	· [R () 10.0	2.50
Acetone	71-43-2	-	υ 0.400	0.125
Benzene	108-86-1	-	U 1.00	0.125
Bromobenzene	74-97-5		u 1.00	0.200
Bromochloromethane	a contract of the second se		υ 0.500	0.250
Bromodichloromethane	75-27-4		u 1.00	0.500
Bromoform	75-25-2		U 1.00	0.500
Bromomethane	74-83-9		υ <u>3.00</u> υ <u>1.00</u>	0,250
Carbon tetrachloride	56-23-5		U 0.500	0.125
Chlorobenzene	108-90-7		U 1.00	0.500
Chloroethane	75-00-3		U 0.300	0.125
Chloroform	67-66-3			0.250
Chloromethane	74-87-3		•	And the second s
cis-1,2-Dichloroethene	156-59-2		U 1.00	0.250
cis-1,3-Dichloropropene	10061-01-5		υ 0.500	0.250
Dibromochloromethane	124-48-1		B) 0.500	0.250
Dibromomethane	74-95-3		U 1.00	0.250
Dichlorodifluoromethane	75-71-8		U 1.00	0.250
Ethylbenzene	100-41-4		υ 1.00	0.250
Hexachlorobutadiene	87-68-3		U 0.600	0.250
Isopropylbenzene	98-82-8		U 1.00	0.250
Methylene chloride	75-09-2		U 1.00	0.250
Methyl t-butyl ether (MTBE)	1634-04-4		U 5.00	0.500
MEK (2-Butanone)	78-93-3		υ 10.0	2.50
MIBK (methyl isobutyl ketone)	108-10-1		υ 10.0	2.50
n-Butylbenzene	104-51-8		U 1.00	0.250
n-Propylbenzene	103-65-1		U 1.00	0.125
m-,p-Xylene	136777-61-2		ບ 2.00	0.500
Naphthalene	91-20-3		U 1.00	0.200
o-Xylene	95-47-6		U 1.00	0.250
	99-87-6		υ 1.00	0.250
p-Isopropyltoluene	135-98-8		U 1.00	0.250
sec-Butylbenzene	100-42-5		υ 1.00	0.125
Styrene	79-01-6	· · ·]	U 1.00	0.250
Trichloroethene	an an an an an ann ann an an an an an an		U 1.00	0.250
tert-Butylbenzene	98-06-6	J	0 T-00	. 0.200

of 4

1

12/12/08

Sample Number:L08110371-01	PrePrep Method:NONE	Instrument:HPMS8
Client ID: TRIP BLANK	Prep Method: 5030B	Prep Date:11/24/2008 23:55
Matrix:Water	Analytical Method:8260B	Cal Date:11/14/2008 18:53
Workgroup Number: WG288953	Analyst MES	Run Date:11/24/2008 23:55
Collect Date:11/12/2008 00:01	Dilution:1	File ID:8M350016
Sample Tag:01	Units:ug/L	
Analyte	CAS. Number Result	Qual RL MDL

	a se an esta de la companya de				
127-18-4	1		ប	1.00	0.250
108-88-3	i		υ.	1.00	: 0.250
156-60-5		1	U	1.00	0.250
10061-02-6			U	1.00	0.500
75-69-4	1	1	υ	1.00	0.250
75-01-4		1	υ	1.00	0.250
% Recoverv :	Lower	Upper		Qual	
98.1	85	115			
96.9	72	119		÷	
105	81	120	;		
103	76	119			
	127-18-4 108-88-3 156-60-5 10061-02-6 75-69-4 75-01-4 % Recovery 98.1 96.9 105	127-18-4 108-88-3 156-60-5 10061-02-6 75-69-4 75-01-4 % Recovery Lower 98.1 85 96.9 72 105 81	127-18-4 108-88-3 156-60-5 10661-02-6 75-69-4 75-01-4 % Recovery Lower Upper 98.1 85 115 96.9 72 119 105 81 120	127-18-4 U 108-88-3 U 156-60-5 U 10061-02-6 U 75-69-4 U 75-01-4 U % Recovery Lower Upper 98.1 85 115 96.9 72 119 105 81 120	127-18-4 U 1.00 108-88-3 U 1.00 156-60-5 U 1.00 10061-02-6 U 1.00 75-69-4 U 1.00 75-01-4 U 1.00 98.1 85 115 96.9 72 119 105 81 120

Undetected; the analyte was analyzed for, but not detected. R Because of quality control deficiences for this analyte, this data may be rejected.

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12/12/08

Sample Number: L08110371-02

Instrument: HPMS9 Prep Method: 5030BInstrument: HPMS9Prep Method: 5030BPrep Date: 11/19/2008 11:18Analytical Method: 8260BCal Date: 11/21/2008 20:29Analyst: TMBRun Date: 11/22/2008 22:53Dilution: 1File ID: 9M66225 Client ID: 595G36-3.0-SO 1 Matrix:Soil Workgroup Number:WG288783 kun Date:11/22 File ID:9M66225 Units:ug/kg Percent Solid:8 Collect Date: 11/12/2008 16:30 File ID:9M66225 Percent Solid:89.1 Sample Tag:01 MDL Analyte CAS. Number Result Qual 81. 7.15 U 3.58 67-64-1 26.8 Acetone Bromobenzene 0.358 3.58 71-43-2 Bromochloromethane Ü 0.358 3.58 108-86-1 3.58 0.358 74-97-5 Ü Bromochloromethane Bromodichloromethane ü 3.58 0.358 75-27-4 . <u>.....</u>..... υ 3.58 0.358 Bromoform Bromomethane 75-25-2 . j... 0.715 74-83-9 ΰ 7,15 Bromomethane 2 -Butanone n-Butylbenzene sec-Butylbenzene tert-Butylbenzene Carbon disulfide Carbon tetrachloride Chlorobenzene Chlorodibromomethane Chloroethane 78-93-3 IJ 7.15 1.79 . . . 104-51-8 ΰ 3.58 0.358 135-98-8 ú 3.58 0.358 . İ. 3.58 0.358 98-06-6 U . . . Ü 3.58 0.358 75-15-0 56-23-5 Ü 3.58 0.358 . . . 108-90-7 ŭ 3.58 0.358 75-00-3 110-75-8 0.358 IJ 3.58 7.15 0.715 Chloroethane 2-Chloroethyl vinyl ether Chloroform Chloromethane ri. 7.15 1.43 τī 3.58 0.358 67-66-3 π 7.15 1.43 Chloromethane 2-Chlorotoluene 4-Chlorotoluene 74-87-3 11 0.358 95-49-8 Ű 3.58 ::: 0.358 106-43-4 U 3.58 4-Chlorotoluene
 4-Chlorotoluene
 106-43-4

 1,2-Dibromo-3-chloropropane
 96-12-8

 1,2-Dibromoethane
 106-93-4

 Dibromomethane
 74-95-3

 05-50-1
 95-50-1
 1.43 U 3.58 ថ 3.58 0.358 3.58 U ò.358 0 200

PrePrep Method:NONE

1.2-Dichlorobenzene	95-50-1	U	3.58	0.358
1.3-Dichlorobenzene	541-73-1	U	3.58	0.358
1.4-Dichlorobenzene	106-46-7	υ	3.58	0.358
Dichlorodifluoromethane	75-71-8	U	7.15	0.715
1.1-Dichloroethane	75-34-3	ບ :	3.58	0.715
1,2-Dichloroethane	107-06-2	υ	3.58	0.358
1,1-Dichloroethene	75-35-4	υ	3.58	0.358
cis-1,2-Dichloroethene	156-59-2	U	3.58	0.358
trans-1,2-Dichloroethene	156-60-5	υ	3.58	0.358
1,2-Dichloropropane	78-87-5	υ	3,58	0.358
1,3-Dichloropropane	142-28-9	υ	3.58	0.358
2,2-Dichloropropane	594-20-7	U	3.58	0.358
cis-1,3-Dichloropropene	10061-01-5	U	3.58	0.358
trans-1, 3-Dichloropropene	10061-02-6	υ	3.58	0.358
1.1-Dichloropropene	563-58-6	U	3.58	0.358
Ethylbenzene	100-41-4	υ	3.58	0.358
2-Hexanone	591-78-6	U	7.15	1.79
Hexachlorobutadiene	87-68-3	ប	3.58	0.358
Isopropylbenzene	98-82-8	U	3.58	0.358
p-Isopropyltoluene	99-87-6	υ	3.58	0.358
4-Methyl-2-pentanone	108-10-1	ប	7.15	1.79
Methylene chloride	75-09-2 3.87		3.58	0.715
Naphthalene	91-20-3	U	7.15	0.358
n-Propylbenzene	103-65-1	U	3.58	0.358
Styrene	100-42-5	υ	3.58	0.358
1,1,1,2-Tetrachloroethane	630-20-6	U	3.58	0.358
1,1,2,2-Tetrachloroethane	79-34-5	υ	3.50	0.358
Tetrachloroethene	127-18-4	U	3.58	0.358
Toluene	108-88-3	U	3.58	0.358
1,2,3-Trichlorobenzene	87-61-6	ប	3.58	0.358
1,2,4-Trichlorobenzene	120-82-1	U	3.58	0.358
1,1,1-Trichloroethane	71-55-6 2.53		3.58	0.358
1,1,2-Trichloroethane	79-00-5	U	3.58	0.358
Trichloroethene	79-01-6 0.637	.F	3.58	0.358
Trichlorofluoromethane	75-69-4	υ	7.15	0.715

of

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Microbac

12/12/08

Sample Number:L08110371-02	PrePrep Method: NONE	Instrument: HPMS9
Client ID: 598636-3.0-80 1	Prep Method:5030B	Prep Date:11/19/2008 11:18
Matrix:Soil	Analytical Method:8260B	Cal Date:11/21/2008 20:29
Workgroup Number:WG288783	Analyst: TMB	Run Date:11/22/2008 22:53
Collect Date:11/12/2008 16:30	Dilution:1	File ID:9%66225
Sample Tag:01	Units:ug/kg	Percent Solid:89.1
		·
Analyte	CAS. Number Result	Qual RL MDL

96-18-4		1 U	3.58	0.715
95-63-6		i U	3.58	0.358
108-67~8		U	3.58	0.358
75-01-4		ប	7.15	0.715
95-47-6		1 11	3.58	0.358
336777-61-3	2	1 0	3.58	V.358
% Recoverv	Lower	Unper	Ouai	
: 103	65	135		
111	52	149		
98.5	84	116		
98.2	84	118		
	96-18-4 95-63-6 108-67-8 75-01-4 95-47-6 136777-61- % Recovery 103 111 98.5	96-18-4 95-63-6 108-67-8 75-01-4 95-47-6 136777-61-2 % Recovery Lower 103 65 111 52 98.5 84	96-18-4 U 95-63-6 U 108-67-8 U 75-01-4 U 95-63-6 U 108-67-8 U 95-63-6 U 98 136 111 52 98 94 98 94	95-63-6 U 3.58 108-67-8 U 3.58 75-01-4 U 7.15 95-47-6 U 3.58 136777-61-2 U 3.58 8 Recovery Lower Upper 103 65 135 111 52 149

U Undetected; the analyte was analyzed for, but not detected. F The analyte was positively identified, but the quantitation was below the RL.

Microbac

12/12/08

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