



**CONESTOGA-ROVERS  
& ASSOCIATES**

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April 29, 2008

Reference No. 012559

Ms. Laura C. Price  
Voluntary Cleanup Unit/Remedial Section  
Bureau of Environmental Remediation  
Kansas Department of Health & Environment  
1000 SW Jackson Street, Suite 410  
Topeka, Kansas 66612-1367

Dear Ms. Price:

Re: Groundwater Monitoring Results (Fourth Quarter 2007)  
Former General Motors Fairfax I Plant Site  
Kansas City, Kansas

## **1.0 INTRODUCTION**

### **1.1 GENERAL**

This monitoring report provides a summary of the groundwater monitoring program implemented by General Motors (GM) at the Former General Motors Fairfax I Plant (Site) in Kansas City, Kansas during the fourth quarter of 2007. This represents the 2<sup>nd</sup> event of the comprehensive groundwater sampling requested by the Kansas Department of Health and Environment (KDHE).

GM entered the former Fairfax I Site into the Voluntary Cleanup and Property Redevelopment Program (VCPRP) in March 14, 2001. Conestoga-Rovers & Associates (CRA) has performed numerous studies at the Site.

Field activities were performed in accordance with the scope of work and procedures outlined in the documents approved by KDHE and titled "Groundwater Monitoring Plan (Revised)" (CRA, July 2007) and "Pre-Design Investigations Work Plan" (CRA, August 2001).

### **1.2 SITE LOCATION AND DESCRIPTION**

The Site is located at 100 Kindelberger Road, Kansas City, Kansas and is located in the floodplain of the Missouri River. Refer to Figure 1.1 for the Site location. The Fairfax I Plant Site is vacant as operations ceased in 1986/87 and the facility was demolished in 1987. A number of monitoring wells, however, remain on the Site. Figure 1.2 presents the Site layout and the approximate locations of known existing monitoring wells. Table 1.1 presents well construction details for the wells in the monitoring well network.

REGISTERED COMPANY  
**ISO 9001**  
ENGINEERING DESIGN



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## **2.0 SCOPE OF WORK**

As outlined in the 2007 Groundwater Monitoring Plan (Revised), the field activities conducted during the week of December 3, 2007 consisted of the following tasks:

- Collection of depth to water measurements at the 32 wells in the monitoring well network;
- Collection of groundwater samples from 31 of the 32 wells in the monitoring well network; and
- Analysis of groundwater samples for target compound list volatile organic compounds (TCL VOCs) and total petroleum hydrocarbon as gasoline range organics (TPH-GRO).

## **3.0 FIELD ACTIVITES**

Field activities completed during this period consisted of a quarterly groundwater sampling event.

### **3.1 HYDRAULIC MONITORING**

Depth to water measurements were collected at the Site on December 3, 2007. Water levels were measured to the nearest 0.01 feet using an electronic water level indicator.

### **3.2 QUARTERLY GROUNDWATER SAMPLING AND ANALYSIS**

Groundwater sampling was conducted from December 4 to 6, 2007. Groundwater samples were collected from 31 of the 32 wells in the monitoring well network. Monitoring well MW-115 did not contain enough water to collect a sample. Field parameters were measured during monitoring well purging and recorded in the field logbook. Purging parameters are summarized in Table 3.1. Monitoring wells were purged using a submersible pump with dedicated tubing and using low-flow minimal drawdown techniques except at monitoring well MW-103A. Due to the presence of a strong odor at the well and the previous presence of light non-aqueous phase liquid (LNAPL) in the well, well MW-103A was purged of three well volumes of water (approximately 7 gallons) and sampled using a disposable bailer; no field parameters were measured. Groundwater samples were analyzed for TCL VOCs and TPH-GRO.

Groundwater samples were packed in a cooler with ice and shipped by overnight courier under chain-of-custody protocols to TestAmerica Laboratories, Inc. (TestAmerica) of North Canton, Ohio.



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#### **4.0 SUMMARY OF GROUNDWATER MONITORING RESULTS**

The following discussion relates to the quarterly groundwater sampling event and does not reference past groundwater sampling events whose results were proved in previous reports to the KDHE.

##### **4.1 HYDRAULIC MONITORING**

A summary of groundwater elevations is presented in Table 4.1. Figure 4.1 presents groundwater contours for the shallow depth monitoring wells, and Figure 4.2 presents groundwater contours for the intermediate depth monitoring wells.

##### **4.2 QUALITY ASSURANCE/QUALITY CONTROL**

An analytical data quality assessment and validation of the analytical results obtained by TestAmerica for the samples collected was conducted by CRA's quality assurance officer. The evaluation of the analytical data was based on a review of the information provided by TestAmerica, including laboratory blank data, as well as recovery data from matrix spike samples, surrogate compounds, and laboratory control samples. The analytical data were assessed for accuracy and precision based on the provided quality assurance data. Based on the results of the data quality assessment and validation, the analytical data are valid and suitable for quantitative purposes. The analytical data quality assessment and validation memorandum is provided in Attachment A.

##### **4.3 ANALYTICAL RESULTS**

A summary of detected analytical data for the current quarter is provided in Table 4.2. A summary of historical VOC detected analytical data for samples collected by CRA is provided in Attachment B (see Table B.1). The tables provide a comparison of the analytical data to both the residential and non-residential KDHE groundwater standards. The discussion of results, however, is limited to only those exceedences for the non-residential scenario. The property associated with the Fairfax I Site is now, and is expected to be in the future, industrially and commercially zoned. Consequently, the residential standards are provided for information purposes only.

Figure 1.2 provides monitoring well locations. Laboratory analytical results are provided in Attachment C.

During the sampling event, 4-methyl-2-pentanone, ethylbenzene, toluene, xylenes, and TPH-GRO were detected in the groundwater samples collected from MW-103A (sample and duplicate sample) at concentrations above the KDHE standards. Tetrachloroethene (PCE) was detected in the groundwater sample collected from MW-110A at a concentration slightly above the KDHE standard. Vinyl chloride and cis-1,2-dichloroethene (cis-1,2-DCE) were detected in



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the groundwater sample collected from MW-111 at concentrations slightly above the KDHE standard.

Results are consistent with previous sampling events. Exceedences of the KDHE groundwater standards were observed at three monitoring wells. Exceedences were observed at one location on the Site property where GM is currently planning and designing remedial actions (e.g., the MW-103A area).<sup>1</sup> One exceedence (PCE) was detected at a well that is off Site and generally upgradient of the Former Fairfax I Plant Site (MW-110A) and in an industrial area south of the Site.<sup>2</sup> The remaining exceedences (vinyl chloride and cis-1,2-DCE) were detected at a well (MW-111) historically considered to also be impacted from off-Site sources.<sup>2</sup>

Should you have any questions on the above, please do not hesitate to contact us.

Yours truly,

CONESTOGA-ROVERS & ASSOCIATES

A handwritten signature in black ink, appearing to read "Phil Harvey".

Phil Harvey

PH/lg/12

Encl.

c.c.: Ken Richards, GM

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<sup>1</sup> For the MW-103A area, refer to the work plan titled "Pre-Design Investigations Work Plan" dated August 2001.

<sup>2</sup> Refer to past reports titled "Comprehensive Groundwater Sampling Report", draft dated November 1998, and "Results of Soil and Groundwater Investigations Near Monitoring Well MW-103A", dated June 2000.

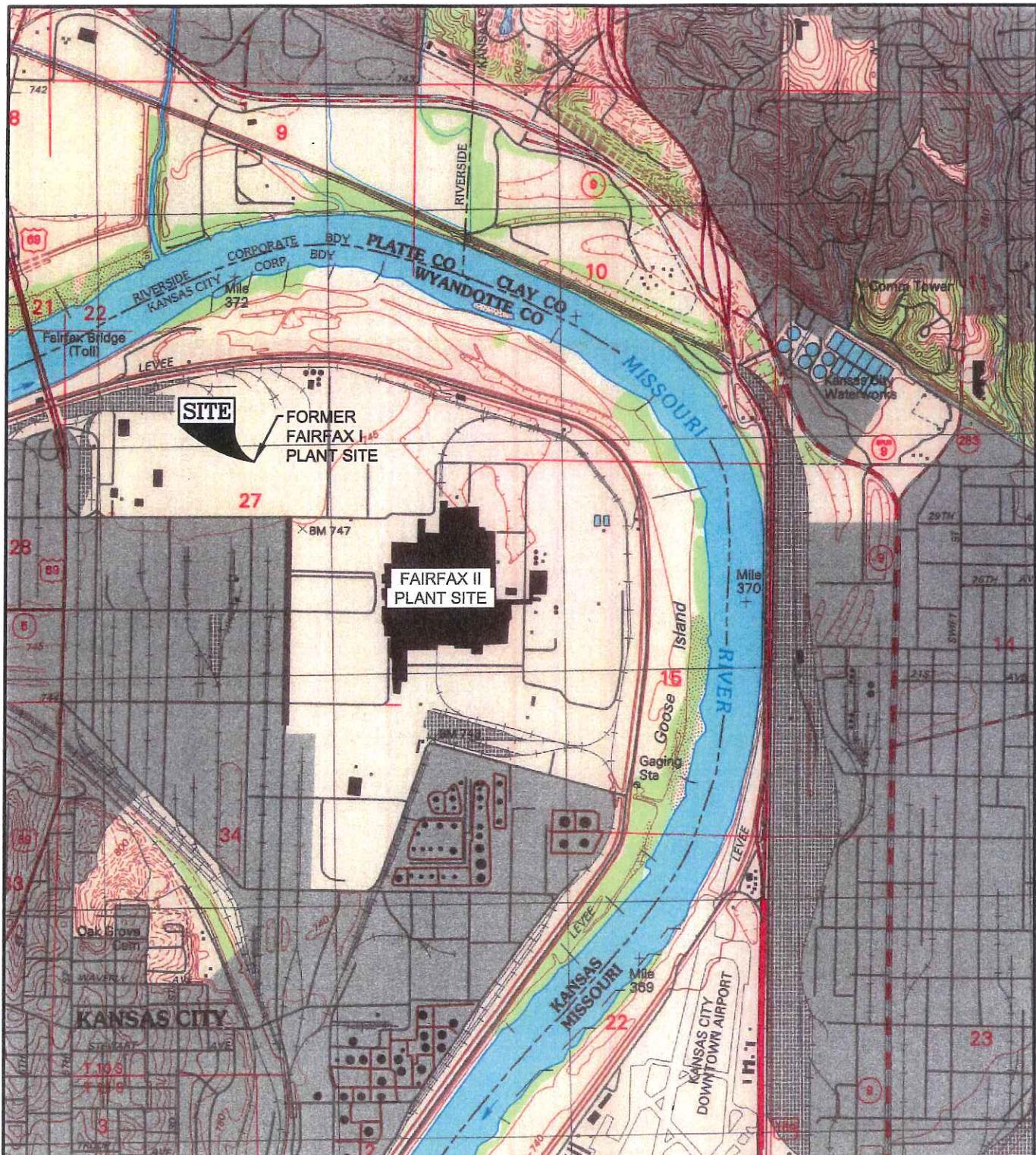
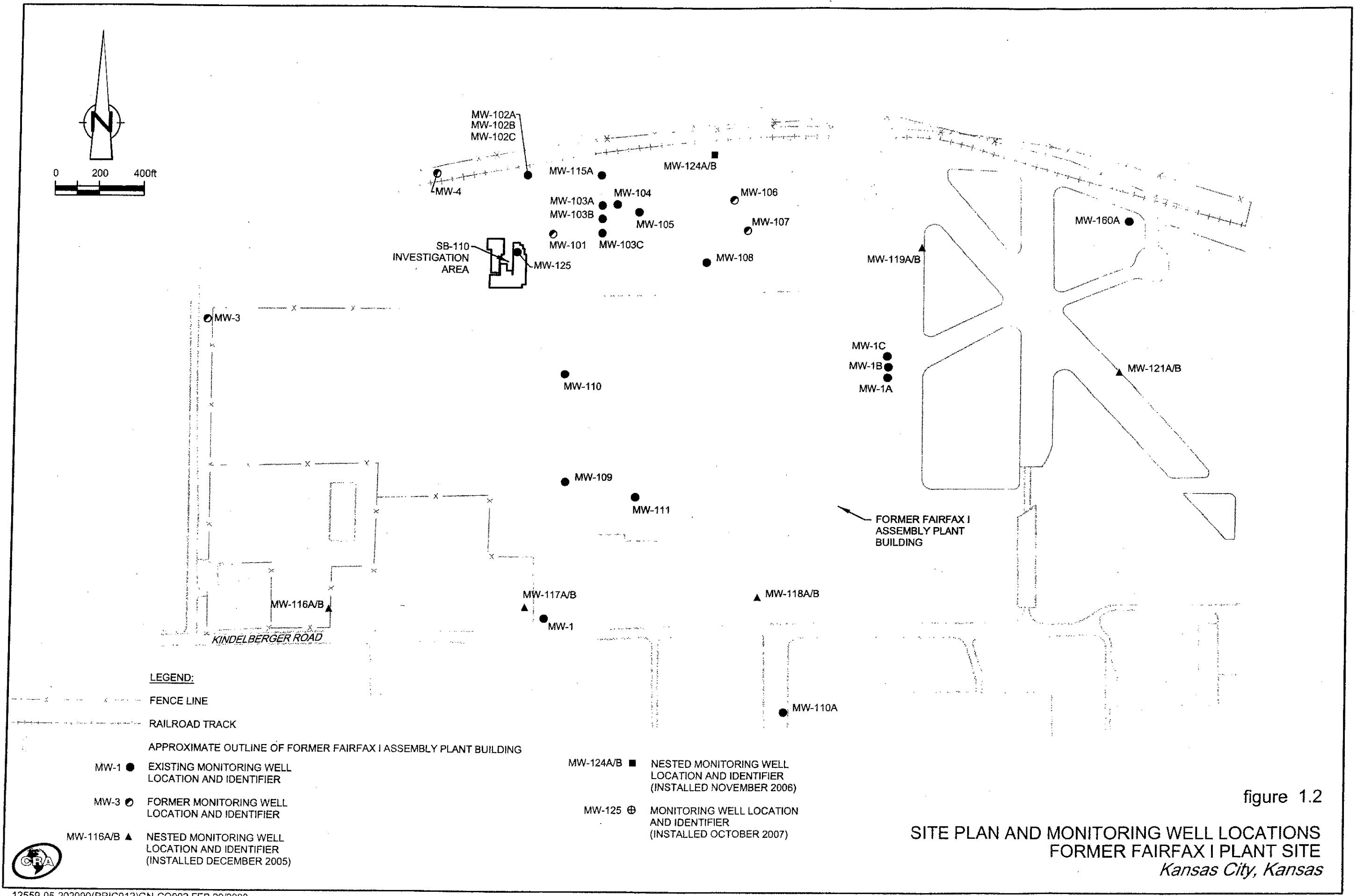


figure 1.1

SITE LOCATION MAP  
FORMER FAIRFAX I PLANT SITE  
*Kansas City, Kansas*



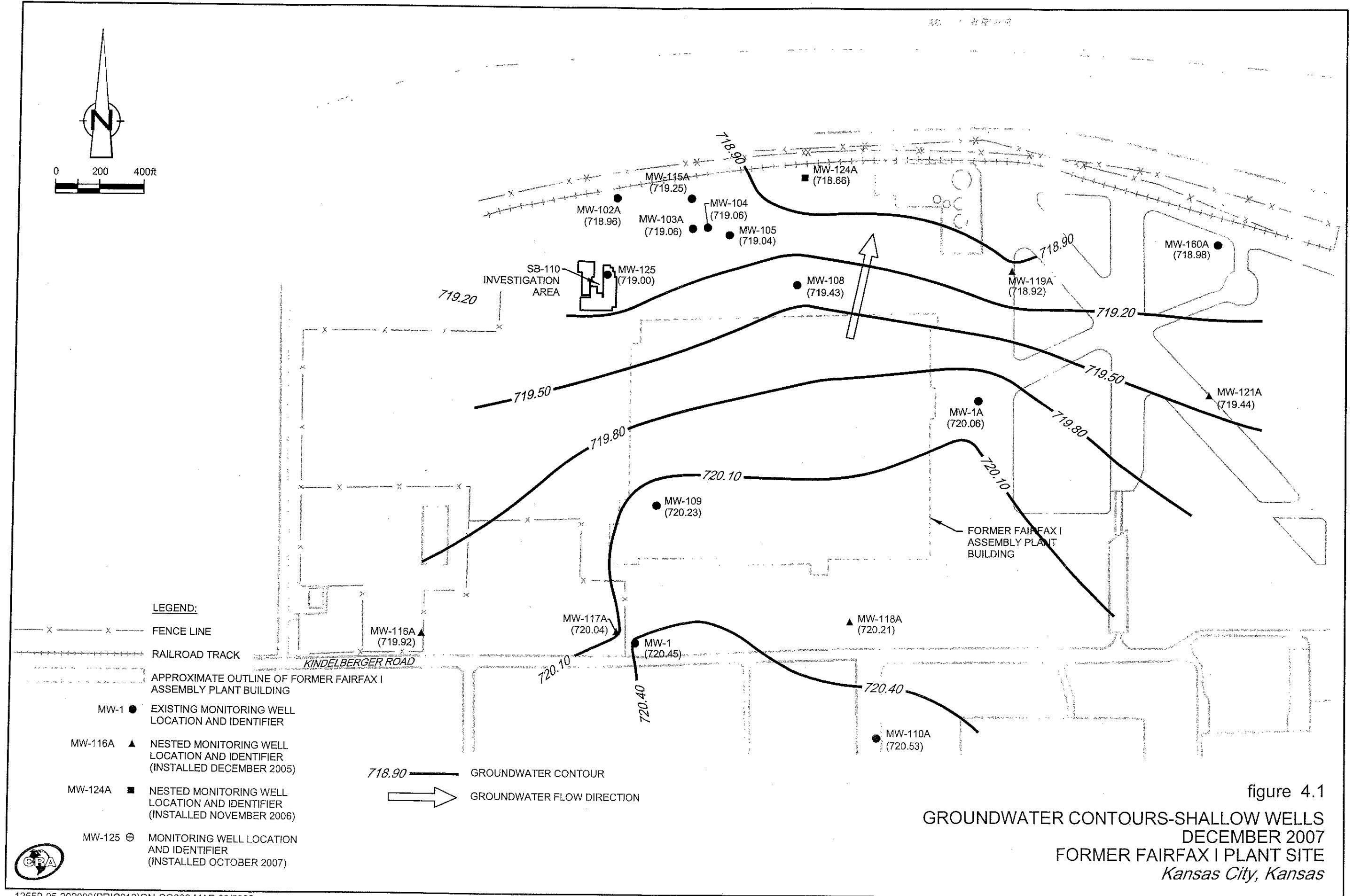


figure 4.1  
GROUNDWATER CONTOURS-SHALLOW WELLS  
DECEMBER 2007  
FORMER FAIRFAX I PLANT SITE  
*Kansas City, Kansas*

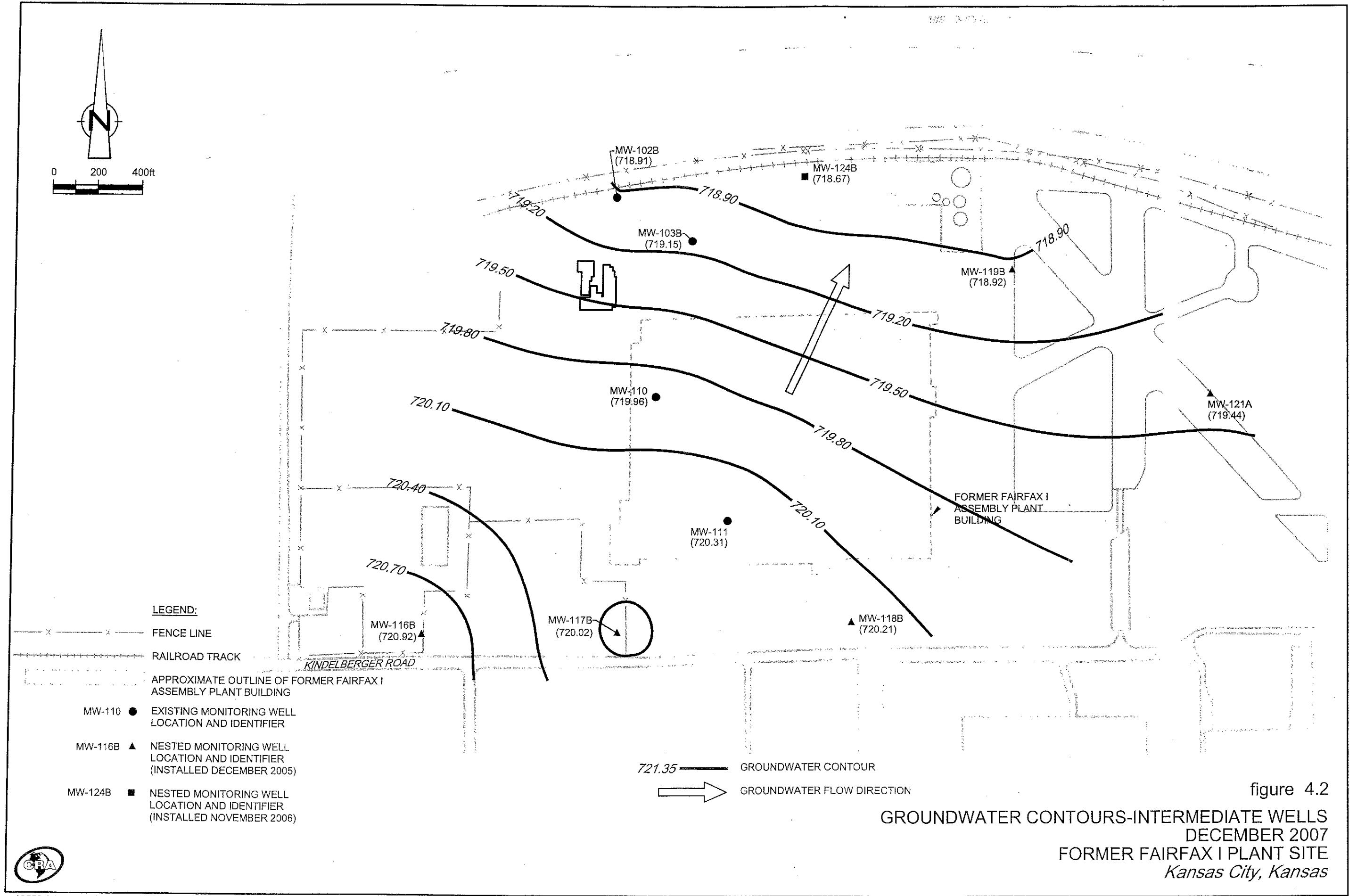


TABLE 1.1

**MONITORING WELL CONSTRUCTION DETAILS**  
**FORMER GM FAIRFAX I PLANT SITE**  
**KANSAS CITY, KANSAS**

Monitoring Well	Reference Elevation			Completion Date	Well Material <sup>2</sup>	Screened Interval			Elevation Bottom (ft AMSL)
	Top of Casing (ft AMSL) <sup>1</sup>	Ground Surface (ft AMSL)	Completion (ft AMSL)			Depth Top	Bottom (ft bgs)	(ft AMSL)	
MW-1	746.43	744.2	9/2/87	PVC	— <sup>4</sup>	30.0	—	—	714.2
MW-1A	745.93	743.7	3/27/85	PVC	15.0	35.0	—	728.7	708.7
MW-1B	745.40	743.6	3/27/85	PVC	65.0	85.0	—	678.6	658.6
MW-1C	746.13	743.6	4/1/85	PVC	87.0	107.0	—	656.6	636.6
MW-102A	748.82	746.1	1/9/88	PVC	—	29.3	—	—	716.9
MW-102B	749.15	746.1	1/13/88	PVC	—	61.2	—	—	684.9
MW-102C	749.20	746.2	1/12/88	PVC	—	94.0	—	—	652.2
MW-103A	747.71	745.1	1/88	SS	—	29.6	—	—	715.5
MW-103B	746.30	745.1	1/88	SS	—	61.4	—	—	683.8
MW-103C	745.90	745.1	1/88	SS	—	102.1	—	—	643.0
MW-104	746.51	744.8	1/10/88	PVC	—	28.4	—	—	716.4
MW-105	746.70	744.1	1/8/88	PVC	—	30.2	—	—	713.9
MW-108	747.17	744.0	1/6/88	PVC	—	28.5	—	—	715.5
MW-109	746.11	743.5	6/89	—	—	30.0	—	—	713.5
MW-110	745.79	743.8	4/90	—	—	61.2	—	—	682.6
MW-110A	743.89	743.5	—	—	—	36.1	—	—	707.4
MW-111	745.96	744.1	4/90	—	—	59.8	—	—	684.3
MW-115A	748.07	745.7	5/12/99	PVC	11.7	26.7	—	734.0	719.0
MW-116A	743.54	743.9	12/16/05	PVC	24.5	34.5	—	719.4	709.4
MW-116B	743.37	743.7	12/20/05	PVC	45.5	58.5	—	698.2	685.2
MW-117A	743.72	744.0	12/13/05	PVC	20.0	30.0	—	724.0	714.0
MW-117B	743.75	744.0	12/15/05	PVC	49.0	59.0	—	695.0	685.0
MW-118A	747.02	744.0	12/12/05	PVC	20.0	30.0	—	724.0	714.0

TABLE 1.1

**MONITORING WELL CONSTRUCTION DETAILS**  
**FORMER GM FAIRFAX I PLANT SITE**  
**KANSAS CITY, KANSAS**

Monitoring Well	Reference Elevation		Completion Date	Well Material <sup>2</sup>	Screened Interval		
	Top of Casting (ft AMSL) <sup>1</sup>	Ground Surface (ft AMSL)			Depth Top	Bottom (ft bgs) <sup>3</sup>	Elevation Top (ft AMSL)
MW-118B	746.86	744.0	12/13/05	PVC	48.5	58.5	695.5
MW-119A	746.69	743.7	12/20/05	PVC	23.5	33.5	720.2
MW-119B	746.65	743.6	12/21/05	PVC	49.0	59.0	694.6
MW-121A	745.53	742.8	12/22/05	PVC	25.0	35.0	717.8
MW-121B	745.80	742.9	12/22/05	PVC	48.5	58.5	694.4
MW-124A	747.22	744.8	11/10/06	PVC	25.0	35.0	719.8
MW-124B	747.15	744.6	11/10/06	PVC	50.0	60.0	694.6
MW-125	745.34	742.3	10/31/07	PVC	25.0	35.0	717.3
MW-160A	746.56	744.3	3/87	—	—	35.4	—
							708.9

<sup>1</sup> ft AMSL - feet above mean sea level

<sup>2</sup> PVC - polyvinyl chloride; SS - stainless steel

<sup>3</sup> ft bgs - feet below ground surface

<sup>4</sup> — information not available

Screened interval data is not available for MW-1, MW-102A/B/C, MW-103A/B/C, MW-104, MW-105, MW-108, MW-10, MW-110, MW-110A, MW-111, and MW-160A. The approximate bottom of screen elevation was calculated using the difference between the surveyed top of casing elevation and the measured well total depth from 12/3/07; the approximate bottom of screen depth below ground surface was calculated using the difference between the surveyed top of casing and ground surface elevations and the measured total depth from 12/3/07.

TABLE 3.1

**SUMMARY OF MONITORING WELL PURGING PARAMETERS**  
**FORMER FAIRFAX I PLANT SITE**  
**KANSAS CITY, KANSAS**

<b>Location</b>	<b>Date</b>	<b>Time Purged (minutes)</b>	<b>Depth to Water (ft BGS)<sup>1</sup></b>	<b>pH (S.U.)<sup>2</sup></b>	<b>Temperature (°C)<sup>3</sup></b>	<b>Conductivity (µS/cm)<sup>4</sup></b>	<b>Dissolved</b>		<b>Turbidity (NTU)<sup>7</sup></b>	<b>Other</b>
							<b>Oxygen (mg/L)<sup>5</sup></b>	<b>ORP (mV)<sup>6</sup></b>		
MW-1	12/4/07	35	26.00	6.63	18.8	1,303	0.37	6.9	2.50	Clear
		40	26.00	6.62	18.8	1,306	0.35	9.6	1.22	Clear
		45	26.00	6.62	18.8	1,306	0.30	10.2	1.12	Clear
MW-1A	12/6/07	30	25.91	6.90	15.2	941	0.53	-98.7	3.50	Clear
		35	25.91	6.87	15.2	943	0.17	-106.0	3.55	Clear
		40	25.91	6.87	15.2	943	0.11	-107.6	2.94	Clear
MW-1B	12/6/07	10	23.43	7.13	14.2	940	0.13	-120.3	58.5	Cloudy
		15	23.43	7.13	14.1	937	0.10	-124.8	30.2	Cloudy
		20	23.43	7.12	14.1	937	0.08	-128.0	20.8	Clear
MW-1C	12/5/07	25	23.43	7.13	14.1	938	0.06	-130.5	7.08	Clear
		30	23.44	7.13	14.1	939	0.07	-132.0	4.03	Clear
		35	26.08	7.07	14.6	943	0.19	-90.1	1.13	Clear
MW-102A	12/5/07	15	26.08	7.07	14.6	943	0.17	-94.6	0.74	Clear
		20	26.08	7.07	14.6	943	0.17	-94.6	0.74	Clear
		25	26.08	7.09	14.6	944	0.14	-99.3	0.71	Clear
MW-102B	12/5/07	10	29.89	6.65	14.6	1,151	0.30	-3.0	56.3	Brown, slightly cloudy
		20	29.89	6.64	14.5	1,153	0.19	4.0	39.9	Brown, slightly cloudy
		30	29.89	6.61	14.5	1,158	0.16	6.0	15.5	Clear
MW-102C	12/5/07	35	29.89	6.60	14.6	1,156	0.13	6.8	7.3	Clear
		40	29.89	6.60	14.6	1,156	0.14	7.2	4.91	Clear
		25	30.25	7.26	14.5	981	0.19	-100.7	1.39	Clear
	10	30	30.25	7.27	14.5	982	0.16	-101.9	0.95	Clear
		15	30.25	7.15	14.2	959	0.09	-109.3	1.65	Clear
		20	30.24	7.15	14.2	959	0.08	-113.5	0.78	Clear
	10	35	30.25	7.29	14.5	981	0.08	-121.9	--	Clear
		15	30.25	7.15	14.3	959	0.09	-109.3	1.65	Clear
		20	30.24	7.15	14.2	959	0.08	-115.8	0.49	Clear

TABLE 3.1

**SUMMARY OF MONITORING WELL PURGING PARAMETERS**  
**FORMER FAIRFAX PLANT SITE**  
**KANSAS CITY, KANSAS**

Location	Date	Time Purged (minutes)	Depth to Water (ft BGS) <sup>1</sup>	pH (S.U.) <sup>2</sup>	Temperature (°C) <sup>3</sup>	Conductivity (µS/cm) <sup>4</sup>	Dissolved Oxygen (mg/L) <sup>5</sup>	ORP (mV) <sup>6</sup>	Turbidity (NTU) <sup>7</sup>	Other
MW-103A 12/6/07										
MW-103B	12/5/07	10	27.15	7.15	14.6	1,131	0.19	-121.4	1.49	Clear
		15	27.15	7.16	14.7	1,130	0.14	-126.6	1.86	Clear
		20	27.15	7.17	14.7	1,127	0.10	-13.5	0.97	Clear
MW-103C	12/5/07	20	26.78	7.16	14.5	991	0.10	-101.3	3.03	Clear
		25	26.78	7.14	14.5	989	0.15	-110.9	2.33	Clear
		30	26.78	7.14	14.5	991	0.10	-114.4	2.00	Clear
MW-104	12/5/07	5	27.65	6.72	14.5	1,817	1.01	-101.3	10.1	Clear
		10	27.66	6.73	14.7	1,837	0.20	-121.8	7.61	Clear
		15	27.66	6.74	14.7	1,838	0.17	-124.9	6.11	Clear
		20	27.66	6.74	14.7	1,833	0.13	-129.1	4.80	Clear
MW-105	12/5/07	15	27.71	6.70	14.6	1,325	0.69	-7.1	0.69	Clear
		20	27.71	6.68	14.5	1,324	0.66	-11.9	0.66	Clear
		25	27.72	6.67	14.5	1,325	0.59	-18.7	0.59	Clear
MW-108	12/5/07	5	27.69	6.66	14.4	1,255	0.47	7.7	1.45	Clear
		10	27.69	6.63	14.4	1,259	0.33	12.7	1.20	Clear
		15	27.69	6.61	14.4	1,260	0.27	16.5	1.12	Clear
MW-109	12/6/07	15	25.91	6.69	15.2	1,308	0.97	-16.7	4.07	Clear
		20	25.91	6.62	15.3	1,340	0.71	-10.0	1.03	Clear
		25	25.91	6.60	15.3	1,362	0.54	-5.2	0.74	Clear
		30	25.91	6.58	15.3	1,368	0.44	-0.9	0.51	Clear
MW-110	12/6/07	5	25.85	7.09	13.8	982	0.27	-98.2	4.25	Clear
		10	25.86	7.09	13.9	989	0.20	-109.2	3.79	Clear
		15	25.86	7.10	13.9	1,000	0.11	-125.1	1.81	Clear
		20	25.86	7.10	13.9	1,001	0.09	-126.6	0.88	Clear

**SUMMARY OF MONITORING WELL PURGING PARAMETERS**  
**FORMER FAIRFAX I PLANT SITE**  
**KANSAS CITY, KANSAS**

Location	Date	Time Purged (minutes)	Depth to Water (ft BGS) <sup>1</sup>	pH (S.U.) <sup>2</sup>	Temperature (°C) <sup>3</sup>	Conductivity (µS/cm) <sup>4</sup>	Dissolved Oxygen (mg/L) <sup>5</sup>	ORP (mV) <sup>6</sup>	Turbidity (NTU) <sup>7</sup>	Other
MW-110A	12/5/07	10	23.42	6.69	16.3	1,185	0.84	-10.3	11.2	Clear
		15	23.42	6.69	16.3	1,191	0.57	-4.5	5.76	Clear
		20	23.43	6.67	16.3	1,196	0.50	-0.5	4.55	Clear
MW-111	12/6/07	5	25.66	6.88	13.4	1,060	0.54	-47.8	8.52	Clear
		10	25.66	6.84	13.8	1,095	0.31	-69.0	10.6	Clear
		15	25.66	6.84	13.8	1,113	0.22	-91.5	6.51	Clear
		20	25.66	6.85	13.8	1,115	0.12	-102.2	4.53	Clear
MW-116A	12/4/07	10	23.68	6.61	17.0	1,524	2.63	130.9	4.33	Clear
		15	23.68	6.60	17.0	1,527	2.71	130.8	3.47	Clear
		20	23.69	6.59	17.0	1,522	2.71	130.4	2.91	Clear
MW-116B	12/4/07	10	23.56	7.07	17.3	1,072	3.63	-79.7	3.84	Clear
		15	23.56	7.04	17.3	1,078	4.31	-97.2	3.89	Clear
		20	23.57	6.98	17.4	1,108	4.25	-108.6	2.29	Clear
		25	23.57	6.94	17.4	1,112	4.15	-110.6	2.04	Clear
		30	23.55	6.91	17.3	1,116	4.09	-110.8	2.10	Clear
MW-117A	12/4/07	15	23.69	6.63	18.3	1,281	1.83	-20.5	62.3	Slightly cloudy
		20	23.69	6.54	18.4	1,281	1.49	-14.2	22.5	Clear
		25	23.70	6.53	18.4	1,278	1.34	-9.9	10.4	Clear
		30	23.70	6.54	18.4	1,278	1.19	-5.2	5.74	Clear
		35	23.70	6.54	18.4	1,276	1.18	-4.2	3.88	Clear
MW-117B	12/4/07	25	23.75	6.67	16.7	1,059	0.62	-99.7	18.7	Clear
		30	23.75	6.75	16.8	1,059	0.61	-100.3	17.4	Clear
		35	23.75	6.79	16.7	1,056	0.49	-110.4	9.73	Clear
		40	23.75	6.82	16.7	1,057	0.41	-115.3	5.48	Clear
		45	23.75	6.84	16.7	1,058	0.37	-117.5	4.49	Clear

**SUMMARY OF MONITORING WELL PURGING PARAMETERS**  
**FORMER FAIRFAX I PLANT SITE**  
**KANSAS CITY, KANSAS**

Location	Date	Time Purged (minutes)	Depth to Water (ft BGS) <sup>1</sup>	pH (S.U.) <sup>2</sup>	Temperature (°C) <sup>3</sup>	Conductivity (µS/cm) <sup>4</sup>	Dissolved Oxygen (mg/L) <sup>5</sup>	ORP (mV) <sup>6</sup>	Turbidity (NTU) <sup>7</sup>	Other
MW-118A	12/4/07	5	26.83	6.74	15.2	1,182	1.03	36.1	148	Brown, cloudy
		15	26.83	6.65	15.4	1,194	0.76	44.9	31.0	Clear
		20	26.83	6.64	15.4	1,196	0.63	48.2	17.3	Clear
		25	26.84	6.64	15.5	1,196	0.56	51.3	11.8	Clear
		30	26.84	6.64	15.5	1,196	0.56	52.6	9.03	Clear
		35	26.84	6.64	15.5	1,195	0.57	54.3	6.58	Clear
		40	26.84	6.63	15.5	1,196	0.57	56.1	4.89	Clear
MW-118B	12/4/07	10	26.68	6.87	15.1	1,000	0.36	-96.2	3.12	Clear
		15	26.68	6.85	15.2	1,000	0.28	-102.2	2.84	Clear
		20	26.69	6.91	15.2	999	0.17	-119.5	2.91	Clear
		25	26.68	6.92	15.2	998	0.12	-125.8	2.71	Clear
MW-119A	12/6/07	30	27.80	6.71	16.5	1,338	0.39	-60.5	0.93	Clear
		35	27.80	6.70	16.5	1,340	0.34	-59.6	0.81	Clear
		40	27.87	6.69	16.5	1,340	0.29	-58.5	0.85	Clear
MW-119B	12/6/07	15	27.81	7.30	14.2	1,001	0.57	-102.3	12.4	Clear
		20	27.81	7.15	14.5	1,026	0.25	-98.9	9.60	Clear
		25	27.81	7.14	15.1	1,066	0.17	-115.2	10.7	Clear
		30	27.81	7.14	14.3	1,040	0.25	-113.2	22.1	Cloudy
		35	27.81	7.14	15.1	1,060	0.49	-119.1	227	Gray, cloudy
		60	27.85	7.14	15.0	1,058	0.13	-129.7	10.2	Clear
		65	27.83	7.14	15.1	1,062	0.12	-132.7	5.64	Clear
		70	27.81	7.14	15.0	1,062	0.11	-134.1	3.53	Clear
MW-121A	12/4/07	10	26.11	6.49	15.9	1,377	0.17	29.1	7.88	Clear
		15	26.11	6.50	15.9	1,371	0.13	23.4	4.24	Clear
		20	26.11	6.48	15.9	1,373	0.10	19.8	2.47	Clear

**SUMMARY OF MONITORING WELL PURGING PARAMETERS**  
**FORMER FAIRFAX I PLANT SITE**  
**KANSAS CITY, KANSAS**

Location	Date	Time Purged (minutes)	Depth to Water (ft BGS) <sup>1</sup>	pH (S.U.) <sup>2</sup>	Temperature (°C) <sup>3</sup>	Conductivity (µS/cm) <sup>4</sup>	Dissolved Oxygen (mg/L) <sup>5</sup>	ORP (mV) <sup>6</sup>	Turbidity (NTU) <sup>7</sup>	Other
MW-121B	12/4/07	15	26.37	6.88	15.3	1,014	0.09	-116.4	2.47	Clear
		20	26.37	6.89	15.3	1,013	0.07	-122.6	1.74	Clear
		25	26.37	6.92	15.3	1,012	0.07	-127.5	1.26	Clear
MW-124A	12/5/07	10	28.57	6.60	14.9	1,369	0.25	-80.9	15.0	Clear
		15	28.57	6.63	15.0	1,376	0.17	-87.7	5.79	Clear
		20	28.57	6.63	15.0	1,397	0.15	-92.6	3.00	Clear
		25	28.58	6.62	15.0	1,381	0.15	-94.6	2.87	Clear
MW-124B	12/5/07	20	28.49	7.03	14.7	1,239	5.35	-109.1	9.50	Clear
		25	28.49	7.02	14.7	1,239	5.29	-109.2	6.99	Clear
		30	28.49	6.99	14.8	1,241	4.56	-111.5	4.51	Clear
MW-125A	12/6/07	5	26.34	6.69	14.5	1,301	0.83	-12.5	45.2	Slightly cloudy
		10	26.34	6.63	14.1	1,254	0.21	-31.5	28.1	Clear
		15	26.34	6.65	14.2	1,243	0.13	-43.5	10.9	Clear
		20	23.34	6.66	14.2	1,242	0.11	-51.8	6.31	Clear
		25	23.34	6.67	14.2	1,239	0.09	-56.8	4.43	Clear
MW-160A	12/4/07	5	27.60	6.54	16.4	1,347	1.12	-21.4	56.7	Brown, cloudy
		10	27.61	6.51	16.5	1,343	0.59	-30.3	18.5	Clear
		15	27.61	6.52	16.5	1,352	0.32	-37.8	10.1	Clear
		20	27.60	6.52	16.4	1,353	0.21	-41.9	6.27	Clear
		25	27.60	6.52	16.4	1,354	0.16	-45.2	4.26	Clear

<sup>1</sup> ft BGS - feet below ground surface

<sup>2</sup> S.U. - standard units

<sup>3</sup> °C - degrees Celsius

<sup>4</sup> µS/cm - microsiemens per centimeter

<sup>5</sup> mg/L - milligrams per liter

<sup>6</sup> mV - millivolts

<sup>7</sup> NTU - nephelometric turbidity units

**TABLE 4.1**  
**SUMMARY OF GROUNDWATER ELEVATIONS**  
**FORMER FAIRFAX I PLANT SITE**  
**KANSAS CITY, KANSAS**

<i>Monitoring Well</i>	<i>Top of Casing Elevation (ft AMSL)<sup>1</sup></i>	<i>September 2007</i>		<i>December 2007</i>	
		<i>Depth to Water (ft BTOC)<sup>2</sup></i>	<i>Groundwater Elevation (ft AMSL)</i>	<i>Depth to Water (ft BTOC)</i>	<i>Groundwater Elevation (ft AMSL)</i>
MW-1	746.43	24.13	722.30	25.98	720.45
MW-1A	745.93	23.85	722.08	25.87	720.06
MW-1B	745.40	23.30	722.10	25.36	720.04
MW-1C	746.13	24.00	722.13	26.06	720.07
MW-102A	748.82	26.45	722.37	29.86	718.96
MW-102B	749.15	26.60	722.55	30.24	718.91
MW-102C	749.20	26.80	722.40	30.24	718.96
MW-103A	747.71	25.21	722.50	28.65	719.06
MW-103B	746.30	23.78	722.52	27.15	719.15
MW-103C	745.90	23.38	722.52	26.75	719.15
MW-104	746.51	24.03	722.48	27.45	719.06
MW-105	746.70	24.41	722.29	27.66	719.04
MW-108	747.17	24.91	722.26	27.74	719.43
MW-109	746.11	23.85	722.26	25.88	720.23
MW-110	745.79	23.43	722.36	25.83	719.96
MW-110A	743.89	22.25	721.64	23.36	720.53
MW-111	745.96	23.86	722.10	25.65	720.31
MW-115A	748.07	25.56	722.51	28.82	719.25
MW-116A	743.54	21.18	722.36	23.62	719.92
MW-116B	743.37	21.00	722.37	22.45	720.92
MW-117A	743.72	21.66	722.06	23.68	720.04
MW-117B	743.75	21.73	722.02	23.73	720.02
MW-118A	747.02	25.86	721.16	26.81	720.21
MW-118B	746.86	25.70	721.16	26.65	720.21
MW-119A	746.69	24.87	721.82	27.77	718.92
MW-119B	746.65	24.84	721.81	27.73	718.92
MW-121A	745.53	24.00	721.53	26.09	719.44
MW-121B	745.80	24.25	721.55	26.36	719.44
MW-124A	747.22	25.13	722.09	28.56	718.66
MW-124B	747.15	25.06	722.09	28.48	718.67
MW-125	745.34	22.54	722.80 <sup>3</sup>	26.34	719.00
MW-160A	746.56	24.52	722.04	27.58	718.98

<sup>1</sup> ft AMSL - feet above mean sea level

<sup>2</sup> ft BTOC - feet below top of casing

<sup>3</sup> MW-125 - depth to water measured November 1, 2007

TABLE 4.2

**SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS**

Sample Location:	MW-1	MW-1A	MW-1B	MW-1B	MW-1C	MW-102A	MW-102B	MW-102C	MW-103A	MW-103A	MW-103B	MW-103B	MW-103C
	GW-120407-JH-005	GW-120607-JH-029	GW-120607-JH-027	GW-120607-JH-028	GW-120507-JH-024	GW-120507-JH-014	GW-120507-JH-015	GW-120507-JH-016	GW-120607-JH-030	GW-120607-JH-031	GW-120607-JH-031	GW-120507-JH-021	GW-120507-JH-020
Sample Date:	12/4/2007	12/6/2007	12/6/2007	12/6/2007	12/5/2007	12/5/2007	12/5/2007	12/5/2007	12/6/2007	12/6/2007	12/6/2007	12/5/2007	12/5/2007
Sample Type:	Residential GW Pathway Scenario <sup>1</sup>	Non-Residential GW Pathway Scenario <sup>1</sup>			Duplicate					Duplicate			
Parameters	a	b											
<b>Volatile Organic Compounds (mg/L)<sup>2</sup></b>													
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	0.07	0.23	0.01 U <sup>3</sup>	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	160 <sup>ab5</sup>	160 <sup>ab</sup>	0.01 U	0.01 U
Chlorobenzene	0.1	0.1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	5 U	5 U	0.001 U	0.001 U
cis-1,2-Dichloroethene	0.07	0.07	0.0005 U	0.00036 J <sup>4</sup>	0.0005 U	2.5 U	2.5 U	0.0064	0.0005 U				
Ethylbenzene	0.7	0.7	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	8 <sup>ab</sup>	7.5 <sup>b</sup>	0.001 U	0.001 U
Methylene chloride	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	5 U	5 U	0.001 U	0.001 U
Tetrachloroethene	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	150 <sup>ab</sup>	140 <sup>ab</sup>	0.00026 J	0.001 U
Toluene	1	1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	2.5 U	2.5 U	0.0005 U	0.0005 U
trans-1,2-Dichloroethene	0.1	0.1	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	5 U	5 U	0.001 U	0.001 U
Trichloroethene	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	5 U	5 U	0.001	0.001 U
Vinyl chloride	0.002	0.002	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	57 <sup>ab</sup>	56 <sup>ab</sup>	0.001 U	0.001 U
Xylene (total)	10	10	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	470 <sup>ab</sup>	540 <sup>ab</sup>	0.1 U	0.1 U
<b>Petroleum Products</b>													
Total Petroleum Hydrocarbons - purgeable (GRO)	0.500	0.500	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U	470 <sup>ab</sup>	540 <sup>ab</sup>	0.1 U	0.1 U

TABLE 4.2

**SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS**

<i>Sample Location:</i>	<i>MW-104</i>	<i>MW-105</i>	<i>MW-108</i>	<i>MW-109</i>	<i>MW-110</i>	<i>MW-110</i>	<i>MW-110A</i>	<i>MW-111</i>	<i>MW-111</i>	<i>MW-116A</i>	<i>MW-116B</i>	<i>MW-117A</i>
	<i>GW-120507-JH-019</i>	<i>GW-120507-JH-018</i>	<i>GW-120507-JH-017</i>	<i>GW-120607-JH-033</i>	<i>GW-120607-JH-037</i>	<i>GW-120607-JH-038</i>	<i>GW-120507-JH-023</i>	<i>GW-120607-JH-035</i>	<i>GW-120607-JH-036</i>	<i>GW-120407-JH-001</i>	<i>GW-120407-JH-002</i>	<i>GW-120407-JH-003</i>
<i>Sample Date:</i>												
<i>Sample Type:</i>												
	<i>Residential</i>	<i>Non-Residential</i>										
	<i>GW Pathway</i>	<i>GW Pathway</i>										
	<i>Scenario<sup>1</sup></i>	<i>Scenario<sup>1</sup></i>										
<i>Parameters</i>	<i>a</i>	<i>b</i>										
<i>Volatile Organic Compounds (mg/L)<sup>2</sup></i>												
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	0.07	0.23	0.01 U	0.033 U	0.033 U	0.01 U	0.01 U	0.01 U				
Chlorobenzene	0.1	0.1	0.001 U	0.0033 U	0.0033 U	0.001 U	0.001 U	0.001 U				
cis-1,2-Dichloroethene	0.07	0.07	0.00022 J	0.0037	0.0005 U	0.02	0.00033 J	0.00035 J	0.0019	0.086 <sup>b</sup>	0.089 <sup>b</sup>	0.0005 U
Ethylbenzene	0.7	0.7	0.001 U	0.0033 U	0.0033 U	0.001 U	0.001 U					
Methylene chloride	0.005	0.005	0.001 U	0.0017 J	0.0028 J	0.001 U	0.001 U	0.001 U				
Tetrachloroethene	0.005	0.005	0.001 U	0.00035 J	0.001 U	0.001 U	0.001 U	0.0096 <sup>b</sup>	0.0033 U	0.0033 U	0.001 U	0.001 U
Toluene	1	1	0.001 U	0.0033 U	0.0033 U	0.00022 J	0.001 U					
trans-1,2-Dichloroethene	0.1	0.1	0.0005 U	0.0005 U	0.0005 U	0.0015	0.0005 U	0.0005 U	0.0062	0.006	0.0005 U	0.0005 U
Trichloroethene	0.005	0.005	0.001 U	0.00075 J	0.001 U	0.001 U	0.001 U	0.0014	0.0033 U	0.0033 U	0.001 U	0.001 U
Vinyl chloride	0.002	0.002	0.001 U	0.001 U	0.001 U	0.00029 J	0.001 U	0.001 U	0.001 U	0.007 <sup>b</sup>	0.0067 <sup>b</sup>	0.001 U
Xylene (total)	10	10	0.001 U	0.0033 U	0.0033 U	0.001 U	0.001 U					
<i>Petroleum Products</i>												
Total Petroleum Hydrocarbons - purgeable (GRO)	0.500	0.500	0.1 U	0.13 U	0.12 U	0.1 U	0.1 U	0.1 U				

TABLE 4.2

**SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS**

<i>Sample Location:</i>	MW-117B	MW-118A	MW-118B	MW-119A	MW-119B	MW-121A	MW-121B	MW-124A	MW-124B	MW-125	MW-160A
<i>Sample ID:</i>	GW-120407-JH-004	GW-120407-JH-006	GW-120407-JH-007	GW-120607-JH-026	GW-120607-JH-025	GW-120407-JH-008	GW-120407-JH-009	GW-120507-JH-011	GW-120507-JH-012	GW-120607-JH-032	GW-120407-JH-010
<i>Sample Date:</i>	12/4/2007	12/4/2007	12/4/2007	12/6/2007	12/6/2007	12/4/2007	12/4/2007	12/5/2007	12/5/2007	12/6/2007	12/4/2007
<i>Sample Type:</i>	<i>Residential</i> <i>GW Pathway Scenario<sup>1</sup></i>	<i>Non-Residential</i> <i>GW Pathway Scenario<sup>1</sup></i>									
<i>Parameters</i>	a	b									
<i>Volatile Organic Compounds (mg/L)<sup>2</sup></i>											
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	0.07	0.23	0.01 U								
Chlorobenzene	0.1	0.1	0.001 U	0.00023 J	0.001 U	0.001 U					
cis-1,2-Dichloroethene	0.07	0.07	0.0013	0.0031	0.0028	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.00022 J	0.00048 J
Ethylbenzene	0.7	0.7	0.001 U								
Methylene chloride	0.005	0.005	0.001 U								
Tetrachloroethene	0.005	0.005	0.001 U	0.00047 J	0.001 U						
Toluene	1	1	0.001 U								
trans-1,2-Dichloroethene	0.1	0.1	0.0005 U								
Trichloroethene	0.005	0.005	0.001 U	0.00064 J	0.001 U						
Vinyl chloride	0.002	0.002	0.001 U	0.001 U	0.00036 J	0.001 U					
Xylene (total)	10	10	0.001 U								
<i>Petroleum Products</i>											
Total Petroleum Hydrocarbons - purgeable (GRO)	0.500	0.500	0.1 U								

<sup>1</sup> Risk-Based Standards for Kansas RSK Manual - 4th Version, Appendix A, June 2007<sup>2</sup> mg/L - milligrams per liter<sup>3</sup> U - not present at or above the associated value<sup>4</sup> J - estimated concentration<sup>5</sup> Box denotes exceedence of KDHE criteria<sup>ab</sup> The letter a and/or b represents the Risk-Based Standard exceeded, respectively

**ATTACHMENT A**

**DATA QUALITY ASSESSMENT AND VALIDATION MEMORANDUM**



**CONESTOGA-ROVERS  
& ASSOCIATES**

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## MEMORANDUM

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TO: Phil Harvey (Chicago)

REF. NO.: 017307-095007

FROM: Susan C. Scrocchi/jbh/7~~SCS~~

DATE: January 16, 2008

Via E-Mail and U.S. Mail

C.C.: Patricia Klick (Chicago), Dennis Hoyt

RE: Data Quality Assessment and Validation  
Quarterly Groundwater Sampling Event  
GM Fairfax  
December 2007

PREVIOUSLY TRANSMITTED  
BY E-MAIL

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The following details a quality assessment and validation of the analytical data resulting from the December 2007 collection of the quarterly groundwater samples from the General Motors (GM) Fairfax Site in Kansas City, Kansas. The sample summary detailing sample identification, sample location, quality control (QC) samples, and analytical parameters is presented in Table 1. Sample analysis was completed at TestAmerica Inc. (TestAmerica), in North Canton, Ohio, in accordance with the methodologies presented in Table 2.

The QC criteria used to assess the data were established by the methods and the guidance document, "USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review", United States Environmental Protection Agency (USEPA) 540/R-99/008, October 1999.

These guidelines are collectively referred to as "Guidelines" in this memorandum.

### Sample Quantitation

The laboratory reported detected concentrations of volatile organic compounds (VOC) and Total Petroleum Hydrocarbon-Gasoline Range Organic (TPH-GRO), below the laboratory's practical quantitation limit (PQL)/report limit (RL) but above the laboratory's method detection limit (MDL). The laboratory flagged these sample concentrations with a "J". These concentrations should be qualified as estimated (J) values unless qualified otherwise in this memorandum.

### Sample Preservation and Holding Times

Sample holding time periods and preservation requirements are summarized in the analytical methods. All sample extractions and/or analyses were performed within the specified holding times.

All samples were properly preserved and cooled to 4°C(±2°C) after collection.

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Method Blank Samples

Method blank samples are prepared from a purified sample matrix and are processed concurrently with investigative samples to assess the presence and the magnitude of sample contamination introduced during sample analysis. Method blank samples are analyzed at a minimum frequency of one per analytical batch and target analytes should be non-detect.

Method blanks were analyzed at the recommended frequency. Methylene chloride was present at 0.58 parts per billion (ppb) on December 11, 2007. All associated sample results were non-detect and would not have been impacted.

Surrogate Compounds - Organic Analyses

Individual sample performance for organic analyses was monitored by assessing the results of surrogate compound percent recoveries. Surrogate percent recoveries are reviewed against the laboratory developed control limits provided in the analytical report.

All surrogate recoveries met the method criteria, demonstrating acceptable analytical efficiency for these analyses.

Matrix Spike/Matrix Spike Duplicate (MS/MSD)Analyses

To assess the long term accuracy and precision of the analytical methods on various matrices, MS/MSD percent recoveries and relative percent differences (RPD) of the concentrations were determined. The organic MS/MSD percent recovery and RPD control limits are established by the laboratory.

All MS/MSD recoveries were acceptable indicating good analytical accuracy and precision.

Laboratory Control Sample/Laboratory control Sample Duplicate (LCS/LCSD)

The LCS analysis serves as a monitor of the overall performance in all steps of the sample analysis and are analyzed with each sample batch. The LCS percent recoveries were evaluated against method and laboratory established control limits.

The LCS/LCSD percent recoveries were all within the laboratory control limits indicating acceptable analytical accuracy and precision.

Target Compound Identification

To minimize erroneous compound identification during organic analyses, qualitative criteria including compound retention time and mass spectra (if applicable) were evaluated according to identification criteria established by the methods. The organic compounds reported adhered to the specified identification criteria.

Field Quality Assurance/Quality Control (QA/QC)

The field QA/QC consisted of three equipment blank sample, four field duplicate sample sets and two trip blanks.

**CRA MEMORANDUM**

To assess the cleanliness of sample containers and the presence of field contamination, the equipment blank samples identified in Table 1 were collected and analyzed.

Chloroform, acetone and TPH-GRO were present at low concentrations in the rinse blanks. All associated sample results were with similar concentrations were qualified as non-detect (see Table 3).

Overall precision for the sampling event and laboratory procedures was monitored using the results of the field duplicate sample sets. The RPDs associated with these duplicate samples must be less than 50 percent for water. If the reported concentration in either the investigative sample or its duplicate is less than five times the RL, the evaluation criteria is one times the RL value for water.

All field duplicate results were acceptable indicating good field and analytical precision.

To monitor potential cross-contamination of VOC during aqueous sample transportation and storage, a trip blank was submitted to the laboratory for VOC analysis with each shipping cooler containing multiple samples.

All trip blank results were non-detect for the compounds of interest with the exception of a low concentration of acetone. All associated sample results were non-detect and would not have been impacted.

**Overall Assessment**

The data were found to exhibit acceptable levels of accuracy and precision, based on the provided information, and may be used with the qualifications noted.

**TABLE 1**  
**SAMPLE COLLECTION AND ANALYSIS SUMMARY**  
**QUARTERLY GROUNDWATER SAMPLING EVENT**  
**GENERAL MOTORS FAIRFAX**  
**KANSAS CITY, KANSAS**  
**DECEMBER 2007**

**Analysis/Parameters**

<b>Sample I.D.</b>	<b>Location I.D.</b>	<b>Collection Date (mm/dd/yy)</b>	<b>Collection Time (hr:min)</b>	<b>TCL VOCs</b>	<b>TPH-GRO</b>	<b>Comments</b>
GW-120407-JH-001	MW-116A	12/04/07	9:00	X	X	
GW-120407-JH-002	MW-116B	12/04/07	9:55	X	X	
GW-120407-JH-003	MW-117A	12/04/07	11:00	X	X	
GW-120407-JH-004	MW-117B	12/04/07	11:30	X	X	
GW-120407-JH-005	MW-1	12/04/07	13:20	X	X	
GW-120407-JH-006	MW-118A	12/04/07	14:40	X	X	
GW-120407-JH-007	MW-118B	12/04/07	14:55	X	X	
GW-120407-JH-008	MW-121A	12/04/07	15:45	X	X	
GW-120407-JH-009	MW-121B	12/04/07	16:00	X	X	
GW-120407-JH-010	MW-160A	12/04/07	16:50	X	X	
GW-120507-JH-011	MW-124A	12/05/07	8:55	X	X	
GW-120507-JH-012	MW-124B	12/05/07	9:15	X	X	
RW-120507-JH-013	Equipment Blank	12/05/07	9:40	X	X	
GW-120507-JH-014	MW-102A	12/05/07	10:50	X	X	
GW-120507-JH-015	MW-102B	12/05/07	11:10	X	X	
Trip Blank-A	Trip Blank	12/05/07	-	X		
GW-120507-JH-016	MW-102C	12/05/07	11:30	X	X	
GW-120507-JH-017	MW-108	12/05/07	13:05	X	X	
GW-120507-JH-018	MW-105	12/05/07	13:50	X	X	
GW-120507-JH-019	MW-104	12/05/07	14:20	X	X	
GW-120507-JH-020	MW-103C	12/05/07	14:50	X	X	
GW-120507-JH-021	MW-103B	12/05/07	15:05	X	X	
RW-120507-JH-022	Equipment Blank	12/05/07	15:20	X	X	
GW-120507-JH-023	MW-110A	12/05/07	16:05	X	X	
GW-120507-JH-024	MW-1C	12/05/07	16:45	X	X	
GW-120607-JH-025	MW-119B	12/06/07	9:15	X	X	
GW-120607-JH-026	MW-119A	12/06/07	9:30	X	X	
GW-120607-JH-027	MW-1B	12/06/07	10:30	X	X	
GW-120607-JH-028	MW-1B	12/06/07	10:45	X	X	
GW-120607-JH-029	MW-1A	12/06/07	11:15	X	X	
GW-120607-JH-030	MW-103A	12/06/07	12:00	X	X	
GW-120607-JH-031	MW-103A	12/06/07	12:30	X	X	Field duplicate of GW-120607-JH-027
GW-120607-JH-032	MW-125A	12/06/07	13:40	X	X	Field duplicate of GW-120607-JH-030
GW-120607-JH-033	MW-109	12/06/07	14:20	X	X	
RW-120607-JH-034	Equipment Blank	12/06/07	14:30	X	X	
GW-120607-JH-035	MW-111	12/06/07	15:10	X	X	Field duplicate of GW-120607-JH-035
GW-120607-JH-036	MW-111	12/06/07	15:30	X	X	
GW-120607-JH-037	MW-110	12/06/07	15:55	X	X	Field duplicate of GW-120607-JH-037
GW-120607-JH-038	MW-110	12/06/07	16:30	X	X	Field duplicate of GW-120607-JH-038
Trip Blank B Cooler 68187	Trip Blank	12/06/07	-	X		
Trip Blank C Cooler 68182	Trip Blank	12/06/07	-	X		

Notes:

- Not applicable.

GRO Gasoline Range Organic.

TCL Target Compound List.

TPH Total Petroleum Hydrocarbon.

VOC Volatile Organic Compound.

**TABLE 2**  
**SUMMARY OF ANALYTICAL METHODS**  
**QUARTERLY GROUNDWATER SAMPLING EVENT**  
**GENERAL MOTORS FAIRFAX**  
**KANSAS CITY, KANSAS**  
**DECEMBER 2007**

<i>Parameter</i>	<i>Method</i> <sup>1</sup>
TCL VOCs	SW-846 8260B
TPH-GRO	SW-846 8015

**Notes:**

<sup>1</sup> "Test Methods for Solid Waste/Physical Chemical Methods", SW-846, 3rd Edition, September 1986 (with all subsequent revisions).

GRO      Gasoline Range Organic.  
TCL      Target Compound List.  
TPH      Total Petroleum Hydrocarbon.  
VOCs     Volatile Organic Compounds.

**TABLE 3**  
**QUALIFIED SAMPLE RESULTS DUE TO ANALYTE CONCENTRATIONS IN THE RINSE BLANKS**  
**QUARTERLY GROUNDWATER SAMPLING EVENT**  
**GENERAL MOTORS FAIRFAX**  
**KANSAS CITY, KANSAS**  
**DECEMBER 2007**

Parameter	Rinse Blank Date	Analyte	Blank Result	Sample ID	Sample Result	Qualified Sample Result	Units
TPH-GRO	12/4-6/07	TPH - purgeable (GRO)	45	GW-120407-JH-001	60 J	100 U	ug/L
				GW-120407-JH-002	33 J	100 U	ug/L
				GW-120407-JH-003	45 J	100 U	ug/L
				GW-120407-JH-004	45 J	100 U	ug/L
				GW-120407-JH-005	30 J	100 U	ug/L
				GW-120407-JH-006	41 J	100 U	ug/L
				GW-120407-JH-007	40 J	100 U	ug/L
				GW-120407-JH-008	36 J	100 U	ug/L
				GW-120407-JH-009	33 J	100 U	ug/L
				GW-120407-JH-010	33 J	100 U	ug/L
				GW-120507-JH-011	36 J	100 U	ug/L
				GW-120507-JH-012	35 J	100 U	ug/L
				GW-120507-JH-014	32 J	100 U	ug/L
				GW-120507-JH-015	30 J	100 U	ug/L
				GW-120507-JH-016	33 J	100 U	ug/L
				GW-120507-JH-017	28 J	100 U	ug/L
				GW-120507-JH-020	31 J	100 U	ug/L
				GW-120507-JH-021	31 J	100 U	ug/L
				GW-120507-JH-023	33 J	100 U	ug/L
				GW-120607-JH-025	31 J	100 U	ug/L
				GW-120607-JH-026	33 J	100 U	ug/L
				GW-120607-JH-028	30 J	100 U	ug/L
				GW-120607-JH-032	50 J	100 U	ug/L
				GW-120607-JH-033	56 J	100 U	ug/L
				GW-120607-JH-035	130	130 U	ug/L
				GW-120607-JH-036	120	120 U	ug/L
				GW-120607-JH-037	29 J	100 U	ug/L

Notes:

GRO Gasoline Range Organic.  
 J Estimated.  
 TPH Total Petroleum Hydrocarbon.  
 U Not detected.

**ATTACHMENT B**

**HISTORICAL SUMMARY OF GROUNDWATER ANALYTICAL DATA**

TABLE B.1

**HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS**

Parameters	Units	Residential		Non-Residential		GW Pathway Scenario / GW Pathway Scenario <sup>1</sup>	MW-1	MW-1	MW-1	MW-1	MW-1	
		a	b	GW-100398-JM-001 10/3/1998	GW-122104-DS-006 12/21/2004	GW-091107-NR-006 9/11/2007	GW-120407-H-005 12/4/2007	GW-100398-JM-001 10/3/1998	GW-120407-H-005 12/4/2007	GW-100398-JM-001 10/3/1998		
<i>Volatile Organic Compounds</i>												
1,2-Dichloroethene (total)	mg/L	—	—	—	—	—	—	—	—	—	—	—
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.005 U <sup>3</sup>	0.005 U	0.01 U	0.01 U	0.005 U	0.01 U	0.005 U	0.005 U	0.005 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.005 U	0.005 U	0.01 U	0.01 U	0.008	0.01 U	0.008	0.008	0.008
Acetone	mg/L	0.26	0.93	0.005 U	0.005 U	0.01 U	0.01 U	0.005 U	0.01 U	0.005 U	0.005 U	0.005 U
Benzene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.001 U	0.00036 <sup>4</sup>	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.002 U	0.002 U	0.001 U	0.001 U	0.002 U	0.001 U	0.002 U	0.002 U	0.002 U
Tetrachloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Toluene	mg/L	1	1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.001 U	0.001 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U
Trichloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Xylene (total)	mg/L	10	10	0.001 U	0.002 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U

TABLE B.1

**HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS**

Parameters	Units	Residential		Non-Residential		GW Pathway Scenario <sup>1</sup> GW Pathway Scenario <sup>1</sup>	MW-1A	MW-1B	MW-1A	MW-1B
		a	b	GW-12104-DS-013	GW-091307-NP-029	GW-120607-JH-029	GW-091307-NP-028	GW-120607-JH-027	GW-091307-NP-028	GW-120607-JH-027
<i>Volatile Organic Compounds</i>										
1,2-Dichloroethene (total)	mg/L <sup>2</sup>	—	—	—	—	—	—	—	—	—
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Acetone	mg/L	0.26	0.93	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Benzene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.00074 J	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.00087 J	0.0004 J	0.00036 J	0.0005 U	0.0005 U	0.0005 U	0.0005 U
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Tetrachloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Toluene	mg/L	1	1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.001 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U
Trichloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Xylene (total)	mg/L	10	10	0.002 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U

HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX 1 PLANT SITE  
KANSAS CITY, KANSAS

Sample Location:		MW-1B	MW-1C	MW-1C	MW-1C	MW-1C
Sample ID:		GW-120607-JH-028	GW-100398-JM-009	GW-091307-NT-027	GW-120507-JH-024	GW-051229-JH-102
Sample Date:	12/6/2007	10/3/1998	9/13/2007	12/5/2007	5/12/1999	
Sample Type:	Duplicate					
Parameters						
Volatile Organic Compounds						
1,2-Dichloroethene (total) <sup>2</sup>	mg/L	—	—	—	—	0.001 U
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.01 U	0.005 U	0.02 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.01 U	0.041	0.01 U
Acetone	mg/L	0.26	0.93	0.01 U	0.005 U	0.1 U
Benzene	mg/L	0.005	0.005	0.001 U	0.001	0.001 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.001	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U	0.001	0.005 U
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.001	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.001	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.001	0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.0005 U	0.0005 U	0.0005 U
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.001	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.002	0.005 U
Tetrachloroethene	mg/L	0.005	0.005	0.001 U	0.001	0.001 U
Toluene	mg/L	1	1	0.001 U	0.21	0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.0005 U	0.0005 U	0.0005 U
Trichloroethene	mg/L	0.005	0.005	0.001 U	0.001	0.001 U
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.001	0.001 U
Xylene (total)	mg/L	10	10	0.001 U	0.11	0.001 U

HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS

Parameters	Units	Residential		Non-Residential		GW Pathway Scenario 1 GW Pathway Scenario <sup>1</sup>	MW-102A	MW-102B	MW-102A	MW-102B
		a	b	GW-062706-JH-014	GW-091307-NF-020	GW-120507-HH-014	GW-091307-NF-021	GW-120507-HH-015	GW-102A	GW-102B
<i>Volatile Organic Compounds</i>										
1,2-Dichloroethene (total)	mg/L <sup>2</sup>	—	—	—	—	—	—	—	—	—
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Acetone	mg/L	0.26	0.93	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Benzene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.00071 J	0.00071 J	0.00071 J	0.00071 J	0.00071 J	0.00071 J
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Tetrachloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Toluene	mg/L	1	1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U
Trichloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Xylene (total)	mg/L	10	10	0.001 U	0.00071 J	0.00071 J	0.00071 J	0.00071 J	0.00071 J	0.00071 J

TABLE B.1

**HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS**

Sample Location:	MW-102C	MW-102C	MW-103A	MW-103A	MW-103A
Sample ID:	GW-001307-NH-022	GW-120507-JH-016	GW-10399-JH-003	GW-051399-JH-105	GW-051399-JH-106
Sample Date:	9/13/2007	12/5/2007	10/3/1998	5/13/1999	5/13/1999
Sample Type:	Duplicate				
Parameters	a	b			
<i>Volatile Organic Compounds</i>					
1,2-Dichloroethene (total)	--	--	--	--	0.003
2-Butanone (Methyl Ethyl Ketone)	0.82	2.8	0.01 U	0.01 U	0.8 J
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	0.07	0.23	0.01 U	0.01 U	29 <sup>ab</sup>
Acetone	0.26	0.93	0.01 U	0.01 U	120 <sup>ab</sup>
Benzene	0.005	0.005	0.001 U	0.001 U	0.14 <sup>ab</sup>
Bromodichloromethane	0.08	0.08	0.001 U	0.001 U	0.001 U
Carbon disulfide	0.009	0.03	0.001 U	0.001 U	0.005 U
Chlorobenzene	0.1	0.1	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)	0.08	0.08	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	0.02	0.04	0.001 U	0.001 U	0.001 U
cis-1,2-Dichloroethene	0.07	0.07	0.0005 U	0.0005 U	0.003
Ethylbenzene	0.7	0.7	0.001 U	0.001 U	0.57
Methylene chloride	0.005	0.005	0.001 U	0.002 U	0.005 U
Tetrachloroethene	0.005	0.005	0.001 U	0.001 U	0.001 U
Toluene	1	1	0.001 U	0.001 U	9 <sup>ab</sup>
trans-1,2-Dichloroethene	0.1	0.1	0.0005 U	0.0005 U	0.001 U
Trichloroethene	0.005	0.005	0.001 U	0.001 U	0.001 U
Vinyl chloride	0.002	0.002	0.001 U	0.001 U	0.001 U
Xylene (total)	10	10	0.001 U	0.001 U	3.8
					3.9

TABLE B.1

**HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS**

Parameters	Units	Residential		Non-Residential		GW Pathway Scenario <sup>1</sup>	MW-103A	MW-103A	MW-103A	MW-103A	MW-103B	
		a	b	9/12/2007	9/12/2007	Duplicate	12/6/2007	12/6/2007	9/12/2007	Duplicate	9/12/2007	
<i>Volatile Organic Compounds</i>												
1,2-Dichloroethene (total)	mg/L	—	—	—	—	—	—	—	—	—	—	—
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	50 U	50 U	50 U	50 U	50 U	50 U	50 U	50 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	30 J <sup>a</sup>	33 J <sup>a</sup>	33 J <sup>a</sup>	160 <sup>b</sup>	0.01 U				
Acetone	mg/L	0.26	0.93	50 U	50 U	50 U	50 U	50 U	50 U	50 U	50 U	0.01 U
Benzene	mg/L	0.005	0.005	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U	0.001 U
Bromodichloromethane	mg/L	0.08	0.08	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U	0.001 U
Carbon disulfide	mg/L	0.009	0.03	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U	0.00091 J
Chlorobenzene	mg/L	0.1	0.1	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U	0.001 U
cis-1,2-Dichloroethylene	mg/L	0.07	0.07	2.5 U	2.5 U	2.5 U	2.5 U	2.5 U	2.5 U	2.5 U	2.5 U	0.0034
Ethylbenzene	mg/L	0.7	0.7	7.4 <sup>b</sup>	7.4 <sup>b</sup>	7.4 <sup>b</sup>	7.8 <sup>b</sup>	7.8 <sup>b</sup>	7.8 <sup>b</sup>	7.8 <sup>b</sup>	7.8 <sup>b</sup>	0.001 U
Methylene chloride	mg/L	0.005	0.005	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U	0.001 U
Tetrachloroethylene	mg/L	0.005	0.005	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U	0.001 U
Toluene	mg/L	1	1	120 <sup>b</sup>	130 <sup>b</sup>	130 <sup>b</sup>	150 <sup>b</sup>	150 <sup>b</sup>	150 <sup>b</sup>	150 <sup>b</sup>	150 <sup>b</sup>	0.001 U
trans-1,2-Dichloroethylene	mg/L	0.1	0.1	2.5 U	2.5 U	2.5 U	2.5 U	2.5 U	2.5 U	2.5 U	2.5 U	0.0005 U
Trichloroethylene	mg/L	0.005	0.005	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U	0.001 U
Vinyl chloride	mg/L	0.002	0.002	5 U	5 U	5 U	5 U	5 U	5 U	5 U	5 U	0.001 U
Xylene (total)	mg/L	10	10	63 <sup>b</sup>	67 <sup>b</sup>	67 <sup>b</sup>	57 <sup>b</sup>	57 <sup>b</sup>	57 <sup>b</sup>	57 <sup>b</sup>	57 <sup>b</sup>	0.001 U

HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS

Sample Location:	MW-103B	MW-103C	MW-103C	MW-103C
Sample ID:	GW-120507-JH-021	GW-100399-JM-005	GW-100398-JM-007	GW-120507-JH-020
Sample Date:	12/5/2007	10/3/1998	10/3/1998	9/12/2007
Sample Type:	Duplicate			
Parameters	Residential GW Pathway Scenario <sup>1</sup>	Non-Residential GW Pathway Scenario <sup>1</sup>		
Units	a	b		
<i>Volatile Organic Compounds</i>				
1,1,2-Dichloroethane (total)	mg/L	—	—	—
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.01 U
4-Methyl-1,2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.01 U
Acetone	mg/L	0.26	0.93	0.01 U
Benzene	mg/L	0.005	0.005	0.004
Bromodichloromethane	mg/L	0.08	0.08	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U
Chlorobenzene	mg/L	0.1	0.1	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.0064
Ethylbenzene	ng/L	0.7	0.7	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.001 U
Tetrachloroethene	mg/L	0.005	0.005	0.001 U
Toluene	mg/L	1	1	0.00026 J
trans-1,2-Dichloroethene	ng/L	0.1	0.1	0.0005 U
Trichloroethene	mg/L	0.005	0.005	0.001 U
Vinyl chloride	mg/L	0.002	0.002	0.001 U
Xylene (total)	mg/L	10	10	0.001 U
			0.38	0.16

HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS

Parameters	Sample Location:	Sample ID:	Sample Date:	Sample Type:	Units	Residential GW Pathway Scenario <sup>1</sup>	Non-Residential GW Pathway Scenario <sup>1</sup>	MW-104	MW-104	MW-104	MW-104	MW-104	MW-104
<i>Volatile Organic Compounds</i>					mg/L <sup>2</sup>	a	b	—	—	—	—	—	—
1,2-Dichloroethene (total)					mg/L	—	—	0.001 U	0.02 U	0.01 U	0.01 U	0.01 U	0.01 U
2-Butanone (Methyl Ethyl Ketone)					mg/L	0.82	2.8	—	—	—	—	—	—
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)					mg/L	0.07	0.23	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Acetone					mg/L	0.26	0.93	0.1 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Benzene					mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Bromodichloromethane					mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Carbon disulfide					mg/L	0.009	0.03	0.005 U	0.005 U	0.001 U	0.001 U	0.001 U	0.001 U
Chlorobenzene					mg/L	0.1	0.1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)					mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)					mg/L	0.02	0.04	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
cis-1,2-Dichloroethene					mg/L	0.07	0.07	0.001 U	0.0021 U	0.0005 U	0.00025 J	0.00022 J	0.00022 J
Ethylbenzene					mg/L	0.7	0.7	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Methylene chloride					mg/L	0.005	0.005	0.005 U	0.005 U	0.001 U	0.001 U	0.001 U	0.001 U
Tetrachloroethene					mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Toluene					mg/L	1	1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
trans-1,2-Dichloroethene					mg/L	0.1	0.1	0.001 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U
Trichloroethene					mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Vinyl chloride					mg/L	0.002	0.002	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Xylene (total)					mg/L	10	10	0.003 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U

TABLE B.1

**HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAIR I PLANT SITE  
KANSAS CITY, KANSAS**

Parameters	Units	Residential		Non-Residential		GW Pathway Scenario <sup>1</sup> GW Pathway Scenario <sup>1</sup>	MW-105	MW-105	MW-105	MW-105
		a	b	GW-062706-HH-003	GW-062706-HH-009					
<i>Volatile Organic Compounds</i>										
1,2-Dichloroethene (total)	mg/L <sup>2</sup>	--	--	0.001 U	--	--	--	--	--	--
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.02 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Acetone	mg/L	0.26	0.93	0.1 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Benzene	ng/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.005 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.001 U	0.0063	0.0049	0.0031	0.001 U	0.001 U	0.001 U
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.005 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Tetrachloroethene	mg/L	0.005	0.005	0.001	0.001 U	0.001 U	0.001 U	0.001 U	0.00063 J	0.00063 J
Toluene	mg/L	1	1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.001 U	0.001 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U
Trichloroethene	mg/L	0.005	0.005	0.005 U	0.001 U	0.00058 J	0.00043 J	0.00043 J	0.00043 J	0.00043 J
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Xylene (total)	mg/L	10	10	0.003 U	0.002 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U

**TABLE B.1**  
**HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA**  
**FORMER FAIRFAX I PLANT SITE**  
**KANSAS CITY, KANSAS**

Parameters	Units	Residential		Non-Residential		GW Pathway Scenario <sup>1</sup>	GW Pathway Scenario <sup>1</sup>
		a	b	MW-108	MW-108		
<b>Volatile Organic Compounds</b>							
1,2-Dichloroethene (total)	mg/L <sup>2</sup>	--	--	--	--	--	--
2-Butanone (Methyl Ethyl Ketone)	ng/L	0.82	2.8	0.01 U	0.005 U	0.01 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	ng/L	0.07	0.23	0.01 U	0.005 U	0.01 U	0.01 U
Acetone	mg/L	0.26	0.93	0.01 U	0.005 U	0.01 U	0.01 U
Benzene	mg/L	0.005	0.005	0.001 U	0.004 J	0.001 U	0.001 U
Bromodichloromethane	ng/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U	0.001 U	0.001 U	0.001 U
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.001 U	0.001 U	0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.0037	0.001 U	0.0005 U	0.0005 U
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.001 U	0.001 U	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.002 U	0.001 U	0.001 U
Tetrachloroethene	mg/L	0.005	0.005	0.00035 J	0.001 U	0.001 U	0.001 U
Toluene	mg/L	1	1	0.001 U	0.001 U	0.001 U	0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.0005 U	0.001 U	0.0005 U	0.0005 U
Trichloroethene	mg/L	0.005	0.005	0.00075 J	0.001 U	0.001 U	0.001 U
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.001 U	0.001 U	0.001 U
Xylyne (total)	mg/L	10	10	0.001 U	0.001 U	0.001 U	0.001 U

TABLE B.1

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**HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS**

Sample Location:	MW-109	MW-109	MW-109	MW-109	MW-109
Sample ID:	GW-100398-JM-023	GW-12204-DS-005	GW-091307-NYL-025	GW-120607-JH-033	GW-100398-JM-025
Sample Date:	10/3/1998	12/21/2004	9/13/2007	12/6/2007	10/31/1998
<b>Sample Type:</b>					
<b>Residential GW Pathway Scenario<sup>1</sup></b>		<b>Non-Residential GW Pathway Scenario<sup>1</sup></b>			
Parameters	Units	a	b	c	d
<b>Volatile Organic Compounds</b>					
1,2-Dichloroethene (total)	mg/L <sup>2</sup>	—	—	—	—
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.32	2.8	0.005 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.005 U	0.01 U
Acetone	mg/L	0.26	0.93	0.005 U	0.0022 J
Benzene	mg/L	0.005	0.005	0.001 U	0.001 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U	0.001 U
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.018	0.02
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.002 U	0.001 U
Tetrachloroethene	mg/L	0.005	0.005	0.001 U	0.001 U
Toluene	mg/L	1	1	0.001 U	0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.002	0.0013
Trichloroethene	mg/L	0.005	0.005	0.001 U	0.001 U
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.0012
Xylene (total)	mg/L	10	10	0.001 U	0.00029 J

HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS

Parameters	Units	Residential		Non-Residential		GW Pathway Scenario <sup>1</sup> GW Pathway Scenario <sup>1</sup>	MW-110	MW-110	MW-110	MW-110	MW-110	
		a	b	GW-12214-DS-012	12/21/2004		GW-122104-DS DUP	1/18/2006	GW-091307-JH-008	9/13/2007	GW-091307-JH-023	12/6/2007
<i>Volatile Organic Compounds</i>												
1,2-Dichloroethene (total)	mg/L	—	—	—	—	—	—	—	—	—	—	—
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Acetone	mg/L	0.26	0.93	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Benzene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U	0.001 U	0.00049 J	0.00028 J	0.00028 J	0.00028 J	0.00028 J	0.00028 J	0.00028 J
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.00074 J	0.00067 J	0.00067 J	0.00066 J	0.00066 J	0.00066 J	0.00066 J	0.00066 J	0.00066 J
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Tetrachloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Toluene	mg/L	1	1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.001 U	0.001 U	0.001 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U
Trichloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Xylene (total)	mg/L	10	10	0.002 U	0.002 U	0.002 U	0.002 U	0.002 U	0.002 U	0.002 U	0.002 U	0.002 U

HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS

Parameters	Units	Residential		Non-Residential		GW Pathway Scenario <sup>1</sup> GW Pathway Scenario <sup>1</sup>	MW-T10A	MW-T10A	MW-T10A	MW-T10A	MW-T10A
		a	b	GW-12607-H-038	GW-100398-JM-017	GW-122104-DS-008	GW-091407-NT-033	GW-091407-NT-034			
Sample ID:				12/6/2007	10/3/1998	12/21/2004	9/14/2007	9/14/2007			
Sample Date:				Duplicate			Duplicate	Duplicate			
Sample Type:											
<b>Volatile Organic Compounds</b>											
1,2-Dichloroethene (total)	mg/L <sup>2</sup>	—	—	—	—	—	—	—	—	—	—
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.01 U	0.005 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.01 U	0.005 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Acetone	mg/L	0.26	0.93	0.01 U	0.005 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Benzene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.002	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
cis-1,2-Dichloroethylene	mg/L	0.07	0.07	0.00035 J	0.002	0.004	0.004	0.0026	0.0026	0.0027	0.0027
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.002 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Tetrachloroethylene	mg/L	0.005	0.005	0.001 U	0.029 <sup>3b</sup>	0.0048	0.0048	0.0073 <sup>3b</sup>	0.0073 <sup>3b</sup>	0.0073 <sup>3b</sup>	0.0073 <sup>3b</sup>
Toluene	mg/L	1	1	0.001 U	0.023	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.0005 U	0.00021 J	0.001 U	0.0005 U	0.00021 J	0.00021 J	0.0005 U	0.0005 U
Trichloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.0014	0.0014
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Xylene (total)	mg/L	10	10	0.001 U	0.013	0.002 U	0.002 U	0.002 U	0.002 U	0.001 U	0.001 U

TABLE B.1

**HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS**

Sample Location:	MW-110A	MW-111	MW-111	MW-111	MW-111
Sample ID:	GW-120307-JH-023	GW-122104-DS-011	GW-091307-NT-024	GW-120607-JH-025	GW-120607-JH-036
Sample Date:	12/5/2007	12/21/2004	9/13/2007	12/6/2007	12/6/2007
Sample Type:	Duplicate				
Parameters	Residential GW Pathway Scenario <sup>1</sup>	Non-Residential GW Pathway Scenario <sup>1</sup>			
Units	a	b			
<i>Volatile Organic Compounds</i>					
1,2-Dichloroethene (total)	mg/L	--	--	--	--
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.01 U	0.05 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.01 U	0.05 U
Acetone	mg/L	0.26	0.93	0.01 U	0.05 U
Benzene	mg/L	0.005	0.005	0.001 U	0.005 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.005 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U	0.005 U
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.005 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.005 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.005 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.0019	0.15 <sup>ab</sup>
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.005 U
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.005 U
Tetrachloroethene	mg/L	0.005	0.005	0.0096 <sup>ab</sup>	0.005 U
Toluene	mg/L	1	1	0.001 U	0.005 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.0005 U	0.011
Trichloroethene	mg/L	0.005	0.005	0.0014	0.005 U
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.0017 <sup>bc</sup>
Xylene (total)	mg/L	10	10	0.001 U	0.0025 U

TABLE B.1

**HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS**

Sample Location:	MW-115A	MW-116A	MW-116A	MW-116A	MW-116A
Sample ID:	GW-051399-JH-104	GW-091207-NF-019	GW-011706-JH-001	GW-062706-JH-008	GW-091107-NF-002
Sample Date:	5/13/1999	9/12/2007	1/17/2006	6/27/2006	9/11/2007
Sample Type:					
<i>Parameters</i>					
<i>Volatile Organic Compounds</i>					
1,2-Dichloroethene (total)	—	—	0.0001 U	—	—
2-Butanone (Methyl Ethyl Ketone)	0.82	2.8	0.02 U	0.01 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	0.07	0.23	0.01 U	0.00044 J	0.01 U
Acetone	0.26	0.93	0.1 U	0.01 U	0.01 U
Benzene	0.005	0.005	0.001 U	0.001 U	0.001 U
Bromodichloromethane	0.08	0.08	0.001 U	0.001 U	0.001 U
Carbon disulfide	0.009	0.03	0.005 U	0.001 U	0.001 U
Chlorobenzene	0.1	0.1	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)	0.08	0.08	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	0.02	0.04	0.001 U	0.001 U	0.001 U
cis-1,2-Dichloroethene	0.07	0.07	0.001 U	0.0005 U	0.0005 U
Ethylbenzene	0.7	0.7	0.001 U	0.0019	0.001 U
Methylene chloride	0.005	0.005	0.005 U	0.001 U	0.001 U
Tetrachloroethene	0.005	0.005	0.001 U	0.001 U	0.001 U
Toluene	1	1	0.001 U	0.016	0.001 U
trans-1,2-Dichloroethene	0.1	0.1	0.001 U	0.0005 U	0.0005 U
Trichloroethene	0.005	0.005	0.001 U	0.001 U	0.001 U
Vinyl chloride	0.002	0.002	0.001 U	0.001 U	0.001 U
Xylene (total)	10	10	0.003 U	0.016	0.001 U

TABLE B.1

**HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA**  
**FORMER FAIRFAX I PLANT SITE**  
**KANSAS CITY, KANSAS**

Sample Location:	MW-116A	MW-116B	MW-116B	MW-116B	MW-116B	MW-116B
Sample ID:	GW-120407-JH-001	GW-011706-JH-002	GW-0622706-JH-007	GW-091107-NR-001	GW-120407-JH-002	MW-116B
Sample Date:	12/4/2007	1/17/2006	6/27/2006	9/11/2007	12/4/2007	GW-120407-JH-002
Sample Type:						
<i>Parameters</i>						
<i>Volatile Organic Compounds</i>						
1,2-Dichloroethene (total)	mg/L	—	—	—	—	—
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.01 U	0.01 U	0.01 U
4-Methyl-1,2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.01 U	0.01 U	0.01 U
Acetone	mg/L	0.26	0.93	0.01 U	0.01 U	0.01 U
Benzene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U	0.001 U	0.001 U
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.001 U	0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.0005 U	0.0005 U	0.0005 U
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.001 U	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U
Tetrachloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U
Toluene	mg/L	1	1	0.00022 U	0.001 U	0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.0005 U	0.0005 U	0.0005 U
Trichloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.001 U	0.001 U
Xylene (total)	mg/L	10	10	0.001 U	0.001 U	0.001 U

TABLE B.1

**HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS**

Parameters	Units	Residential GW Pathway Scenario <sup>1</sup>		Non-Residential GW Pathway Scenario <sup>1</sup>		MW-117A GW-011706-JH-003 1/17/2006	MW-117A GW-062706-JH-004 6/27/2006	MW-117A GW-091107-NT-003 9/11/2007	MW-117A GW-091107-NT-004 9/11/2007	MW-117A GW-120407-JH-003 12/4/2007	
		a	b	a	b						
<b>Volatile Organic Compounds</b>											
1,2-Dichloroethene (total)	mg/L <sup>2</sup>	--	--	--	--	--	--	--	--	--	--
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Acetone	mg/L	0.26	0.93	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Benzene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Tetrachloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Toluene	mg/L	1	1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U
Trichloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Xylene (total)	mg/L	10	10	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U

TABLE B.1

**HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA**  
**FORMER FAIRFAX I PLANT SITE**  
**KANSAS CITY, KANSAS**

Sample Location:	MW-117B	MW-117B	MW-117B	MW-117B	MW-118A
Sample ID:	GW-011706-JH-004	GW-062706-JH-003	GW-091107-NT-005	GW-120407-JH-004	GW-011706-JH-006
Sample Date:	1/17/2006	6/27/2006	9/11/2007	12/4/2007	1/17/2006
Sample Type:					
<i>Parameters</i>					
<i>Volatile Organic Compounds</i>					
1,2-Dichloroethene (total)	—	—	—	—	—
2-Butanone (Methyl Ethyl Ketone)	0.82	2.8	0.01 U	0.01 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	0.07	0.23	0.01 U	0.01 U	0.01 U
Acetone	0.26	0.93	0.01 U	0.01 U	0.01 U
Benzene	0.005	0.005	0.001 U	0.001 U	0.001 U
Bromodichloromethane	0.08	0.08	0.001 U	0.001 U	0.001 U
Carbon disulfide	0.009	0.03	0.001 U	0.001 U	0.001 U
Chlorobenzene	0.1	0.1	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)	0.08	0.08	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	0.02	0.04	0.00019 J	0.001 U	0.001 U
cis-1,2-Dichloroethene	0.07	0.07	0.0013	0.0011	0.0013
Ethylbenzene	0.7	0.7	0.001 U	0.001 U	0.001 U
Methylene chloride	0.005	0.005	0.001 U	0.001 U	0.001 U
Tetrachloroethene	0.005	0.005	0.001 U	0.001 U	0.001 U
Toluene	1	1	0.001 U	0.001 U	0.001 U
trans-1,2-Dichloroethene	0.1	0.1	0.0005 U	0.0005 U	0.0005 U
Trichloroethene	0.005	0.005	0.001 U	0.001 U	0.0016
Vinyl chloride	0.002	0.002	0.001 U	0.001 U	0.001 U
Xylene (total)	10	10	0.001 U	0.001 U	0.001 U

TABLE B.1

**HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS**

Sample Location:	MW-118A	MW-118B	MW-118A	MW-118B	MW-118B
Sample ID:	GW-062706-JH-006	GW-091107-NT-008	GW-120407-JH-006	GW-011706-JH-005	GW-062706-JH-005
Sample Date:	6/27/2006	9/11/2007	12/4/2007	1/17/2006	6/27/2006
Sample Type:					
Residential					
Non-Residential					
GW Pathway Scenario <sup>1</sup> GW Pathway Scenario <sup>1</sup>					
Parameters	Units	a	b	c	d
<i>Volatile Organic Compounds</i>					
1,2-Dichloroethene (total)	mg/L	—	—	—	—
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.01 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.01 U	0.01 U
Acetone	mg/L	0.26	0.93	0.01 U	0.01 U
Benzene	mg/L	0.005	0.005	0.001 U	0.001 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U	0.001 U
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.0035	0.0045
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.001 U
Tetrachloroethene	mg/L	0.005	0.005	0.001 U	0.001 U
Toluene	mg/L	1	1	0.001 U	0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.0005 U	0.00027 J
Trichloroethene	mg/L	0.005	0.005	0.00063 J	0.001 U
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.00038 J
Xylene (total)	mg/L	10	10	0.001 U	0.001 U

HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS

Parameters	Units	Residential		Non-Residential		GW Pathway Scenario <sup>1</sup>	GW Pathway Scenario <sup>1</sup>	MW-118B 9/11/2007	MW-118B 12/4/2007	MW-118A 1/18/2006	MW-119A 1/18/2006	MW-119A 1/18/2006	MW-119A 1/18/2006	MW-119A 6/27/2006	
		a	b	GW-09107-NR-007 9/11/2007	GW-011806-JH-007 12/4/2007	GW-011806-JH-010 1/18/2006	GW-011806-JH-011 1/18/2006								
<i>Volatile Organic Compounds</i>															
1,2-Dichloroethene (total)	mg/L	--	--	--	--	--	--	--	--	--	--	--	--	--	--
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Acetone	mg/L	0.26	0.93	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Benzene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U	0.001 U	0.001 U	0.0003 J	0.0003 J	0.0003 J	0.0003 J	0.0003 J	0.0003 J	0.0003 J	0.0003 J	0.0003 J
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.0016	0.0016	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.0029	0.0029	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Tetrachloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Toluene	mg/L	1	1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U
Trichloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Vinyl chloride	mg/L	0.002	0.002	0.00031 J	0.00031 J	0.00036 J	0.00036 J	0.00036 J	0.00036 J	0.00036 J	0.00036 J	0.00036 J	0.00036 J	0.00036 J	0.00036 J
Xylene (total)	mg/L	10	10	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U

TABLE B.1

**HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS**

Parameters	Units	a	b	Residential GW Pathway Scenario <sup>1</sup>	Non-Residential GW Pathway Scenario <sup>1</sup>	MW-119A GW-091307-NI-030 9/13/2007	MW-119B GW-011806-JH-009 12/6/2007	MW-119B GW-011806-JH-009 1/18/2006	MW-119B GW-062706-JH-016 6/27/2006	MW-119B GW-091307-NI-031 9/13/2007	
<i>Volatile Organic Compounds</i>											
1,2-Dichloroethene (total)	mg/L <sup>2</sup>	—	—	—	—	—	—	—	—	—	—
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Acetone	mg/L	0.26	0.93	0.023 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U	0.01 U
Benzene	ng/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Tetrachloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Toluene	mg/L	1	1	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U	0.0005 U
Trichloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U
Xylene (total)	mg/L	10	10	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U	0.001 U

TABLE B.1

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**HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA**  
**FORMER FAIRFAX 1 PLANT SITE**  
**KANSAS CITY, KANSAS**

Parameters	Units	Residential <sup>1</sup>		Non-Residential <sup>1</sup>		GW Pathway Scenario <sup>1</sup> GW Pathway Scenario <sup>1</sup>
		a	b	GW-01806-JH-013	MW-121A	
<i>Volatile Organic Compounds</i>						
1,2-Dichloroethene (total)	mg/L	—	—	—	—	—
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.01 U	0.01 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.01 U	0.01 U	0.01 U
Acetone	mg/L	0.26	0.93	0.01 U	0.01 U	0.01 U
Benzene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U	0.001 U	0.001 U
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.001 U	0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.0005 U	0.0005 U	0.0005 U
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.001 U	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U
Tetrachloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U
Toluene	mg/L	1	1	0.001 U	0.001 U	0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.0005 U	0.0005 U	0.0005 U
Trichloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.001 U	0.001 U
Xylene (total)	mg/L	10	10	0.001 U	0.001 U	0.001 U

HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS

Sample Location:	MW-121B	MW-121B	MW-121B	MW-121B	MW-121B
Sample ID:	GW-011806-JH-012	GW-062606-JH-001	GW-091407-NT-036	GW-120407-JH-009	MW-124A
Sample Date:	1/18/2006	6/26/2006	9/14/2007	12/4/2007	11/28/2006
Sample Type:					
Parameters	Residential GW Pathway Scenario <sup>1</sup>	Non-Residential GW Pathway Scenario <sup>1</sup>	Residential GW Pathway Scenario <sup>1</sup>	Non-Residential GW Pathway Scenario <sup>1</sup>	Residential GW Pathway Scenario <sup>1</sup>
Units	a	b	a	b	a
<i>Volatile Organic Compounds</i>					
1,2-Dichloroethene (total)	mg/L	—	—	—	—
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.01 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.01 U	0.01 U
Acetone	mg/L	0.26	0.93	0.01 U	0.01 U
Benzene	mg/L	0.005	0.005	0.001 U	0.001 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U	0.001 U
Chlorobenzene	mg/L	0.1	0.1	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.0005 U	0.0005 U
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.001 U
Tetrachloroethene	mg/L	0.005	0.005	0.001 U	0.001 U
Toluene	mg/L	1	1	0.001 U	0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.0005 U	0.0005 U
Trichloroethene	mg/L	0.005	0.005	0.001 U	0.001 U
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.001 U
Xylene (total)	mg/L	10	10	0.001 U	0.001 U

HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS

Parameters	Units	Residential		Non-Residential		GW Pathway Scenario <sup>1</sup>	GW Pathway Scenario <sup>1</sup>
		a	b	MW-124A	MW-124B		
<b>Volatile Organic Compounds</b>	mg/L <sup>2</sup>	—	—	—	—	—	—
1,2-Dichloroethene (total)	mg/L	0.82	2.8	0.01 U	0.01 U	0.0029 J	0.01 U
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.07	0.23	0.01 U	0.01 U	0.01 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.26	0.93	0.01 U	0.01 U	0.01 U	0.01 U
Acetone	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U
Benzene	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U
Bromodichloromethane	mg/L	0.009	0.03	0.001 U	0.001 U	0.001 U	0.001 U
Carbon disulfide	mg/L	0.1	0.1	0.001 U	0.001 U	0.001 U	0.0003 J
Chlorobenzene	mg/L	0.08	0.08	0.001 U	0.001 U	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.02	0.04	0.001 U	0.001 U	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.07	0.07	0.0005 U	0.0005 U	0.0005 U	0.0005 U
cis-1,2-Dichloroethene	mg/L	0.7	0.7	0.001 U	0.001 U	0.001 U	0.001 U
Ethylbenzene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U
Tetrachloroethene	mg/L	1	1	0.001 U	0.001 U	0.001 U	0.001 U
Toluene	ng/L	0.1	0.1	0.0005 U	0.0005 U	0.0005 U	0.0005 U
trans-1,2-Dichloroethene	ng/L	0.005	0.005	0.001 U	0.001 U	0.001 U	0.001 U
Trichloroethene	ng/L	0.002	0.002	0.001 U	0.001 U	0.001 U	0.001 U
Vinyl chloride	ng/L	10	10	0.001 U	0.001 U	0.001 U	0.001 U
Xylene (total)	ng/L	—	—	—	—	—	—

**HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS**

Parameters	Units	Residential		Non-Residential		GW Pathway Scenario 1 GW Pathway Scenario <sup>1</sup>
		a	b	MW-124B 12/5/2007	MW-125 11/1/2007	
<i>Volatile Organic Compounds</i>						
1,2-Dichloroethene (total)	mg/L <sup>2</sup>	—	—	—	—	—
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.82	2.8	0.01 U	0.01 U	0.005 U 0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.07	0.23	0.01 U	0.01 U	0.008 U 0.01 U
Acetone	mg/L	0.26	0.93	0.01 U	0.01 U	0.005 U 0.01 U
Benzene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U 0.001 U
Bromodichloromethane	mg/L	0.08	0.08	0.001 U	0.00017 J	0.001 U 0.001 U
Carbon disulfide	mg/L	0.009	0.03	0.001 U	0.001 U	0.001 U 0.001 U
Chlorobenzene	mg/L	0.1	0.1	0.00023 J	0.0001 U	0.001 U 0.001 U
Chloroform (Trichloromethane)	mg/L	0.08	0.08	0.001 U	0.00033 J	0.001 U 0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.02	0.04	0.001 U	0.001 U	0.001 U 0.001 U
cis-1,2-Dichloroethene	mg/L	0.07	0.07	0.00022 J	0.0005 U	0.00048 J 0.001 U
Ethylbenzene	mg/L	0.7	0.7	0.001 U	0.001 U	0.002 U 0.001 U
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.001 U	0.002 U 0.001 U
Tetrachloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.00047 J 0.001 U
Toluene	mg/L	1	1	0.001 U	0.001 U	0.041 U 0.001 U
trans-1,2-Dichloroethene	mg/L	0.1	0.1	0.0005 U	0.0005 U	0.001 U 0.001 U
Trichloroethene	mg/L	0.005	0.005	0.001 U	0.001 U	0.001 U 0.001 U
Vinyl chloride	mg/L	0.002	0.002	0.001 U	0.001 U	0.001 U 0.001 U
Xylene (total)	mg/L	10	10	0.001 U	0.001 U	0.018 U 0.002 U

HISTORICAL SUMMARY OF DETECTED GROUNDWATER ANALYTICAL DATA  
FORMER FAIRFAX I PLANT SITE  
KANSAS CITY, KANSAS

Parameters	Units	Residential		Non-Residential	
		GW Pathway Scenario <sup>1</sup>	GW Pathway Scenario <sup>1</sup>	GW-160A	GW-160A
Volatile Organic Compounds	mg/L <sup>2</sup>	—	—	—	—
1,2-Dichloroethene (total)	mg/L	0.82	2.8	0.01 U	0.01 U
2-Butanone (Methyl Ethyl Ketone)	mg/L	0.07	0.23	0.01 U	0.01 U
4-Methyl-2-Pentanone (Methyl Isobutyl Ketone)	mg/L	0.26	0.93	0.01 U	0.01 U
Acetone	mg/L	0.005	0.005	0.001 U	0.001 U
Benzene	mg/L	0.08	0.08	0.001 U	0.001 U
Bromodichloromethane	mg/L	0.009	0.03	0.001 U	0.001 U
Carbon disulfide	mg/L	0.1	0.1	0.001 U	0.001 U
Chlorobenzene	mg/L	0.08	0.08	0.001 U	0.001 U
Chloroform (Trichloromethane)	mg/L	0.02	0.04	0.001 U	0.001 U
Chloromethane (Methyl Chloride)	mg/L	0.07	0.07	0.0005 U	0.0005 U
cis-1,2-Dichloroethylene	mg/L	0.7	0.7	0.001 U	0.001 U
Ethylbenzene	mg/L	0.005	0.005	0.001 U	0.001 U
Methylene chloride	mg/L	0.005	0.005	0.001 U	0.001 U
Tetrachloroethene	mg/L	1	1	0.001 U	0.001 U
Toluene	mg/L	0.1	0.1	0.0005 U	0.0005 U
trans-1,2-Dichloroethene	mg/L	0.005	0.005	0.001 U	0.001 U
Trichloroethene	mg/L	0.002	0.002	0.001 U	0.001 U
Vinyl chloride	mg/L	10	10	0.001 U	0.001 U
Xylene (total)	mg/L				

<sup>1</sup>Risk-Based Standards for Kansas RSK Manual - 4th Version, Appendix A, June 2007

<sup>2</sup>mg/L - milligrams per liter

<sup>3</sup>U - not present at or above the associated value

<sup>4</sup>J - estimated concentration

<sup>5</sup>Box denotes exceedence of KDHE criteria

<sup>a,b</sup>The letter a and/or b represents the Risk-Based Standard exceeded, respectively

**ATTACHMENT C**

**LABORATORY ANALYTICAL DATA**

TABLE C.1

**SAMPLE KEY**  
**FORMER FAIRFAX I PLANT SITE**  
**KANSAS CITY, KANSAS**

Sample ID	Well ID	Date	QA/QC <sup>1</sup>	Parameters <sup>2</sup>
GW-120407-JH-001	MW-116A	12/04/07		TCL VOCs, TPH-GRO
GW-120407-JH-002	MW-116B	12/04/07		TCL VOCs, TPH-GRO
GW-120407-JH-003	MW-117A	12/04/07		TCL VOCs, TPH-GRO
GW-120407-JH-004	MW-117B	12/04/07		TCL VOCs, TPH-GRO
GW-120407-JH-005	MW-1	12/04/07		TCL VOCs, TPH-GRO
GW-120407-JH-006	MW-118A	12/04/07		TCL VOCs, TPH-GRO
GW-120407-JH-007	MW-118B	12/04/07		TCL VOCs, TPH-GRO
GW-120407-JH-008	MW-121A	12/04/07		TCL VOCs, TPH-GRO
GW-120407-JH-009	MW-121B	12/04/07	MS/MSD	TCL VOCs, TPH-GRO
GW-120407-JH-010	MW-160A	12/04/07		TCL VOCs, TPH-GRO
GW-120507-JH-011	MW-124A	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-012	MW-124B	12/05/07		TCL VOCs, TPH-GRO
RW-120507-JH-013	Equipment Blank	12/05/07	Equipment Blank	TCL VOCs, TPH-GRO
GW-120507-JH-014	MW-102A	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-015	MW-102B	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-016	MW-102C	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-017	MW-108	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-018	MW-105	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-019	MW-104	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-020	MW-103C	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-021	MW-103B	12/05/07		TCL VOCs, TPH-GRO
RW-120507-JH-022	Equipment Blank	12/05/07	Equipment Blank	TCL VOCs, TPH-GRO
GW-120507-JH-023	MW-110A	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-024	MW-1C	12/05/07		TCL VOCs, TPH-GRO
GW-120607-JH-025	MW-119B	12/06/07		TCL VOCs, TPH-GRO
GW-120607-JH-026	MW-119A	12/06/07	MS/MSD	TCL VOCs, TPH-GRO
GW-120607-JH-027	MW-1B	12/06/07		TCL VOCs, TPH-GRO
GW-120607-JH-028	MW-1B	12/06/07	Duplicate	TCL VOCs, TPH-GRO
GW-120607-JH-029	MW-1A	12/06/07		TCL VOCs, TPH-GRO
GW-120607-JH-030	MW-103A	12/06/07		TCL VOCs, TPH-GRO
GW-120607-JH-031	MW-103A	12/06/07	Duplicate	TCL VOCs, TPH-GRO
GW-120607-JH-032	MW-125A	12/06/07		TCL VOCs, TPH-GRO
GW-120607-JH-033	MW-109	12/06/07		TCL VOCs, TPH-GRO
RW-120607-JH-034	Equipment Blank	12/06/07	Equipment Blank	TCL VOCs, TPH-GRO
GW-120607-JH-035	MW-111	12/06/07		TCL VOCs, TPH-GRO
GW-120607-JH-036	MW-111	12/06/07	Duplicate	TCL VOCs, TPH-GRO
GW-120607-JH-037	MW-110	12/06/07		TCL VOCs, TPH-GRO
GW-120607-JH-038	MW-110	12/06/07	Duplicate	TCL VOCs, TPH-GRO

<sup>1</sup> QA/QC - quality assurance/quality control

MS/MSD - matrix spike/matrix spike duplicate

<sup>2</sup> TCL VOCs - target compound list volatile organic compounds

TPH-GRO - total petroleum hydrocarbon gasoline range organics

**ATTACHMENT C**

**LABORATORY ANALYTICAL DATA**

**TABLE C.1**  
**SAMPLE KEY**  
**FORMER FAIRFAX I PLANT SITE**  
**KANSAS CITY, KANSAS**

<i>Sample ID</i>	<i>Well ID</i>	<i>Date</i>	<i>QA/QC</i> <sup>1</sup>	<i>Parameters</i> <sup>2</sup>
GW-120407-JH-001	MW-116A	12/04/07		TCL VOCs, TPH-GRO
GW-120407-JH-002	MW-116B	12/04/07		TCL VOCs, TPH-GRO
GW-120407-JH-003	MW-117A	12/04/07		TCL VOCs, TPH-GRO
GW-120407-JH-004	MW-117B	12/04/07		TCL VOCs, TPH-GRO
GW-120407-JH-005	MW-1	12/04/07		TCL VOCs, TPH-GRO
GW-120407-JH-006	MW-118A	12/04/07		TCL VOCs, TPH-GRO
GW-120407-JH-007	MW-118B	12/04/07		TCL VOCs, TPH-GRO
GW-120407-JH-008	MW-121A	12/04/07		TCL VOCs, TPH-GRO
GW-120407-JH-009	MW-121B	12/04/07	MS/MSD	TCL VOCs, TPH-GRO
GW-120407-JH-010	MW-160A	12/04/07		TCL VOCs, TPH-GRO
GW-120507-JH-011	MW-124A	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-012	MW-124B	12/05/07		TCL VOCs, TPH-GRO
RW-120507-JH-013	Equipment Blank	12/05/07	Equipment Blank	TCL VOCs, TPH-GRO
GW-120507-JH-014	MW-102A	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-015	MW-102B	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-016	MW-102C	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-017	MW-108	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-018	MW-105	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-019	MW-104	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-020	MW-103C	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-021	MW-103B	12/05/07		TCL VOCs, TPH-GRO
RW-120507-JH-022	Equipment Blank	12/05/07	Equipment Blank	TCL VOCs, TPH-GRO
GW-120507-JH-023	MW-110A	12/05/07		TCL VOCs, TPH-GRO
GW-120507-JH-024	MW-1C	12/05/07		TCL VOCs, TPH-GRO
GW-120607-JH-025	MW-119B	12/06/07		TCL VOCs, TPH-GRO
GW-120607-JH-026	MW-119A	12/06/07	MS/MSD	TCL VOCs, TPH-GRO
GW-120607-JH-027	MW-1B	12/06/07		TCL VOCs, TPH-GRO
GW-120607-JH-028	MW-1B	12/06/07	Duplicate	TCL VOCs, TPH-GRO
GW-120607-JH-029	MW-1A	12/06/07		TCL VOCs, TPH-GRO
GW-120607-JH-030	MW-103A	12/06/07		TCL VOCs, TPH-GRO
GW-120607-JH-031	MW-103A	12/06/07	Duplicate	TCL VOCs, TPH-GRO
GW-120607-JH-032	MW-125A	12/06/07		TCL VOCs, TPH-GRO
GW-120607-JH-033	MW-109	12/06/07		TCL VOCs, TPH-GRO
RW-120607-JH-034	Equipment Blank	12/06/07	Equipment Blank	TCL VOCs, TPH-GRO
GW-120607-JH-035	MW-111	12/06/07		TCL VOCs, TPH-GRO
GW-120607-JH-036	MW-111	12/06/07	Duplicate	TCL VOCs, TPH-GRO
GW-120607-JH-037	MW-110	12/06/07		TCL VOCs, TPH-GRO
GW-120607-JH-038	MW-110	12/06/07	Duplicate	TCL VOCs, TPH-GRO

<sup>1</sup> QA/QC - quality assurance/quality control

MS/MSD - matrix spike/matrix spike duplicate

<sup>2</sup> TCL VOCs - target compound list volatile organic compounds

TPH-GRO - total petroleum hydrocarbon gasoline range organics

# TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

## ANALYTICAL REPORT

PROJECT NO. 12559

GM FAIRFAX SSOW# E095007  
SDG #: 7L06288

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TESTAMERICA LABORATORIES, INC.

*Denise Heckler*  
Denise D. Heckler  
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December 26, 2007

TestAmerica North Canton

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## **CASE NARRATIVE**

7L06288

The following report contains the analytical results for forty water samples and one quality control sample submitted to TestAmerica North Canton by Conestoga-Rovers & Associates, Inc. from the GM Fairfax SSOW# E095007 Site, project number 12559. The samples were received December 06, 2007 and December 07, 2007, according to documented sample acceptance procedures.

This SDG consists of (2) laboratory IDs: A7L060288 and A7L070423.

TestAmerica utilizes USEPA approved methods in all analytical work. The samples presented in this report were analyzed for the parameter(s) listed on the analytical methods summary page in accordance with the method(s) indicated. A summary of QC data for these analyses is included at the back of the report.

TestAmerica North Canton attests to the validity of the laboratory data generated by TestAmerica facilities reported herein. All analyses performed by TestAmerica facilities were done using established laboratory SOPs that incorporate QA/QC procedures described in the applicable methods. TestAmerica's operations groups have reviewed the data for compliance with the laboratory QA/QC plan, and data have been found to be compliant with laboratory protocols unless otherwise noted below.

The test results in this report meet all NELAP requirements for parameters for which accreditation is required or available. Any exceptions to NELAP requirements are noted in this report. Pursuant to NELAP, this report may not be reproduced, except in full, without the written approval of the laboratory.

Please refer to the Quality Control Elements Narrative following this case narrative for additional quality control information.

If you have any questions, please call the Project Manager, Denise D. Heckler, at 330-497-9396.

This report is sequentially paginated. The final page of the report is labeled as "END OF REPORT." The total number of pages in this report is 153.

## **SUPPLEMENTAL QC INFORMATION**

### **SAMPLE RECEIVING**

The temperatures of the coolers upon sample receipt were 2.7 and 4.1°C.

See TestAmerica's Cooler Receipt Form for additional information.

## **CASE NARRATIVE (continued)**

### **GC/MS VOLATILES**

The sample(s) that contain results between the MDL and the RL were flagged with "J". There is a possibility of false positive or mis-identification at these quantitation levels. In analytical methods requiring confirmation of the analyte reported, confirmation was performed only down to the standard reporting limit (SRL). The acceptance criteria for QC samples may not be met at these quantitation levels.

### **PURGEABLE PETROLEUM HYDROCARBONS-8015**

The sample(s) that contain results between the MDL and the RL were flagged with "J". There is a possibility of false positive or mis-identification at these quantitation levels. In analytical methods requiring confirmation of the analyte reported, confirmation was performed only down to the standard reporting limit (SRL). The acceptance criteria for QC samples may not be met at these quantitation levels.

## QUALITY CONTROL ELEMENTS NARRATIVE

TestAmerica North Canton (formerly STL North Canton) conducts a quality assurance/quality control (QA/QC) program designed to provide scientifically valid and legally defensible data. Toward this end, several types of quality control indicators are incorporated into the QA/QC program, which is described in detail in QA Policy, QA-003. These indicators are introduced into the sample testing process to provide a mechanism for the assessment of the analytical data.

### QC BATCH

Environmental samples are taken through the testing process in groups called QUALITY CONTROL BATCHES (QC batches). A QC batch contains up to twenty environmental samples of a similar matrix (water, soil) that are processed using the same reagents and standards. TestAmerica North Canton (formerly STL North Canton) requires that each environmental sample be associated with a QC batch.

Several quality control samples are included in each QC batch and are processed identically to the twenty environmental samples.

For SW846/RCRA methods, QC samples include a METHOD BLANK (MB), a LABORATORY CONTROL SAMPLE (LCS) and, where appropriate, a MATRIX SPIKE/MATRIX SPIKE DUPLICATE (MS/MSD) pair or a MATRIX SPIKE/SAMPLE DUPLICATE (MS/DU) pair. If there is insufficient sample to perform an MS/MSD or an MS/DU, then a LABORATORY CONTROL SAMPLE DUPLICATE (LCSD) is included in the QC batch.

For 600 series/CWA methods, QC samples include a METHOD BLANK (MB), a LABORATORY CONTROL SAMPLE (LCS) and, where appropriate, a MATRIX SPIKE (MS). An MS is prepared and analyzed at a 10% frequency for GC Methods and at a 5% frequency for GC/MS methods.

### LABORATORY CONTROL SAMPLE

The Laboratory Control Sample is a QC sample that is created by adding known concentrations of a full or partial set of target analytes to a matrix similar to that of the environmental samples in the QC batch. Multi peak responders may not be included in the target spike list due to co-elution. The LCS analyte recovery results are used to monitor the analytical process and provide evidence that the laboratory is performing the method within acceptable guidelines. All control analytes indicated by a bold type in the LCS must meet acceptance criteria. Failure to meet the established recovery guidelines requires the repreparation and reanalysis of all samples in the QC batch. Comparison of only the failed parameters from the first batch are evaluated. The only exception to the rework requirement is that if the LCS recoveries are biased high and the associated sample is ND (non-detected) for the parameter(s) of interest, the batch is acceptable.

At times, a Laboratory Control Sample Duplicate (LCSD) is also included in the QC batch. An LCSD is a QC sample that is created and handled identically to the LCS. Analyte recovery data from the LCSD is assessed in the same way as that of the LCS. The LCSD recoveries, together with the LCS recoveries, are used to determine the reproducibility (precision) of the analytical system. Precision data are expressed as relative percent differences (RPDs). If the RPD fails for an LCS/LCSD and yet the recoveries are within acceptance criteria, the batch is still acceptable.

### METHOD BLANK

The Method Blank is a QC sample consisting of all the reagents used in analyzing the environmental samples contained in the QC batch. Method Blank results are used to determine if interference or contamination in the analytical system could lead to the reporting of false positive data or elevated analyte concentrations. All target analytes must be below the reporting limits (RL) or the associated sample(s) must be ND except under the following circumstances:

- Common organic contaminants may be present at concentrations up to 5 times the reporting limits. Common metals contaminants may be present at concentrations up to 2 times the reporting limit, or the reported blank concentration must be twenty fold less than the concentration reported in the associated environmental samples. (See common laboratory contaminants listed in the table.)

Volatile (GC or GC/MS)	Semivolatile (GC/MS)	Metals ICP-MS	Metals ICP Trace
Methylene Chloride, Acetone, 2-Butanone	Phthalate Esters	Copper, Iron, Zinc, Lead, Calcium, Magnesium, Potassium, Sodium, Barium, Chromium, Manganese	Copper, Iron, Zinc, Lead

## QUALITY CONTROL ELEMENTS NARRATIVE (continued)

- Organic blanks will be accepted if compounds detected in the blank are present in the associated samples at levels 10 times the blank level. Inorganic blanks will be accepted if elements detected in the blank are present in the associated samples at 20 times the blank level.
- Blanks will be accepted if the compounds/elements detected are not present in any of the associated environmental samples.

Failure to meet these Method Blank criteria requires the repreparation and reanalysis of all samples in the QC batch.

### **MATRIX SPIKE/MATRIX SPIKE DUPLICATE**

A Matrix Spike and a Matrix Spike Duplicate are a pair of environmental samples to which known concentrations of a full or partial set of target analytes are added. The MS/MSD results are determined in the same manner as the results of the environmental sample used to prepare the MS/MSD. The analyte recoveries and the relative percent differences (RPDs) of the recoveries are calculated and used to evaluate the effect of the sample matrix on the analytical results. Due to the potential variability of the matrix of each sample, the MS/MSD results may not have an immediate bearing on any samples except the one spiked; therefore, the associated batch MS/MSD may not reflect the same compounds as the samples contained in the analytical report. When these MS/MSD results fail to meet acceptance criteria, the data is evaluated. If the LCS is within acceptance criteria, the batch is considered acceptable.

For certain methods, a Matrix Spike/Sample Duplicate (MS/DU) may be included in the QC batch in place of the MS/MSD. For the parameters (i.e. pH, ignitability) where it is not possible to prepare a spiked sample, a Sample Duplicate may be included in the QC batch. However, a Sample Duplicate is less likely to provide usable precision statistics depending on the likelihood of finding concentrations below the standard reporting limit. When the Sample Duplicate result fails to meet acceptance criteria, the data is evaluated.

For certain methods (600 series methods/CWA), a Matrix Spike is required in place of a Matrix Spike/Matrix Spike Duplicate (MS/MSD) or Matrix Spike/Sample Duplicate (MS/DU).

The acceptance criteria do not apply to samples that are diluted.

### **SURROGATE COMPOUNDS**

In addition to these batch-related QC indicators, each organic environmental and QC sample is spiked with surrogate compounds. Surrogates are organic chemicals that behave similarly to the analytes of interest and that are rarely present in the environment. Surrogate recoveries are used to monitor the individual performance of a sample in the analytical system.

If surrogate recoveries are biased high in the LCS, LCSD, or the Method Blank, and the associated sample(s) are ND, the batch is acceptable. Otherwise, if the LCS, LCSD, or Method Blank surrogate(s) fail to meet recovery criteria, the entire sample batch is reprepared and reanalyzed. If the surrogate recoveries are outside criteria for environmental samples, the samples will be reprepared and reanalyzed unless there is objective evidence of matrix interference or if the sample dilution is greater than the threshold outlined in the associated method SOP.

The acceptance criteria do not apply to samples that are diluted. All other surrogate recoveries will be reported.

For the GC/MS BNA methods, the surrogate criterion is that two of the three surrogates for each fraction must meet acceptance criteria. The third surrogate must have a recovery of ten percent or greater.

For the Pesticide and PCB methods, the surrogate criterion is that one of two surrogate compounds must meet acceptance criteria. The second surrogate must have a recovery of 10% or greater.



### **TestAmerica North Canton (formerly STL North Canton) Certifications and Approvals:**

California (#01144CA), Connecticut (#PH-0590), Florida (#E87225),  
Illinois (#200004), Kansas (#E10336), Minnesota (#39-999-348), New Jersey (#OH001), New York (#10975), Ohio VAP (#CL0024), West Virginia (#210), Wisconsin (#999518190), NAVY, ARMY, USDA Soil Permit,

## EXECUTIVE SUMMARY - Detection Highlights

7L06288 : A7L060288

<u>PARAMETER</u>	<u>RESULT</u>	<u>REPORTING LIMIT</u>	<u>UNITS</u>	<u>ANALYTICAL METHOD</u>
GW-120407-JH-001 12/04/07 09:00 001				
TPH (as Gasoline)	60 J	100	ug/L	SW846 8015B
Toluene	0.22 J	1.0	ug/L	SW846 8260B
GW-120407-JH-002 12/04/07 09:55 002				
TPH (as Gasoline)	33 J	100	ug/L	SW846 8015B
GW-120407-JH-003 12/04/07 11:00 003				
TPH (as Gasoline)	45 J	100	ug/L	SW846 8015B
GW-120407-JH-004 12/04/07 11:30 004				
TPH (as Gasoline)	45 J	100	ug/L	SW846 8015B
cis-1,2-Dichloroethene	1.3	0.50	ug/L	SW846 8260B
GW-120407-JH-005 12/04/07 13:20 005				
TPH (as Gasoline)	30 J	100	ug/L	SW846 8015B
GW-120407-JH-006 12/04/07 14:40 006				
TPH (as Gasoline)	41 J	100	ug/L	SW846 8015B
cis-1,2-Dichloroethene	3.1	0.50	ug/L	SW846 8260B
Trichloroethene	0.64 J	1.0	ug/L	SW846 8260B
GW-120407-JH-007 12/04/07 14:55 007				
TPH (as Gasoline)	40 J	100	ug/L	SW846 8015B
cis-1,2-Dichloroethene	2.8	0.50	ug/L	SW846 8260B
Vinyl chloride	0.36 J	1.0	ug/L	SW846 8260B
GW-120407-JH-008 12/04/07 15:45 008				
TPH (as Gasoline)	36 J	100	ug/L	SW846 8015B
GW-120407-JH-009 12/04/07 16:00 009				
TPH (as Gasoline)	33 J	100	ug/L	SW846 8015B

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## EXECUTIVE SUMMARY - Detection Highlights

7L06288 : A7L060288

PARAMETER	RESULT	REPORTING LIMIT	UNITS	ANALYTICAL METHOD
<b>GW-120407-JH-010 12/04/07 16:50 010</b>				
TPH (as Gasoline)	33 J	100	ug/L	SW846 8015B
<b>GW-120507-JH-011 12/05/07 08:55 011</b>				
TPH (as Gasoline)	36 J	100	ug/L	SW846 8015B
<b>GW-120507-JH-012 12/05/07 09:15 012</b>				
TPH (as Gasoline)	35 J	100	ug/L	SW846 8015B
Chlorobenzene	0.23 J	1.0	ug/L	SW846 8260B
cis-1,2-Dichloroethene	0.22 J	0.50	ug/L	SW846 8260B
<b>RW-120507-JH-013 12/05/07 09:40 013</b>				
TPH (as Gasoline)	31 J	100	ug/L	SW846 8015B
Chloroform	0.27 J	1.0	ug/L	SW846 8260B
<b>GW-120507-JH-014 12/05/07 10:50 014</b>				
TPH (as Gasoline)	32 J	100	ug/L	SW846 8015B
<b>GW-120507-JH-015 12/05/07 11:10 015</b>				
TPH (as Gasoline)	30 J	100	ug/L	SW846 8015B

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## EXECUTIVE SUMMARY - Detection Highlights

7L06288 : A7L070423

<u>PARAMETER</u>	<u>RESULT</u>	<u>REPORTING LIMIT</u>	<u>UNITS</u>	<u>ANALYTICAL METHOD</u>
<b>GW-120507-JH-016 12/05/07 11:30 001</b>				
TPH (as Gasoline)	33 J	100	ug/L	SW846 8015B
<b>GW-120507-JH-017 12/05/07 13:05 002</b>				
TPH (as Gasoline)	28 J	100	ug/L	SW846 8015B
<b>GW-120507-JH-018 12/05/07 13:50 003</b>				
cis-1,2-Dichloroethene	3.7	0.50	ug/L	SW846 8260B
Tetrachloroethene	0.35 J	1.0	ug/L	SW846 8260B
Trichloroethene	0.75 J	1.0	ug/L	SW846 8260B
<b>GW-120507-JH-019 12/05/07 14:20 004</b>				
cis-1,2-Dichloroethene	0.22 J	0.50	ug/L	SW846 8260B
<b>GW-120507-JH-020 12/05/07 14:50 005</b>				
TPH (as Gasoline)	31 J	100	ug/L	SW846 8015B
<b>GW-120507-JH-021 12/05/07 15:05 006</b>				
TPH (as Gasoline)	31 J	100	ug/L	SW846 8015B
cis-1,2-Dichloroethene	6.4	0.50	ug/L	SW846 8260B
Toluene	0.26 J	1.0	ug/L	SW846 8260B
Vinyl chloride	1.0	1.0	ug/L	SW846 8260B
<b>RW-120507-JH-022 12/05/07 15:20 007</b>				
TPH (as Gasoline)	33 J	100	ug/L	SW846 8015B
Acetone	5.6 J	10	ug/L	SW846 8260B
<b>GW-120507-JH-023 12/05/07 16:05 008</b>				
TPH (as Gasoline)	33 J	100	ug/L	SW846 8015B
cis-1,2-Dichloroethene	1.9	0.50	ug/L	SW846 8260B
Tetrachloroethene	9.6	1.0	ug/L	SW846 8260B
Trichloroethene	1.4	1.0	ug/L	SW846 8260B

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## EXECUTIVE SUMMARY - Detection Highlights

7L06288 : A7L070423

PARAMETER	RESULT	REPORTING LIMIT	UNITS	ANALYTICAL METHOD
<b>GW-120607-JH-025 12/06/07 09:15 010</b>				
TPH (as Gasoline)	31 J	100	ug/L	SW846 8015B
<b>GW-120607-JH-026 12/06/07 09:30 011</b>				
TPH (as Gasoline)	33 J	100	ug/L	SW846 8015B
<b>GW-120607-JH-028 12/06/07 10:45 013</b>				
TPH (as Gasoline)	30 J	100	ug/L	SW846 8015B
<b>GW-120607-JH-029 12/06/07 11:15 014</b>				
cis-1,2-Dichloroethene	0.36 J	0.50	ug/L	SW846 8260B
<b>GW-120607-JH-030 12/06/07 12:00 015</b>				
TPH (as Gasoline)	470000	100000	ug/L	SW846 8015B
Ethylbenzene	8000	5000	ug/L	SW846 8260B
4-Methyl-2-pentanone	160000	50000	ug/L	SW846 8260B
Toluene	150000	5000	ug/L	SW846 8260B
Xylenes (total)	57000	5000	ug/L	SW846 8260B
<b>GW-120607-JH-031 12/06/07 12:30 016</b>				
TPH (as Gasoline)	540000	100000	ug/L	SW846 8015B
Ethylbenzene	7500	5000	ug/L	SW846 8260B
4-Methyl-2-pentanone	160000	50000	ug/L	SW846 8260B
Toluene	140000	5000	ug/L	SW846 8260B
Xylenes (total)	56000	5000	ug/L	SW846 8260B
<b>GW-120607-JH-032 12/06/07 13:40 017</b>				
TPH (as Gasoline)	50 J	100	ug/L	SW846 8015B
cis-1,2-Dichloroethene	0.48 J	0.50	ug/L	SW846 8260B
Tetrachloroethene	0.47 J	1.0	ug/L	SW846 8260B
<b>GW-120607-JH-033 12/06/07 14:20 018</b>				
TPH (as Gasoline)	56 J	100	ug/L	SW846 8015B
cis-1,2-Dichloroethene	20	0.50	ug/L	SW846 8260B
trans-1,2-Dichloroethene	1.5	0.50	ug/L	SW846 8260B
Vinyl chloride	0.29 J	1.0	ug/L	SW846 8260B

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## EXECUTIVE SUMMARY - Detection Highlights

7L06288 : A7L070423

<u>PARAMETER</u>	<u>RESULT</u>	<u>REPORTING LIMIT</u>	<u>UNITS</u>	<u>ANALYTICAL METHOD</u>
<b>RW-120607-JH-034 12/06/07 14:30 019</b>				
TPH (as Gasoline)	45 J	100	ug/L	SW846 8015B
Chloroform	0.55 J	1.0	ug/L	SW846 8260B
<b>GW-120607-JH-035 12/06/07 15:10 020</b>				
TPH (as Gasoline)	130	100	ug/L	SW846 8015B
cis-1,2-Dichloroethene	88	1.7	ug/L	SW846 8260B
trans-1,2-Dichloroethene	6.2	1.7	ug/L	SW846 8260B
Methylene chloride	1.7 J	3.3	ug/L	SW846 8260B
Vinyl chloride	7.0	3.3	ug/L	SW846 8260B
<b>GW-120607-JH-036 12/06/07 15:30 021</b>				
TPH (as Gasoline)	120	100	ug/L	SW846 8015B
cis-1,2-Dichloroethene	89	1.7	ug/L	SW846 8260B
trans-1,2-Dichloroethene	6.0	1.7	ug/L	SW846 8260B
Methylene chloride	2.8 J	3.3	ug/L	SW846 8260B
Vinyl chloride	6.7	3.3	ug/L	SW846 8260B
<b>GW-120607-JH-037 12/06/07 15:55 022</b>				
TPH (as Gasoline)	29 J	100	ug/L	SW846 8015B
cis-1,2-Dichloroethene	0.33 J	0.50	ug/L	SW846 8260B
<b>GW-120607-JH-038 12/06/07 16:30 023</b>				
cis-1,2-Dichloroethene	0.35 J	0.50	ug/L	SW846 8260B
<b>TRIP BLANK B COOLER 68187 12/06/07 024</b>				
Acetone	1.9 J	10	ug/L	SW846 8260B

## **ANALYTICAL METHODS SUMMARY**

7L06288

<u>PARAMETER</u>	<u>ANALYTICAL METHOD</u>
Volatile Organics by GC/MS	SW846 8260B
Volatile Petroleum Hydrocarbons	SW846 8015B

**References:**

SW846 "Test Methods for Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 and its updates.

## SAMPLE SUMMARY

7L06288 : A7L060288

WO #	SAMPLE#	CLIENT SAMPLE ID	SAMPLED DATE	SAMP TIME
KDLKE	001	GW-120407-JH-001	12/04/07	09:00
KDLKW	002	GW-120407-JH-002	12/04/07	09:55
KDLK1	003	GW-120407-JH-003	12/04/07	11:00
KDLK3	004	GW-120407-JH-004	12/04/07	11:30
KDLK6	005	GW-120407-JH-005	12/04/07	13:20
KDLK8	006	GW-120407-JH-006	12/04/07	14:40
KDLLA	007	GW-120407-JH-007	12/04/07	14:55
KDLLJ	008	GW-120407-JH-008	12/04/07	15:45
KDLLN	009	GW-120407-JH-009	12/04/07	16:00
KDLLQ	010	GW-120407-JH-010	12/04/07	16:50
KDLLV	011	GW-120507-JH-011	12/05/07	08:55
KDLL4	012	GW-120507-JH-012	12/05/07	09:15
KDLL8	013	RW-120507-JH-013	12/05/07	09:40
KDLMD	014	GW-120507-JH-014	12/05/07	10:50
KDLMF	015	GW-120507-JH-015	12/05/07	11:10
KDLMH	016	TRIP BLANK-A	12/04/07	

### NOTE (S) :

- The analytical results of the samples listed above are presented on the following pages.
- All calculations are performed before rounding to avoid round-off errors in calculated results.
- Results noted as "ND" were not detected at or above the stated limit.
- This report must not be reproduced, except in full, without the written approval of the laboratory.
- Results for the following parameters are never reported on a dry weight basis: color, corrosivity, density, flashpoint, ignitability, layers, odor, paint filter test, pH, porosity pressure, reactivity, redox potential, specific gravity, spot tests, solids, solubility, temperature, viscosity, and weight.

(Continued on next page)

## SAMPLE SUMMARY

7L06288 : A7L070423

WO #	SAMPLE#	CLIENT SAMPLE ID	SAMPLED DATE	SAMP TIME
KDQPJ	001	GW-120507-JH-016	12/05/07	11:30
KDQPP	002	GW-120507-JH-017	12/05/07	13:05
KDQPQ	003	GW-120507-JH-018	12/05/07	13:50
KDQPT	004	GW-120507-JH-019	12/05/07	14:20
KDQPV	005	GW-120507-JH-020	12/05/07	14:50
KDQPX	006	GW-120507-JH-021	12/05/07	15:05
KDQP1	007	RW-120507-JH-022	12/05/07	15:20
KDQP2	008	GW-120507-JH-023	12/05/07	16:05
KDQP4	009	GW-120507-JH-024	12/05/07	16:45
KDQP6	010	GW-120607-JH-025	12/06/07	09:15
KDQQA	011	GW-120607-JH-026	12/06/07	09:30
KDQQC	012	GW-120607-JH-027	12/06/07	10:30
KDQQE	013	GW-120607-JH-028	12/06/07	10:45
KDQQG	014	GW-120607-JH-029	12/06/07	11:15
KDQQH	015	GW-120607-JH-030	12/06/07	12:00
KDQQK	016	GW-120607-JH-031	12/06/07	12:30
KDQQM	017	GW-120607-JH-032	12/06/07	13:40
KDQQN	018	GW-120607-JH-033	12/06/07	14:20
KDQQP	019	RW-120607-JH-034	12/06/07	14:30
KDQQR	020	GW-120607-JH-035	12/06/07	15:10
KDQQT	021	GW-120607-JH-036	12/06/07	15:30
KDQQW	022	GW-120607-JH-037	12/06/07	15:55
KDQQO	023	GW-120607-JH-038	12/06/07	16:30
KDQQ2	024	TRIP BLANK B COOLER 68187	12/06/07	
KDQRR	025	TRIP BLANK C COOLER 68182	12/06/07	

**NOTE (S) :**

- The analytical results of the samples listed above are presented on the following pages.
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- Results noted as "ND" were not detected at or above the stated limit.
- This report must not be reproduced, except in full, without the written approval of the laboratory.
- Results for the following parameters are never reported on a dry weight basis: color, corrosivity, density, flashpoint, ignitability, layers, odor, paint filter test, pH, porosity pressure, reactivity, redox potential, specific gravity, spot tests, solids, solubility, temperature, viscosity, and weight.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120407-JH-001

## GC/MS Volatiles

Lot-Sample #....: A7L060288-001    Work Order #....: KDLKE1AA    Matrix.....: WG  
 Date Sampled....: 12/04/07 09:00    Date Received...: 12/06/07  
 Prep Date.....: 12/11/07    Analysis Date...: 12/11/07  
 Prep Batch #....: 7346476  
 Dilution Factor: 1    Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	0.22 J	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY
		LIMITS
Dibromofluoromethane	94	(73 - 122)
1,2-Dichloroethane-d4	96	(61 - 128)
Toluene-d8	96	(76 - 110)
4-Bromofluorobenzene	93	(74 - 116)

(Continued on next page)

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120407-JH-001

GC/MS Volatiles

Lot-Sample #...: A7L060288-001 Work Order #...: KDLKE1AA Matrix.....: WG

NOTE (S) :

J Estimated result. Result is less than RL.

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120407-JH-001

GC Volatiles

Lot-Sample #....: A7L060288-001 Work Order #....: KDLKE1AC Matrix.....: WG  
Date Sampled...: 12/04/07 09:00 Date Received...: 12/06/07  
Prep Date.....: 12/07/07 Analysis Date...: 12/07/07  
Prep Batch #....: 7344082  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	60 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	97	(10 - 150)		

NOTE (S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120407-JH-002

## GC/MS Volatiles

Lot-Sample #....: A7L060288-002    Work Order #....: KDLKW1AA    Matrix.....: WG  
 Date Sampled....: 12/04/07 09:55    Date Received...: 12/06/07  
 Prep Date.....: 12/12/07    Analysis Date...: 12/12/07  
 Prep Batch #....: 7346476  
 Dilution Factor: 1    Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28
SURROGATE		PERCENT RECOVERY	RECOVERY LIMITS	
Dibromofluoromethane	91	(73 - 122)		
1,2-Dichloroethane-d4	92	(61 - 128)		
Toluene-d8	98	(76 - 110)		
4-Bromofluorobenzene	95	(74 - 116)		

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120407-JH-002

GC Volatiles

Lot-Sample #....: A7L060288-002 Work Order #....: KDLKW1AC Matrix.....: WG  
Date Sampled...: 12/04/07 09:55 Date Received..: 12/06/07  
Prep Date.....: 12/07/07 Analysis Date...: 12/07/07  
Prep Batch #....: 7344082  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	33 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	100	(10 - 150)		

NOTE (S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120407-JH-003

## GC/MS Volatiles

Lot-Sample #....: A7L060288-003 Work Order #....: KDLK11AA Matrix.....: WG  
 Date Sampled...: 12/04/07 11:00 Date Received...: 12/06/07  
 Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
 Prep Batch #....: 7346476  
 Dilution Factor: 1 Method.....: SW846 8260B

<u>PARAMETER</u>	<u>RESULT</u>	<u>REPORTING</u>		
		<u>LIMIT</u>	<u>UNITS</u>	<u>MDL</u>
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28
<u>SURROGATE</u>		<u>PERCENT</u>	<u>RECOVERY</u>	
		<u>RECOVERY</u>	<u>LIMITS</u>	
Dibromofluoromethane		96	(73 - 122)	
1,2-Dichloroethane-d4		98	(61 - 128)	
Toluene-d8		97	(76 - 110)	
4-Bromofluorobenzene		94	(74 - 116)	

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120407-JH-003

GC Volatiles

Lot-Sample #....: A7L060288-003 Work Order #....: KDLK11AC Matrix.....: WG  
Date Sampled....: 12/04/07 11:00 Date Received...: 12/06/07  
Prep Date.....: 12/08/07 Analysis Date...: 12/08/07  
Prep Batch #....: 7344082  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	45 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	99	(10 - 150)		

NOTE (S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120407-JH-004

## GC/MS Volatiles

Lot-Sample #....: A7L060288-004    Work Order #....: KDLK31AA    Matrix.....: WG  
 Date Sampled....: 12/04/07 11:30    Date Received...: 12/06/07  
 Prep Date.....: 12/12/07    Analysis Date...: 12/12/07  
 Prep Batch #....: 7346476  
 Dilution Factor: 1    Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	1.3	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY	
		LIMITS	
Dibromofluoromethane	95	(73 - 122)	
1,2-Dichloroethane-d4	96	(61 - 128)	
Toluene-d8	97	(76 - 110)	
4-Bromofluorobenzene	96	(74 - 116)	

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120407-JH-004

GC Volatiles

Lot-Sample #....: A7L060288-004 Work Order #....: KDLK31AC Matrix.....: WG  
Date Sampled...: 12/04/07 11:30 Date Received...: 12/06/07  
Prep Date.....: 12/08/07 Analysis Date...: 12/08/07  
Prep Batch #....: 7344082  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		MDL
		LIMIT	UNITS	
TPH (as Gasoline)	45 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	100	(10 - 150)		

NOTE(S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120407-JH-005

## GC/MS Volatiles

Lot-Sample #....: A7L060288-005    Work Order #....: KDLK61AA    Matrix.....: WG  
 Date Sampled....: 12/04/07 13:20    Date Received...: 12/06/07  
 Prep Date.....: 12/12/07    Analysis Date...: 12/12/07  
 Prep Batch #....: 7346476  
 Dilution Factor: 1    Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY	
		LIMITS	
Dibromofluoromethane	93	(73 - 122)	
1,2-Dichloroethane-d4	94	(61 - 128)	
Toluene-d8	94	(76 - 110)	
4-Bromofluorobenzene	93	(74 - 116)	

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120407-JH-005**

**GC Volatiles**

Lot-Sample #....: A7L060288-005 Work Order #....: KDLK61AC Matrix.....: WG  
Date Sampled...: 12/04/07 13:20 Date Received..: 12/06/07  
Prep Date.....: 12/08/07 Analysis Date...: 12/08/07  
Prep Batch #....: 7344082  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	30 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	96	(10 - 150)		

**NOTE (S) :**

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120407-JH-006

## GC/MS Volatiles

Lot-Sample #...: A7L060288-006 Work Order #...: KDLK81AA Matrix.....: WG  
 Date Sampled...: 12/04/07 14:40 Date Received...: 12/06/07  
 Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
 Prep Batch #...: 7346476  
 Dilution Factor: 1 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	3.1	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	0.64 J	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY	
		LIMITS	
Dibromofluoromethane	97	(73 - 122)	
1,2-Dichloroethane-d4	98	(61 - 128)	
Toluene-d8	98	(76 - 110)	
4-Bromofluorobenzene	95	(74 - 116)	

(Continued on next page)

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120407-JH-006

GC/MS Volatiles

Lot-Sample #....: A7L060288-006 Work Order #....: KDLK81AA Matrix.....: WG

NOTE (S) :

J Estimated result. Result is less than RL.

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120407-JH-006

GC Volatiles

Lot-Sample #....: A7L060288-006 Work Order #....: KDLK81AC Matrix.....: WG  
Date Sampled...: 12/04/07 14:40 Date Received...: 12/06/07  
Prep Date.....: 12/08/07 Analysis Date...: 12/08/07  
Prep Batch #....: 7344082  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	41 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	99	(10 - 150)		

NOTE(S) :

J Estimated result. Result is less than RL.

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120407-JH-007

GC/MS Volatiles

Lot-Sample #....: A7L060288-007    Work Order #....: KDLLA1AA    Matrix.....: WG  
 Date Sampled...: 12/04/07 14:55    Date Received...: 12/06/07  
 Prep Date.....: 12/12/07    Analysis Date...: 12/12/07  
 Prep Batch #....: 7346476  
 Dilution Factor: 1    Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	2.8	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	0.36 J	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS	
		(73 - 122)	(61 - 128)
Dibromofluoromethane	97	(73 - 122)	(61 - 128)
1,2-Dichloroethane-d4	98	(61 - 128)	
Toluene-d8	96	(76 - 110)	
4-Bromofluorobenzene	96	(74 - 116)	

(Continued on next page)

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120407-JH-007

GC/MS Volatiles

Lot-Sample #...: A7L060288-007 Work Order #: K DLLA1AA Matrix.....: WG

NOTE (S) :

J Estimated result. Result is less than RL.

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120407-JH-007**

**GC Volatiles**

**Lot-Sample #....: A7L060288-007 Work Order #....: KDLLA1AC Matrix.....: WG  
Date Sampled....: 12/04/07 14:55 Date Received...: 12/06/07  
Prep Date.....: 12/08/07 Analysis Date...: 12/08/07  
Prep Batch #....: 7344082  
Dilution Factor: 1 Method.....: SW846 8015B**

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	40 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	95	(10 - 150)		

**NOTE (S) :**

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120407-JH-008

## GC/MS Volatiles

Lot-Sample #....: A7L060288-008    Work Order #....: KDLLJ1AA    Matrix.....: WG  
 Date Sampled....: 12/04/07 15:45    Date Received...: 12/06/07  
 Prep Date.....: 12/12/07    Analysis Date...: 12/12/07  
 Prep Batch #....: 7346476  
 Dilution Factor: 1    Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY	
		LIMITS	
Dibromofluoromethane	95	(73 - 122)	
1,2-Dichloroethane-d4	97	(61 - 128)	
Toluene-d8	97	(76 - 110)	
4-Bromofluorobenzene	96	(74 - 116)	

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120407-JH-008**

**GC Volatiles**

**Lot-Sample #....: A7L060288-008    Work Order #....: KDLLJ1AC    Matrix.....: WG**  
**Date Sampled....: 12/04/07 15:45    Date Received...: 12/06/07**  
**Prep Date.....: 12/08/07    Analysis Date...: 12/08/07**  
**Prep Batch #....: 7344082**  
**Dilution Factor: 1                  Method.....: SW846 8015B**

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
<b>TPH (as Gasoline)</b>	<b>36 J</b>	<b>100</b>	<b>ug/L</b>	<b>28</b>
<b>SURROGATE</b>	<b>PERCENT</b>	<b>RECOVERY</b>		
Trifluorotoluene	RECOVERY	LIMITS		
	99	(10 - 150)		

**NOTE (S) :**

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120407-JH-009

## GC/MS Volatiles

Lot-Sample #....: A7L060288-009 Work Order #....: KDLLN1AA Matrix.....: WG  
 Date Sampled....: 12/04/07 16:00 Date Received...: 12/06/07  
 Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
 Prep Batch #....: 7346476  
 Dilution Factor: 1 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		MDL
		LIMIT	UNITS	
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY
		LIMITS
Dibromofluoromethane	94	(73 - 122)
1,2-Dichloroethane-d4	94	(61 - 128)
Toluene-d8	96	(76 - 110)
4-Bromofluorobenzene	96	(74 - 116)

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120407-JH-009

GC Volatiles

Lot-Sample #....: A7L060288-009 Work Order #....: KDLLN1AE Matrix.....: WG  
Date Sampled....: 12/04/07 16:00 Date Received...: 12/06/07  
Prep Date.....: 12/08/07 Analysis Date...: 12/08/07  
Prep Batch #....: 7344082  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	33 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	100	(10 - 150)		

NOTE(S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120407-JH-010

## GC/MS Volatiles

Lot-Sample #...: A7L060288-010 Work Order #...: KDLLQ1AA Matrix.....: WG  
 Date Sampled...: 12/04/07 16:50 Date Received..: 12/06/07  
 Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
 Prep Batch #...: 7346476  
 Dilution Factor: 1 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS
Dibromofluoromethane	98	(73 - 122)
1,2-Dichloroethane-d4	99	(61 - 128)
Toluene-d8	94	(76 - 110)
4-Bromofluorobenzene	96	(74 - 116)

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120407-JH-010**

**GC Volatiles**

**Lot-Sample #....:** A7L060288-010    **Work Order #....:** KDLLQ1AC    **Matrix.....:** WG  
**Date Sampled....:** 12/04/07 16:50    **Date Received..:** 12/06/07  
**Prep Date.....:** 12/08/07    **Analysis Date..:** 12/08/07  
**Prep Batch #....:** 7344082  
**Dilution Factor:** 1    **Method.....:** SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	33 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	91	(10 - 150)		

**NOTE(S) :**

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120507-JH-011

## GC/MS Volatiles

Lot-Sample #....: A7L060288-011 Work Order #....: KDLLV1AA Matrix.....: WG  
 Date Sampled...: 12/05/07 08:55 Date Received...: 12/06/07  
 Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
 Prep Batch #....: 7346476  
 Dilution Factor: 1 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY	
		LIMITS	
Dibromofluoromethane	97	(73 - 122)	
1,2-Dichloroethane-d4	96	(61 - 128)	
Toluene-d8	96	(76 - 110)	
4-Bromofluorobenzene	95	(74 - 116)	

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120507-JH-011**

**GC Volatiles**

Lot-Sample #....: A7L060288-011 Work Order #....: KDLLV1AC Matrix.....: WG  
Date Sampled...: 12/05/07 08:55 Date Received...: 12/06/07  
Prep Date.....: 12/08/07 Analysis Date...: 12/08/07  
Prep Batch #....: 7344082  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	36 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	93	(10 - 150)		

**NOTE(S) :**

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120507-JH-012

## GC/MS Volatiles

Lot-Sample #....: A7L060288-012    Work Order #....: KDLL41AA    Matrix.....: WG  
 Date Sampled....: 12/05/07 09:15    Date Received...: 12/06/07  
 Prep Date.....: 12/12/07    Analysis Date...: 12/12/07  
 Prep Batch #....: 7346476  
 Dilution Factor: 1    Method.....: SW846 .8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	0.23 J	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	0.22 J	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY	
		LIMITS	
Dibromofluoromethane	96	(73	- 122)
1,2-Dichloroethane-d4	95	(61	- 128)
Toluene-d8	95	(76	- 110)
4-Bromofluorobenzene	94	(74	- 116)

(Continued on next page)

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120507-JH-012**

**GC/MS Volatiles**

**Lot-Sample #...: A7L060288-012 Work Order #...: KDLL41AA Matrix.....: WG**

**NOTE (S) :**

J Estimated result. Result is less than RL.

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120507-JH-012**

**GC Volatiles**

Lot-Sample #...: A7L060288-012 Work Order #...: KDLL41AC Matrix.....: WG  
Date Sampled...: 12/05/07 09:15 Date Received...: 12/06/07  
Prep Date.....: 12/08/07 Analysis Date...: 12/08/07  
Prep Batch #...: 7344082  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		MDL
		LIMIT	UNITS	
TPH (as Gasoline)	35 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	95	(10 - 150)		

**NOTE (S) :**

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: RW-120507-JH-013

## GC/MS Volatiles

Lot-Sample #....: A7L060288-013 Work Order #....: KDLL81AA Matrix.....: WG  
 Date Sampled...: 12/05/07 09:40 Date Received...: 12/06/07  
 Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
 Prep Batch #....: 7346476  
 Dilution Factor: 1 Method.....: SW846 8260B

<u>PARAMETER</u>	<u>RESULT</u>	<u>REPORTING</u>		<u>MDL</u>
		<u>LIMIT</u>	<u>UNITS</u>	
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
<b>Chloroform</b>	<b>0.27 J</b>	<b>1.0</b>	<b>ug/L</b>	<b>0.16</b>
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

<u>SURROGATE</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY</u>
		<u>LIMITS</u>
Dibromofluoromethane	93	(73 - 122)
1,2-Dichloroethane-d4	95	(61 - 128)
Toluene-d8	95	(76 - 110)
4-Bromofluorobenzene	95	(74 - 116)

(Continued on next page)

Conestoga-Rovers & Associates, Inc.

Client Sample ID: RW-120507-JH-013

GC/MS Volatiles

Lot-Sample #...: A7L060288-013 Work Order #...: KDLL81AA Matrix.....: WG

NOTE (S) :

J Estimated result. Result is less than RL.

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: RW-120507-JH-013**

**GC Volatiles**

**Lot-Sample #....: A7L060288-013    Work Order #....: KDLL81AC    Matrix.....: WG**  
**Date Sampled....: 12/05/07 09:40    Date Received...: 12/06/07**  
**Prep Date.....: 12/08/07    Analysis Date...: 12/08/07**  
**Prep Batch #....: 7344082**  
**Dilution Factor: 1                  Method.....: SW846 8015B**

<b>PARAMETER</b>	<b>RESULT</b>	<b>REPORTING</b>		
		<b>LIMIT</b>	<b>UNITS</b>	<b>MDL</b>
<b>TPH (as Gasoline)</b>	<b>31 J</b>	<b>100</b>	<b>ug/L</b>	<b>28</b>
<b>SURROGATE</b>	<b>PERCENT</b>	<b>RECOVERY</b>		
Trifluorotoluene	RECOVERY	LIMITS		
	103	(10 - 150)		

**NOTE (S) :**

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120507-JH-014

## GC/MS Volatiles

Lot-Sample #....: A7L060288-014 Work Order #....: KDLMD1AA Matrix.....: WG  
 Date Sampled....: 12/05/07 10:50 Date Received...: 12/06/07  
 Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
 Prep Batch #....: 7346476  
 Dilution Factor: 1 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		MDL
		LIMIT	UNITS	
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	RECOVERY	PERCENT	RECOVERY
		LIMITS	
Dibromofluoromethane	95	(73 - 122)	
1,2-Dichloroethane-d4	98	(61 - 128)	
Toluene-d8	97	(76 - 110)	
4-Bromofluorobenzene	94	(74 - 116)	

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120507-JH-014**

**GC Volatiles**

**Lot-Sample #....: A7L060288-014 Work Order #....: KDLMD1AC Matrix.....: WG  
Date Sampled....: 12/05/07 10:50 Date Received...: 12/06/07  
Prep Date.....: 12/08/07 Analysis Date...: 12/08/07  
Prep Batch #....: 7344082  
Dilution Factor: 1 Method.....: SW846 8015B**

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
<b>TPH (as Gasoline)</b>	<b>32 J</b>	<b>100</b>	<b>ug/L</b>	<b>28</b>
<b>SURROGATE</b>	<b>PERCENT</b>	<b>RECOVERY</b>		
Trifluorotoluene	RECOVERY	LIMITS		
	101	(10 - 150)		

**NOTE (S) :**

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120507-JH-015

## GC/MS Volatiles

Lot-Sample #....: A7L060288-015 Work Order #....: KDLMF1AA Matrix.....: WG  
 Date Sampled....: 12/05/07 11:10 Date Received...: 12/06/07  
 Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
 Prep Batch #....: 7346476  
 Dilution Factor: 1 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28
SURROGATE		PERCENT RECOVERY	RECOVERY LIMITS	
Dibromofluoromethane	95	(73 - 122)		
1,2-Dichloroethane-d4	97	(61 - 128)		
Toluene-d8	96	(76 - 110)		
4-Bromofluorobenzene	98	(74 - 116)		

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120507-JH-015**

**GC Volatiles**

**Lot-Sample #....:** A7L060288-015    **Work Order #....:** KDLMF1AC    **Matrix.....:** WG  
**Date Sampled....:** 12/05/07 11:10    **Date Received...:** 12/06/07  
**Prep Date.....:** 12/08/07    **Analysis Date...:** 12/08/07  
**Prep Batch #....:** 7344082  
**Dilution Factor:** 1    **Method.....:** SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	30 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	95	(10 - 150)		

**NOTE(S) :**

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: TRIP BLANK-A

## GC/MS Volatiles

Lot-Sample #....: A7L060288-016    Work Order #....: KDLMH1AA    Matrix.....: WQ  
 Date Sampled...: 12/04/07    Date Received...: 12/06/07  
 Prep Date.....: 12/12/07    Analysis Date...: 12/12/07  
 Prep Batch #...: 7346476  
 Dilution Factor: 1    Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY	
		LIMITS	
Dibromofluoromethane	96	(73 - 122)	
1,2-Dichloroethane-d4	98	(61 - 128)	
Toluene-d8	96	(76 - 110)	
4-Bromofluorobenzene	95	(74 - 116)	

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120507-JH-016

## GC/MS Volatiles

Lot-Sample #....: A7L070423-001   Work Order #....: KDQPJ1AA      Matrix.....: WG  
 Date Sampled...: 12/05/07 11:30   Date Received...: 12/07/07  
 Prep Date.....: 12/15/07      Analysis Date...: 12/15/07  
 Prep Batch #....: 7351272  
 Dilution Factor: 1              Method.....: SW846 8260B

<u>PARAMETER</u>	<u>RESULT</u>	<u>REPORTING</u>		
		<u>LIMIT</u>	<u>UNITS</u>	<u>MDL</u>
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

<u>SURROGATE</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY</u>
		<u>LIMITS</u>
Dibromofluoromethane	89	(73 - 122)
1,2-Dichloroethane-d4	92	(61 - 128)
Toluene-d8	89	(76 - 110)
4-Bromofluorobenzene	84	(74 - 116)

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120507-JH-016

GC Volatiles

Lot-Sample #....: A7L070423-001 Work Order #....: KDQPJ1AC Matrix.....: WG  
Date Sampled...: 12/05/07 11:30 Date Received...: 12/07/07  
Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
Prep Batch #...: 7346409  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		MDL
		LIMIT	UNITS	
TPH (as Gasoline)	33 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	84	(10 - 150)		

NOTE (S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120507-JH-017

## GC/MS Volatiles

Lot-Sample #....: A7L070423-002   Work Order #....: KDQPP1AA   Matrix.....: WG  
 Date Sampled...: 12/05/07 13:05   Date Received...: 12/07/07  
 Prep Date.....: 12/15/07   Analysis Date...: 12/15/07  
 Prep Batch #....: 7351272  
 Dilution Factor: 1   Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28
SURROGATE		PERCENT RECOVERY	RECOVERY LIMITS	
Dibromofluoromethane	90	(73 - 122)		
1,2-Dichloroethane-d4	93	(61 - 128)		
Toluene-d8	91	(76 - 110)		
4-Bromofluorobenzene	86	(74 - 116)		

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120507-JH-017

GC Volatiles

Lot-Sample #...: A7L070423-002 Work Order #...: KDQPP1AC Matrix.....: WG  
Date Sampled...: 12/05/07 13:05 Date Received...: 12/07/07  
Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
Prep Batch #...: 7346409  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	28 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	85	(10 - 150)		

NOTE (S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120507-JH-018

## GC/MS Volatiles

Lot-Sample #....: A7L070423-003 Work Order #....: KDQPQ1AA Matrix.....: WG  
 Date Sampled....: 12/05/07 13:50 Date Received...: 12/07/07  
 Prep Date.....: 12/15/07 Analysis Date...: 12/15/07  
 Prep Batch #....: 7351272  
 Dilution Factor: 1 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	3.7	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	0.35 J	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	0.75 J	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY	
		LIMITS	
Dibromofluoromethane	91	(73 - 122)	
1,2-Dichloroethane-d4	91	(61 - 128)	
Toluene-d8	91	(76 - 110)	
4-Bromofluorobenzene	85	(74 - 116)	

(Continued on next page)

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120507-JH-018**

**GC/MS Volatiles**

**Lot-Sample #...: A7L070423-003 Work Order #...: KDQPQ1AA Matrix.....: WG**

**NOTE (S) :**

**J Estimated result. Result is less than RL.**

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120507-JH-018**

**GC Volatiles**

**Lot-Sample #....: A7L070423-003   Work Order #....: KDQPQ1AC      Matrix.....: WG**  
**Date Sampled....: 12/05/07 13:50 Date Received...: 12/07/07**  
**Prep Date.....: 12/12/07      Analysis Date...: 12/12/07**  
**Prep Batch #....: 7346409**  
**Dilution Factor: 1      Method.....: SW846 8015B**

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	ND	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	84	(10 - 150)		

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120507-JH-019

## GC/MS Volatiles

Lot-Sample #....: A7L070423-004    Work Order #....: KDQPT1AA    Matrix.....: WG  
 Date Sampled...: 12/05/07 14:20    Date Received...: 12/07/07  
 Prep Date.....: 12/15/07    Analysis Date...: 12/15/07  
 Prep Batch #....: 7351272  
 Dilution Factor: 1    Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	0.22 J	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28
SURROGATE		PERCENT	RECOVERY	
		RECOVERY	LIMITS	
Dibromofluoromethane	90		(73 - 122)	
1,2-Dichloroethane-d4	91		(61 - 128)	
Toluene-d8	91		(76 - 110)	
4-Bromofluorobenzene	88		(74 - 116)	

(Continued on next page)

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120507-JH-019**

**GC/MS Volatiles**

**Lot-Sample #...: A7L070423-004 Work Order #...: KDQPT1AA Matrix.....: WG**

**NOTE (S) :**

J Estimated result. Result is less than RL.

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120507-JH-019

GC Volatiles

Lot-Sample #....: A7L070423-004 Work Order #....: KDQPT1AC Matrix.....: WG  
Date Sampled....: 12/05/07 14:20 Date Received...: 12/07/07  
Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
Prep Batch #....: 7346409  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	ND	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	86	(10 - 150)		

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120507-JH-020

## GC/MS Volatiles

Lot-Sample #....: A7L070423-005   Work Order #....: KDQPV1AA      Matrix.....: WG  
 Date Sampled....: 12/05/07 14:50   Date Received...: 12/07/07  
 Prep Date.....: 12/15/07      Analysis Date...: 12/15/07  
 Prep Batch #....: 7351272  
 Dilution Factor: 1              Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY
		LIMITS
Dibromofluoromethane	90	(73 - 122)
1,2-Dichloroethane-d4	92	(61 - 128)
Toluene-d8	90	(76 - 110)
4-Bromofluorobenzene	87	(74 - 116)

Conenstoga-Rovers & Associates, Inc.

Client Sample ID: GW-120507-JH-020

GC Volatiles

Lot-Sample #....: A7L070423-005 Work Order #....: KDQPV1AC Matrix.....: WG  
Date Sampled...: 12/05/07 14:50 Date Received...: 12/07/07  
Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
Prep Batch #....: 7346409  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		MDL
		LIMIT	UNITS	
TPH (as Gasoline)	31 J	100	ug/L	28
<hr/>				
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	85	(10 - 150)		

NOTE (S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120507-JH-021

## GC/MS Volatiles

Lot-Sample #....: A7L070423-006   Work Order #....: KDQPX1AA      Matrix.....: WG  
 Date Sampled...: 12/05/07 15:05   Date Received..: 12/07/07  
 Prep Date.....: 12/15/07      Analysis Date...: 12/15/07  
 Prep Batch #....: 7351272  
 Dilution Factor: 1              Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	<b>6.4</b>	<b>0.50</b>	ug/L	<b>0.17</b>
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	<b>0.26 J</b>	<b>1.0</b>	ug/L	<b>0.13</b>
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	<b>1.0</b>	<b>1.0</b>	ug/L	<b>0.22</b>
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS	
		(73 - 122)	(61 - 128)
Dibromofluoromethane	88	(73 - 122)	
1,2-Dichloroethane-d4	91	(61 - 128)	
Toluene-d8	91	(76 - 110)	
4-Bromofluorobenzene	89	(74 - 116)	

(Continued on next page)

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120507-JH-021

GC/MS Volatiles

Lot-Sample #...: A7L070423-006 Work Order #...: KDQPX1AA Matrix.....: WG

NOTE (S) :

J Estimated result. Result is less than RL.

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120507-JH-021

GC Volatiles

Lot-Sample #....: A7L070423-006 Work Order #....: KDQPX1AC Matrix.....: WG  
Date Sampled....: 12/05/07 15:05 Date Received...: 12/07/07  
Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
Prep Batch #....: 7346409  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	31 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	83	(10 - 150)		

NOTE(S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: RW-120507-JH-022

## GC/MS Volatiles

Lot-Sample #...: A7L070423-007 Work Order #...: KDQP11AA Matrix.....: WG  
 Date Sampled...: 12/05/07 15:20 Date Received...: 12/07/07  
 Prep Date.....: 12/15/07 Analysis Date...: 12/15/07  
 Prep Batch #...: 7351272  
 Dilution Factor: 1 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	5.6 J	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY	
		LIMITS	
Dibromofluoromethane	92	(73 - 122)	
1,2-Dichloroethane-d4	96	(61 - 128)	
Toluene-d8	90	(76 - 110)	
4-Bromofluorobenzene	88	(74 - 116)	

(Continued on next page)

Conestoga-Rovers & Associates, Inc.

Client Sample ID: RW-120507-JH-022

GC/MS Volatiles

Lot-Sample #...: A7L070423-007 Work Order #...: KDQP11AA Matrix.....: WG

NOTE (S) :

J Estimated result. Result is less than RL.

Conestoga-Rovers & Associates, Inc.

Client Sample ID: RW-120507-JH-022

GC Volatiles

Lot-Sample #....: A7L070423-007 Work Order #....: KDQP11AC Matrix.....: WG  
Date Sampled....: 12/05/07 15:20 Date Received...: 12/07/07  
Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
Prep Batch #....: 7346409  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	33 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	84	(10 - 150)		

NOTE (S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120507-JH-023

## GC/MS Volatiles

Lot-Sample #....: A7L070423-008 Work Order #....: KDQP21AA Matrix.....: WG  
 Date Sampled...: 12/05/07 16:05 Date Received...: 12/07/07  
 Prep Date.....: 12/15/07 Analysis Date...: 12/15/07  
 Prep Batch #....: 7351272  
 Dilution Factor: 1 Method.....: SW846 8260B

<u>PARAMETER</u>	<u>RESULT</u>	<u>REPORTING</u>		
		<u>LIMIT</u>	<u>UNITS</u>	<u>MDL</u>
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	<b>1.9</b>	<b>0.50</b>	<b>ug/L</b>	<b>0.17</b>
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	<b>9.6</b>	<b>1.0</b>	<b>ug/L</b>	<b>0.29</b>
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	<b>1.4</b>	<b>1.0</b>	<b>ug/L</b>	<b>0.17</b>
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

<u>SURROGATE</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY</u>
		<u>LIMITS</u>
Dibromofluoromethane	89	(73 - 122)
1,2-Dichloroethane-d4	93	(61 - 128)
Toluene-d8	92	(76 - 110)
4-Bromofluorobenzene	88	(74 - 116)

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120507-JH-023

GC Volatiles

Lot-Sample #....: A7L070423-008 Work Order #....: KDQP21AC Matrix.....: WG  
Date Sampled....: 12/05/07 16:05 Date Received...: 12/07/07  
Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
Prep Batch #....: 7346409  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		MDL
		LIMIT	UNITS	
TPH (as Gasoline)	33 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	85	(10 - 150)		

NOTE (S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120507-JH-024

## GC/MS Volatiles

Lot-Sample #...: A7L070423-009   Work Order #...: KDQP41AA   Matrix.....: WG  
 Date Sampled...: 12/05/07 16:45   Date Received..: 12/07/07  
 Prep Date.....: 12/15/07   Analysis Date...: 12/15/07  
 Prep Batch #...: 7351272  
 Dilution Factor: 1   Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY	
		LIMITS	
Dibromofluoromethane	94	(73 - 122)	
1,2-Dichloroethane-d4	94	(61 - 128)	
Toluene-d8	91	(76 - 110)	
4-Bromofluorobenzene	88	(74 - 116)	

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120507-JH-024

GC Volatiles

Lot-Sample #....: A7L070423-009 Work Order #....: KDQP41AC Matrix.....: WG  
Date Sampled...: 12/05/07 16:45 Date Received...: 12/07/07  
Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
Prep Batch #....: 7346440  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		MDL
		LIMIT	UNITS	
TPH (as Gasoline)	ND	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	98	(10 - 150)		

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120607-JH-025

## GC/MS Volatiles

Lot-Sample #....: A7L070423-010 Work Order #....: KDQP61AA Matrix.....: WG  
 Date Sampled....: 12/06/07 09:15 Date Received...: 12/07/07  
 Prep Date.....: 12/15/07 Analysis Date...: 12/15/07  
 Prep Batch #....: 7351272  
 Dilution Factor: 1 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY
		LIMITS
Dibromofluoromethane	94	(73 - 122)
1,2-Dichloroethane-d4	97	(61 - 128)
Toluene-d8	92	(76 - 110)
4-Bromofluorobenzene	89	(74 - 116)

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120607-JH-025

## GC Volatiles

Lot-Sample #...: A7L070423-010 Work Order #...: KDQP61AC Matrix.....: WG  
Date Sampled...: 12/06/07 09:15 Date Received...: 12/07/07  
Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
Prep Batch #...: 7346440  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	31 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	105	(10 - 150)		

## NOTE (S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120607-JH-026

## GC/MS Volatiles

Lot-Sample #....: A7L070423-011 Work Order #....: KDQQA1AA Matrix.....: WG  
 Date Sampled....: 12/06/07 09:30 Date Received...: 12/07/07  
 Prep Date.....: 12/15/07 Analysis Date...: 12/15/07  
 Prep Batch #....: 7351272  
 Dilution Factor: 1 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28
SURROGATE		PERCENT	RECOVERY	
		RECOVERY	LIMITS	
Dibromofluoromethane		90	(73 - 122)	
1,2-Dichloroethane-d4		94	(61 - 128)	
Toluene-d8		89	(76 - 110)	
4-Bromofluorobenzene		88	(74 - 116)	

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120607-JH-026

GC Volatiles

Lot-Sample #...: A7L070423-011 Work Order #...: KDQQA1AC Matrix.....: WG  
Date Sampled...: 12/06/07 09:30 Date Received...: 12/07/07  
Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
Prep Batch #...: 7346440  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	33 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	99	(10 - 150)		

NOTE (S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120607-JH-027

## GC/MS Volatiles

Lot-Sample #....: A7L070423-012 Work Order #....: KDQQC1AA Matrix.....: WG  
 Date Sampled...: 12/06/07 10:30 Date Received...: 12/07/07  
 Prep Date.....: 12/18/07 Analysis Date...: 12/18/07  
 Prep Batch #....: 7353354  
 Dilution Factor: 1 Method.....: SW846 8260B

<u>PARAMETER</u>	<u>RESULT</u>	<u>REPORTING</u>		
		<u>LIMIT</u>	<u>UNITS</u>	<u>MDL</u>
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

<u>SURROGATE</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>
Dibromofluoromethane	94	(73 - 122)
1,2-Dichloroethane-d4	105	(61 - 128)
Toluene-d8	90	(76 - 110)
4-Bromofluorobenzene	92	(74 - 116)

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120607-JH-027**

**GC Volatiles**

**Lot-Sample #....: A7L070423-012    Work Order #....: KDQQC1AC    Matrix.....: WG**  
**Date Sampled....: 12/06/07 10:30    Date Received...: 12/07/07**  
**Prep Date.....: 12/12/07    Analysis Date...: 12/12/07**  
**Prep Batch #....: 7346440**  
**Dilution Factor: 1    Method.....: SW846 8015B**

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	ND	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	99	(10 - 150)		

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120607-JH-028

## GC/MS Volatiles

Lot-Sample #....: A7L070423-013 Work Order #....: KDQQE1AA Matrix.....: WG  
 Date Sampled....: 12/06/07 10:45 Date Received...: 12/07/07  
 Prep Date.....: 12/18/07 Analysis Date...: 12/18/07  
 Prep Batch #....: 7353354  
 Dilution Factor: 1 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28
 SURROGATE		PERCENT	RECOVERY	
		RECOVERY	LIMITS	
Dibromofluoromethane	94		(73 - 122)	
1,2-Dichloroethane-d4	105		(61 - 128)	
Toluene-d8	90		(76 - 110)	
4-Bromofluorobenzene	89		(74 - 116)	

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120607-JH-028

GC Volatiles

Lot-Sample #....: A7L070423-013 Work Order #....: KDQQE1AC Matrix.....: WG  
Date Sampled...: 12/06/07 10:45 Date Received...: 12/07/07  
Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
Prep Batch #....: 7346440  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	30 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	103	(10 - 150)		

NOTE (S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120607-JH-029

## GC/MS Volatiles

Lot-Sample #....: A7L070423-014 Work Order #....: KDQQG1AA Matrix.....: WG  
 Date Sampled...: 12/06/07 11:15 Date Received...: 12/07/07  
 Prep Date.....: 12/18/07 Analysis Date...: 12/18/07  
 Prep Batch #...: 7353354  
 Dilution Factor: 1 Method.....: SW846 8260B

<u>PARAMETER</u>	<u>RESULT</u>	REPORTING		
		<u>LIMIT</u>	<u>UNITS</u>	<u>MDL</u>
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	0.36 J	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

<u>SURROGATE</u>	<u>PERCENT RECOVERY</u>	RECOVERY
		<u>LIMITS</u>
Dibromofluoromethane	96	(73 - 122)
1,2-Dichloroethane-d4	105	(61 - 128)
Toluene-d8	89	(76 - 110)
4-Bromofluorobenzene	90	(74 - 116)

(Continued on next page)

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120607-JH-029

GC/MS Volatiles

Lot-Sample #...: A7L070423-014 Work Order #...: KDQQG1AA Matrix.....: WG

NOTE(S) :

J Estimated result. Result is less than RL.

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120607-JH-029**

**GC Volatiles**

Lot-Sample #....: A7L070423-014    Work Order #....: KDQQG1AC    Matrix.....: WG  
Date Sampled....: 12/06/07 11:15    Date Received...: 12/07/07  
Prep Date.....: 12/12/07    Analysis Date...: 12/12/07  
Prep Batch #....: 7346440  
Dilution Factor: 1    Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	ND	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	96	(10 ~ 150)		

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120607-JH-030

## GC/MS Volatiles

Lot-Sample #...: A7L070423-015 Work Order #...: KDQQH1AA Matrix.....: WG  
 Date Sampled...: 12/06/07 12:00 Date Received...: 12/07/07  
 Prep Date.....: 12/18/07 Analysis Date...: 12/18/07  
 Prep Batch #...: 7353354  
 Dilution Factor: 5000 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	50000	ug/L	5500
Benzene	ND	5000	ug/L	650
Bromodichloromethane	ND	5000	ug/L	750
Bromoform	ND	5000	ug/L	3200
Bromomethane	ND	5000	ug/L	2000
2-Butanone	ND	50000	ug/L	2800
Carbon disulfide	ND	5000	ug/L	650
Carbon tetrachloride	ND	5000	ug/L	650
Chlorobenzene	ND	5000	ug/L	750
Chloroethane	ND	5000	ug/L	1400
Chloroform	ND	5000	ug/L	800
Chloromethane	ND	5000	ug/L	1500
Dibromochloromethane	ND	5000	ug/L	900
1,1-Dichloroethane	ND	5000	ug/L	750
1,2-Dichloroethane	ND	5000	ug/L	1100
1,1-Dichloroethene	ND	5000	ug/L	950
cis-1,2-Dichloroethene	ND	2500	ug/L	850
trans-1,2-Dichloroethene	ND	2500	ug/L	950
1,2-Dichloropropane	ND	5000	ug/L	900
cis-1,3-Dichloropropene	ND	5000	ug/L	700
trans-1,3-Dichloropropene	ND	5000	ug/L	950
Ethylbenzene	8000	5000	ug/L	850
2-Hexanone	ND	50000	ug/L	2000
Methylene chloride	ND	5000	ug/L	1600
4-Methyl-2-pentanone	160000	50000	ug/L	1600
Styrene	ND	5000	ug/L	550
1,1,2,2-Tetrachloroethane	ND	5000	ug/L	900
Tetrachloroethene	ND	5000	ug/L	1400
Toluene	150000	5000	ug/L	650
1,1,1-Trichloroethane	ND	5000	ug/L	1100
1,1,2-Trichloroethane	ND	5000	ug/L	1400
Trichloroethene	ND	5000	ug/L	850
Vinyl chloride	ND	5000	ug/L	1100
Xylenes (total)	57000	5000	ug/L	1400
SURROGATE		PERCENT RECOVERY	RECOVERY LIMITS	
Dibromofluoromethane	95	(73 - 122)		
1,2-Dichloroethane-d4	106	(61 - 128)		
Toluene-d8	95	(76 - 110)		
4-Bromofluorobenzene	94	(74 - 116)		

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120607-JH-030

GC Volatiles

Lot-Sample #....: A7L070423-015 Work Order #....: KDQQH1AC Matrix.....: WG  
Date Sampled....: 12/06/07 12:00 Date Received...: 12/07/07  
Prep Date.....: 12/11/07 Analysis Date...: 12/11/07  
Prep Batch #....: 7346440  
Dilution Factor: 1000 Method.....: SW846 8015B

PARAMETER	REPORTING		
	RESULT	LIMIT	UNITS
TPH (as Gasoline)	470000	100000	ug/L
RECOVERY			Limits
SURROGATE	PERCENT	RECOVERY	(10 - 150)
Trifluorotoluene	93		

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120607-JH-031

## GC/MS Volatiles

Lot-Sample #....: A7L070423-016 Work Order #....: KDQQK1AA Matrix.....: WG  
 Date Sampled...: 12/06/07 12:30 Date Received...: 12/07/07  
 Prep Date.....: 12/18/07 Analysis Date...: 12/18/07  
 Prep Batch #....: 7353354  
 Dilution Factor: 5000 Method.....: SW846 8260B

<u>PARAMETER</u>	<u>RESULT</u>	<u>REPORTING</u>		
		<u>LIMIT</u>	<u>UNITS</u>	<u>MDL</u>
Acetone	ND	50000	ug/L	5500
Benzene	ND	5000	ug/L	650
Bromodichloromethane	ND	5000	ug/L	750
Bromoform	ND	5000	ug/L	3200
Bromomethane	ND	5000	ug/L	2000
2-Butanone	ND	50000	ug/L	2800
Carbon disulfide	ND	5000	ug/L	650
Carbon tetrachloride	ND	5000	ug/L	650
Chlorobenzene	ND	5000	ug/L	750
Chloroethane	ND	5000	ug/L	1400
Chloroform	ND	5000	ug/L	800
Chloromethane	ND	5000	ug/L	1500
Dibromochloromethane	ND	5000	ug/L	900
1,1-Dichloroethane	ND	5000	ug/L	750
1,2-Dichloroethane	ND	5000	ug/L	1100
1,1-Dichloroethene	ND	5000	ug/L	950
cis-1,2-Dichloroethene	ND	2500	ug/L	850
trans-1,2-Dichloroethene	ND	2500	ug/L	950
1,2-Dichloropropane	ND	5000	ug/L	900
cis-1,3-Dichloropropene	ND	5000	ug/L	700
trans-1,3-Dichloropropene	ND	5000	ug/L	950
Ethylbenzene	7500	5000	ug/L	850
2-Hexanone	ND	50000	ug/L	2000
Methylene chloride	ND	5000	ug/L	1600
4-Methyl-2-pentanone	160000	50000	ug/L	1600
Styrene	ND	5000	ug/L	550
1,1,2,2-Tetrachloroethane	ND	5000	ug/L	900
Tetrachloroethene	ND	5000	ug/L	1400
Toluene	140000	5000	ug/L	650
1,1,1-Trichloroethane	ND	5000	ug/L	1100
1,1,2-Trichloroethane	ND	5000	ug/L	1400
Trichloroethene	ND	5000	ug/L	850
Vinyl chloride	ND	5000	ug/L	1100
Xylenes (total)	56000	5000	ug/L	1400
<u>SURROGATE</u>		<u>PERCENT</u>	<u>RECOVERY</u>	
		<u>RECOVERY</u>	<u>LIMITS</u>	
Dibromofluoromethane	94		(73 - 122)	
1,2-Dichloroethane-d4	104		(61 - 128)	
Toluene-d8	93		(76 - 110)	
4-Bromofluorobenzene	96		(74 - 116)	

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120607-JH-031**

**GC Volatiles**

**Lot-Sample #....: A7L070423-016   Work Order #....: KDQQK1AC      Matrix.....: WG  
Date Sampled...: 12/06/07 12:30 Date Received...: 12/07/07  
Prep Date.....: 12/11/07      Analysis Date...: 12/11/07  
Prep Batch #....: 7346440  
Dilution Factor: 1000      Method.....: SW846 8015B**

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
<b>TPH (as Gasoline)</b>	<b>540000</b>	<b>100000</b>	<b>ug/L</b>	<b>28000</b>
SURROGATE	PERCENT	RECOVERY	LIMITS	
Trifluorotoluene	RECOVERY	(10 - 150)		
	99			

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120607-JH-032

## GC/MS Volatiles

Lot-Sample #....: A7L070423-017 Work Order #....: KDQQM1AA Matrix.....: WG  
 Date Sampled...: 12/06/07 13:40 Date Received..: 12/07/07  
 Prep Date.....: 12/18/07 Analysis Date...: 12/18/07  
 Prep Batch #...: 7353354  
 Dilution Factor: 1 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	0.48 J	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	0.47 J	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS
Dibromofluoromethane	97	(73 - 122)
1,2-Dichloroethane-d4	106	(61 - 128)
Toluene-d8	90	(76 - 110)
4-Bromofluorobenzene	93	(74 - 116)

(Continued on next page)

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120607-JH-032**

**GC/MS Volatiles**

**Lot-Sample #....: A7L070423-017 Work Order #....: KDQQM1AA Matrix.....: WG**

**NOTE (S) :**

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J Estimated result. Result is less than RL.

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120607-JH-032

GC Volatiles

Lot-Sample #...: A7L070423-017 Work Order #...: KDQQM1AC Matrix.....: WG  
Date Sampled...: 12/06/07 13:40 Date Received...: 12/07/07  
Prep Date.....: 12/11/07 Analysis Date...: 12/11/07  
Prep Batch #...: 7346440  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	50 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	94	(10 - 150)		

NOTE (S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120607-JH-033

## GC/MS Volatiles

Lot-Sample #....: A7L070423-018    Work Order #....: KDQQN1AA    Matrix.....: WG  
 Date Sampled....: 12/06/07 14:20    Date Received...: 12/07/07  
 Prep Date.....: 12/18/07    Analysis Date...: 12/18/07  
 Prep Batch #....: 7353354  
 Dilution Factor: 1    Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	20	0.50	ug/L	0.17
trans-1,2-Dichloroethene	1.5	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	0.29 J	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY	
		LIMITS	
Dibromofluoromethane	93	(73 - 122)	
1,2-Dichloroethane-d4	102	(61 - 128)	
Toluene-d8	88	(76 - 110)	
4-Bromofluorobenzene	89	(74 - 116)	

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**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120607-JH-033**

**GC/MS Volatiles**

**Lot-Sample #....: A7L070423-018 Work Order #....: KDQQN1AA Matrix.....: WG**

**NOTE (S) :**

J Estimated result. Result is less than RL.

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120607-JH-033**

**GC Volatiles**

**Lot-Sample #....:** A7L070423-018    **Work Order #....:** KDQQN1AC    **Matrix.....:** WG  
**Date Sampled....:** 12/06/07 14:20    **Date Received...:** 12/07/07  
**Prep Date.....:** 12/11/07    **Analysis Date...:** 12/11/07  
**Prep Batch #....:** 7346440  
**Dilution Factor:** 1    **Method.....:** SW846 8015B

<b>PARAMETER</b>	<b>RESULT</b>	<b>REPORTING</b>		
		<b>LIMIT</b>	<b>UNITS</b>	<b>MDL</b>
<b>TPH (as Gasoline)</b>	56 J	100	ug/L	28
<b>SURROGATE</b>	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	95	(10 - 150)		

**NOTE (S) :**

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: RW-120607-JH-034

## GC/MS Volatiles

Lot-Sample #...: A7L070423-019 Work Order #...: KDQQP1AA Matrix.....: WG  
 Date Sampled...: 12/06/07 14:30 Date Received...: 12/07/07  
 Prep Date.....: 12/18/07 Analysis Date...: 12/18/07  
 Prep Batch #...: 7353354  
 Dilution Factor: 1 Method.....: SW846 8260B

<u>PARAMETER</u>	<u>RESULT</u>	<u>REPORTING LIMIT</u>	<u>UNITS</u>	<u>MDL</u>
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
<b>Chloroform</b>	<b>0.55 J</b>	<b>1.0</b>	<b>ug/L</b>	<b>0.16</b>
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

<u>SURROGATE</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>
Dibromofluoromethane	95	(73 - 122)
1,2-Dichloroethane-d4	107	(61 - 128)
Toluene-d8	90	(76 - 110)
4-Bromofluorobenzene	89	(74 - 116)

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**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: RW-120607-JH-034**

**GC/MS Volatiles**

**Lot-Sample #....: A7L070423-019 Work Order #....: KDQQP1AA Matrix.....: WG**

**NOTE (S) :**

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J Estimated result. Result is less than RL.

Conestoga-Rovers & Associates, Inc.

Client Sample ID: RW-120607-JH-034

GC Volatiles

Lot-Sample #....: A7L070423-019 Work Order #....: KDQQP1AC Matrix.....: WG  
Date Sampled...: 12/06/07 14:30 Date Received...: 12/07/07  
Prep Date.....: 12/11/07 Analysis Date...: 12/11/07  
Prep Batch #...: 7346440  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	45 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	96	(10 - 150)		

NOTE(S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120607-JH-035

## GC/MS Volatiles

Lot-Sample #....: A7L070423-020 Work Order #....: KDQQR1AA Matrix.....: WG  
 Date Sampled...: 12/06/07 15:10 Date Received...: 12/07/07  
 Prep Date.....: 12/18/07 Analysis Date...: 12/18/07  
 Prep Batch #....: 7353354  
 Dilution Factor: 3.33 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	33	ug/L	3.7
Benzene	ND	3.3	ug/L	0.43
Bromodichloromethane	ND	3.3	ug/L	0.50
Bromoform	ND	3.3	ug/L	2.1
Bromomethane	ND	3.3	ug/L	1.4
2-Butanone	ND	33	ug/L	1.9
Carbon disulfide	ND	3.3	ug/L	0.43
Carbon tetrachloride	ND	3.3	ug/L	0.43
Chlorobenzene	ND	3.3	ug/L	0.50
Chloroethane	ND	3.3	ug/L	0.97
Chloroform	ND	3.3	ug/L	0.53
Chloromethane	ND	3.3	ug/L	1.0
Dibromochloromethane	ND	3.3	ug/L	0.60
1,1-Dichloroethane	ND	3.3	ug/L	0.50
1,2-Dichloroethane	ND	3.3	ug/L	0.73
1,1-Dichloroethene	ND	3.3	ug/L	0.63
cis-1,2-Dichloroethene	88	1.7	ug/L	0.57
trans-1,2-Dichloroethene	6.2	1.7	ug/L	0.63
1,2-Dichloropropane	ND	3.3	ug/L	0.60
cis-1,3-Dichloropropene	ND	3.3	ug/L	0.47
trans-1,3-Dichloropropene	ND	3.3	ug/L	0.63
Ethylbenzene	ND	3.3	ug/L	0.57
2-Hexanone	ND	33	ug/L	1.4
Methylene chloride	1.7 J	3.3	ug/L	1.1
4-Methyl-2-pentanone	ND	33	ug/L	1.1
Styrene	ND	3.3	ug/L	0.37
1,1,2,2-Tetrachloroethane	ND	3.3	ug/L	0.60
Tetrachloroethene	ND	3.3	ug/L	0.97
Toluene	ND	3.3	ug/L	0.43
1,1,1-Trichloroethane	ND	3.3	ug/L	0.73
1,1,2-Trichloroethane	ND	3.3	ug/L	0.90
Trichloroethene	ND	3.3	ug/L	0.57
Vinyl chloride	7.0	3.3	ug/L	0.73
Xylenes (total)	ND	3.3	ug/L	0.93

SURROGATE	PERCENT RECOVERY	RECOVERY
		LIMITS
Dibromofluoromethane	93	(73 - 122)
1,2-Dichloroethane-d4	103	(61 - 128)
Toluene-d8	94	(76 - 110)
4-Bromofluorobenzene	87	(74 - 116)

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Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120607-JH-035

GC/MS Volatiles

Lot-Sample #...: A7L070423-020 Work Order #...: KDQQR1AA Matrix.....: WG

NOTE (S) :

J Estimated result. Result is less than RL.

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120607-JH-035

GC Volatiles

Lot-Sample #....: A7L070423-020 Work Order #....: KDQQR1AC Matrix.....: WG  
Date Sampled...: 12/06/07 15:10 Date Received...: 12/07/07  
Prep Date.....: 12/11/07 Analysis Date...: 12/11/07  
Prep Batch #...: 7346440  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	130	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	98	(10 - 150)		

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120607-JH-036

## GC/MS Volatiles

Lot-Sample #....: A7L070423-021 Work Order #....: KDQQT1AA Matrix.....: WG  
 Date Sampled...: 12/06/07 15:30 Date Received..: 12/07/07  
 Prep Date.....: 12/18/07 Analysis Date...: 12/18/07  
 Prep Batch #....: 7353354  
 Dilution Factor: 3.33 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	33	ug/L	3.7
Benzene	ND	3.3	ug/L	0.43
Bromodichloromethane	ND	3.3	ug/L	0.50
Bromoform	ND	3.3	ug/L	2.1
Bromomethane	ND	3.3	ug/L	1.4
2-Butanone	ND	33	ug/L	1.9
Carbon disulfide	ND	3.3	ug/L	0.43
Carbon tetrachloride	ND	3.3	ug/L	0.43
Chlorobenzene	ND	3.3	ug/L	0.50
Chloroethane	ND	3.3	ug/L	0.97
Chloroform	ND	3.3	ug/L	0.53
Chloromethane	ND	3.3	ug/L	1.0
Dibromochloromethane	ND	3.3	ug/L	0.60
1,1-Dichloroethane	ND	3.3	ug/L	0.50
1,2-Dichloroethane	ND	3.3	ug/L	0.73
1,1-Dichloroethene	ND	3.3	ug/L	0.63
cis-1,2-Dichloroethene	89	1.7	ug/L	0.57
trans-1,2-Dichloroethene	6.0	1.7	ug/L	0.63
1,2-Dichloropropane	ND	3.3	ug/L	0.60
cis-1,3-Dichloropropene	ND	3.3	ug/L	0.47
trans-1,3-Dichloropropene	ND	3.3	ug/L	0.63
Ethylbenzene	ND	3.3	ug/L	0.57
2-Hexanone	ND	33	ug/L	1.4
Methylene chloride	2.8 J	3.3	ug/L	1.1
4-Methyl-2-pentanone	ND	33	ug/L	1.1
Styrene	ND	3.3	ug/L	0.37
1,1,2,2-Tetrachloroethane	ND	3.3	ug/L	0.60
Tetrachloroethene	ND	3.3	ug/L	0.97
Toluene	ND	3.3	ug/L	0.43
1,1,1-Trichloroethane	ND	3.3	ug/L	0.73
1,1,2-Trichloroethane	ND	3.3	ug/L	0.90
Trichloroethene	ND	3.3	ug/L	0.57
Vinyl chloride	6.7	3.3	ug/L	0.73
Xylenes (total)	ND	3.3	ug/L	0.93

SURROGATE	PERCENT RECOVERY	RECOVERY
		LIMITS
Dibromofluoromethane	95	(73 - 122)
1,2-Dichloroethane-d4	103	(61 - 128)
Toluene-d8	93	(76 - 110)
4-Bromofluorobenzene	92	(74 - 116)

(Continued on next page)

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120607-JH-036

GC/MS Volatiles

Lot-Sample #....: A7L070423-021 Work Order #....: KDQQT1AA Matrix.....: WG

NOTE(S) :

J Estimated result. Result is less than RL.

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120607-JH-036**

**GC Volatiles**

Lot-Sample #...: A7L070423-021 Work Order #...: KDQQT1AC Matrix.....: WG  
Date Sampled...: 12/06/07 15:30 Date Received...: 12/07/07  
Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
Prep Batch #...: 7346440  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		MDL
		LIMIT	UNITS	
TPH (as Gasoline)	120	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	106	(10 - 150)		

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120607-JH-037

## GC/MS Volatiles

Lot-Sample #....: A7L070423-022 Work Order #....: KDQQW1AA Matrix.....: WG  
 Date Sampled....: 12/06/07 15:55 Date Received...: 12/07/07  
 Prep Date.....: 12/18/07 Analysis Date...: 12/18/07  
 Prep Batch #....: 7353354  
 Dilution Factor: 1 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	0.33 J	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS	
		(	)
Dibromofluoromethane	94	(73	- 122)
1,2-Dichloroethane-d4	106	(61	- 128)
Toluene-d8	93	(76	- 110)
4-Bromofluorobenzene	89	(74	- 116)

(Continued on next page)

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120607-JH-037

GC/MS Volatiles

Lot-Sample #...: A7L070423-022 Work Order #...: KDQQW1AA Matrix.....: WG

NOTE(S) :

J Estimated result. Result is less than RL.

Conestoga-Rovers & Associates, Inc.

Client Sample ID: GW-120607-JH-037

GC Volatiles

Lot-Sample #...: A7L070423-022 Work Order #...: KDQQW1AC Matrix.....: WG  
Date Sampled...: 12/06/07 15:55 Date Received..: 12/07/07  
Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
Prep Batch #...: 7346440  
Dilution Factor: 1 Method.....: SW846 8015B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	29 J	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	99	(10 - 150)		

NOTE (S) :

J Estimated result. Result is less than RL.

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: GW-120607-JH-038

## GC/MS Volatiles

Lot-Sample #....: A7L070423-023 Work Order #....: KDQQ01AA Matrix.....: WG  
 Date Sampled....: 12/06/07 16:30 Date Received...: 12/07/07  
 Prep Date.....: 12/18/07 Analysis Date...: 12/18/07  
 Prep Batch #....: 7353354  
 Dilution Factor: 1 Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING LIMIT	UNITS	MDL
Acetone	ND	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	0.35 J	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS
Dibromofluoromethane	94	(73 - 122)
1,2-Dichloroethane-d4	106	(61 - 128)
Toluene-d8	91	(76 - 110)
4-Bromofluorobenzene	91	(74 - 116)

(Continued on next page)

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120607-JH-038**

**GC/MS Volatiles**

**Lot-Sample #...: A7L070423-023 Work Order #...: KDQQ01AA Matrix.....: WG**

**NOTE (S) :**

J Estimated result. Result is less than RL.

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: GW-120607-JH-038**

**GC Volatiles**

**Lot-Sample #....: A7L070423-023 Work Order #....: KDQQ01AC Matrix.....: WG  
Date Sampled....: 12/06/07 16:30 Date Received...: 12/07/07  
Prep Date.....: 12/12/07 Analysis Date...: 12/12/07  
Prep Batch #....: 7346440  
Dilution Factor: 1 Method.....: SW846 8015B**

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
TPH (as Gasoline)	ND	100	ug/L	28
SURROGATE	PERCENT	RECOVERY		
Trifluorotoluene	RECOVERY	LIMITS		
	100	(10 - 150)		

## Conestoga-Rovers &amp; Associates, Inc.

Client Sample ID: TRIP BLANK B COOLER 68187

## GC/MS Volatiles

Lot-Sample #...: A7L070423-024    Work Order #...: KDQQ21AA    Matrix.....: WG  
 Date Sampled...: 12/06/07    Date Received...: 12/07/07  
 Prep Date.....: 12/18/07    Analysis Date...: 12/18/07  
 Prep Batch #...: 7353354  
 Dilution Factor: 1            Method.....: SW846 8260B

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	MDL
Acetone	1.9 J	10	ug/L	1.1
Benzene	ND	1.0	ug/L	0.13
Bromodichloromethane	ND	1.0	ug/L	0.15
Bromoform	ND	1.0	ug/L	0.64
Bromomethane	ND	1.0	ug/L	0.41
2-Butanone	ND	10	ug/L	0.57
Carbon disulfide	ND	1.0	ug/L	0.13
Carbon tetrachloride	ND	1.0	ug/L	0.13
Chlorobenzene	ND	1.0	ug/L	0.15
Chloroethane	ND	1.0	ug/L	0.29
Chloroform	ND	1.0	ug/L	0.16
Chloromethane	ND	1.0	ug/L	0.30
Dibromochloromethane	ND	1.0	ug/L	0.18
1,1-Dichloroethane	ND	1.0	ug/L	0.15
1,2-Dichloroethane	ND	1.0	ug/L	0.22
1,1-Dichloroethene	ND	1.0	ug/L	0.19
cis-1,2-Dichloroethene	ND	0.50	ug/L	0.17
trans-1,2-Dichloroethene	ND	0.50	ug/L	0.19
1,2-Dichloropropane	ND	1.0	ug/L	0.18
cis-1,3-Dichloropropene	ND	1.0	ug/L	0.14
trans-1,3-Dichloropropene	ND	1.0	ug/L	0.19
Ethylbenzene	ND	1.0	ug/L	0.17
2-Hexanone	ND	10	ug/L	0.41
Methylene chloride	ND	1.0	ug/L	0.33
4-Methyl-2-pentanone	ND	10	ug/L	0.32
Styrene	ND	1.0	ug/L	0.11
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	0.18
Tetrachloroethene	ND	1.0	ug/L	0.29
Toluene	ND	1.0	ug/L	0.13
1,1,1-Trichloroethane	ND	1.0	ug/L	0.22
1,1,2-Trichloroethane	ND	1.0	ug/L	0.27
Trichloroethene	ND	1.0	ug/L	0.17
Vinyl chloride	ND	1.0	ug/L	0.22
Xylenes (total)	ND	1.0	ug/L	0.28
SURROGATE	PERCENT RECOVERY	RECOVERY		
		LIMITS		
Dibromofluoromethane	94	(73 - 122)		
1,2-Dichloroethane-d4	106	(61 - 128)		
Toluene-d8	89	(76 - 110)		
4-Bromofluorobenzene	91	(74 - 116)		

(Continued on next page)

Conestoga-Rovers & Associates, Inc.

Client Sample ID: TRIP BLANK B COOLER 68187

GC/MS Volatiles

Lot-Sample #...: A7L070423-024 Work Order #...: KDQQ21AA Matrix.....: WG

NOTE (S) :

J Estimated result. Result is less than RL.

**Conestoga-Rovers & Associates, Inc.**

**Client Sample ID: TRIP BLANK C COOLER 68182**

**GC Volatiles**

**Lot-Sample #....: A7L070423-025    Work Order #....: KDQRR1AA    Matrix.....: WG**  
**Date Sampled....: 12/06/07    Date Received...: 12/07/07**  
**Prep Date.....: 12/12/07    Analysis Date...: 12/12/07**  
**Prep Batch #....: 7346440**  
**Dilution Factor: 1              Method.....: SW846 8015B**

<b>PARAMETER</b>	<b>RESULT</b>	<b>REPORTING</b>		
		<b>LIMIT</b>	<b>UNITS</b>	<b>MDL</b>
TPH (as Gasoline)	ND	100	ug/L	28
SURROGATE	PERCENT	RECOVERY	LIMITS	
Trifluorotoluene	RECOVERY 99	(10 - 150)		

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***QUALITY CONTROL  
SECTION***

**METHOD BLANK REPORT**

**GC/MS Volatiles**

**Client Lot #...:** 7L06288  
**MB Lot-Sample #:** A7L120000-476

**Work Order #...:** KD1PG1AA

**Matrix.....:** WATER

**Analysis Date..:** 12/11/07  
**Dilution Factor:** 1

**Prep Date.....:** 12/11/07  
**Prep Batch #...:** 7346476

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	METHOD
Acetone	ND	10	ug/L	SW846 8260B
Benzene	ND	1.0	ug/L	SW846 8260B
Bromodichloromethane	ND	1.0	ug/L	SW846 8260B
Bromoform	ND	1.0	ug/L	SW846 8260B
Bromomethane	ND	1.0	ug/L	SW846 8260B
2-Butanone	ND	10	ug/L	SW846 8260B
Carbon disulfide	ND	1.0	ug/L	SW846 8260B
Carbon tetrachloride	ND	1.0	ug/L	SW846 8260B
Chlorobenzene	ND	1.0	ug/L	SW846 8260B
Chloroethane	ND	1.0	ug/L	SW846 8260B
Chloroform	ND	1.0	ug/L	SW846 8260B
Chloromethane	ND	1.0	ug/L	SW846 8260B
Dibromochloromethane	ND	1.0	ug/L	SW846 8260B
1,1-Dichloroethane	ND	1.0	ug/L	SW846 8260B
1,2-Dichloroethane	ND	1.0	ug/L	SW846 8260B
1,1-Dichloroethene	ND	1.0	ug/L	SW846 8260B
cis-1,2-Dichloroethene	ND	0.50	ug/L	SW846 8260B
trans-1,2-Dichloroethene	ND	0.50	ug/L	SW846 8260B
1,2-Dichloropropane	ND	1.0	ug/L	SW846 8260B
cis-1,3-Dichloropropene	ND	1.0	ug/L	SW846 8260B
trans-1,3-Dichloropropene	ND	1.0	ug/L	SW846 8260B
Ethylbenzene	ND	1.0	ug/L	SW846 8260B
2-Hexanone	ND	10	ug/L	SW846 8260B
<b>Methylene chloride</b>	<b>0.58 J</b>	<b>1.0</b>	<b>ug/L</b>	<b>SW846 8260B</b>
4-Methyl-2-pentanone	ND	10	ug/L	SW846 8260B
Styrene	ND	1.0	ug/L	SW846 8260B
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	SW846 8260B
Tetrachloroethene	ND	1.0	ug/L	SW846 8260B
Toluene	ND	1.0	ug/L	SW846 8260B
1,1,1-Trichloroethane	ND	1.0	ug/L	SW846 8260B
1,1,2-Trichloroethane	ND	1.0	ug/L	SW846 8260B
Trichloroethene	ND	1.0	ug/L	SW846 8260B
Vinyl chloride	ND	1.0	ug/L	SW846 8260B
Xylenes (total)	ND	1.0	ug/L	SW846 8260B

SURROGATE	RECOVERY	PERCENT	RECOVERY
		LIMITS	LIMITS
Dibromofluoromethane	96	(73 - 122)	
1,2-Dichloroethane-d4	96	(61 - 128)	
Toluene-d8	101	(76 - 110)	
4-Bromofluorobenzene	96	(74 - 116)	

(Continued on next page)

METHOD BLANK REPORT

GC/MS Volatiles

Client Lot #....: 7L06288

Work Order #....: KD1PG1AA

Matrix.....: WATER

NOTE(S) :

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

J Estimated result. Result is less than RL.

**METHOD BLANK REPORT**

**GC/MS Volatiles**

**Client Lot #....:** 7L06288  
**MB Lot-Sample #:** A7L170000-272

**Work Order #....:** KEAMW1AA

**Matrix.....:** WATER

**Analysis Date..:** 12/15/07  
**Dilution Factor:** 1

**Prep Date.....:** 12/15/07  
**Prep Batch #....:** 7351272

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	METHOD
Acetone	ND	10	ug/L	SW846 8260B
Benzene	ND	1.0	ug/L	SW846 8260B
Bromodichloromethane	ND	1.0	ug/L	SW846 8260B
Bromoform	ND	1.0	ug/L	SW846 8260B
Bromomethane	ND	1.0	ug/L	SW846 8260B
2-Butanone	ND	10	ug/L	SW846 8260B
Carbon disulfide	ND	1.0	ug/L	SW846 8260B
Carbon tetrachloride	ND	1.0	ug/L	SW846 8260B
Chlorobenzene	ND	1.0	ug/L	SW846 8260B
Chloroethane	ND	1.0	ug/L	SW846 8260B
Chloroform	ND	1.0	ug/L	SW846 8260B
Chloromethane	ND	1.0	ug/L	SW846 8260B
Dibromochloromethane	ND	1.0	ug/L	SW846 8260B
1,1-Dichloroethane	ND	1.0	ug/L	SW846 8260B
1,2-Dichloroethane	ND	1.0	ug/L	SW846 8260B
1,1-Dichloroethene	ND	1.0	ug/L	SW846 8260B
cis-1,2-Dichloroethene	ND	0.50	ug/L	SW846 8260B
trans-1,2-Dichloroethene	ND	0.50	ug/L	SW846 8260B
1,2-Dichloropropane	ND	1.0	ug/L	SW846 8260B
cis-1,3-Dichloropropene	ND	1.0	ug/L	SW846 8260B
trans-1,3-Dichloropropene	ND	1.0	ug/L	SW846 8260B
Ethylbenzene	ND	1.0	ug/L	SW846 8260B
2-Hexanone	ND	10	ug/L	SW846 8260B
Methylene chloride	ND	1.0	ug/L	SW846 8260B
4-Methyl-2-pentanone	ND	10	ug/L	SW846 8260B
Styrene	ND	1.0	ug/L	SW846 8260B
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	SW846 8260B
Tetrachloroethene	ND	1.0	ug/L	SW846 8260B
Toluene	ND	1.0	ug/L	SW846 8260B
1,1,1-Trichloroethane	ND	1.0	ug/L	SW846 8260B
1,1,2-Trichloroethane	ND	1.0	ug/L	SW846 8260B
Trichloroethene	ND	1.0	ug/L	SW846 8260B
Vinyl chloride	ND	1.0	ug/L	SW846 8260B
Xylenes (total)	ND	1.0	ug/L	SW846 8260B

SURROGATE	RECOVERY	PERCENT	RECOVERY
		LIMITS	LIMITS
Dibromofluoromethane	90	(73 - 122)	
1,2-Dichloroethane-d4	91	(61 - 128)	
Toluene-d8	91	(76 - 110)	
4-Bromofluorobenzene	88	(74 - 116)	

(Continued on next page)

METHOD BLANK REPORT

GC/MS Volatiles

Client Lot #....: 7L06288

Work Order #....: KEAMW1AA

Matrix.....: WATER

NOTE (S) :

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

**METHOD BLANK REPORT**

**GC/MS Volatiles**

**Client Lot #...:** 7L06288  
**MB Lot-Sample #:** A7L190000-354

**Work Order #...:** KEFNV1AA      **Matrix.....:** WATER  
**Prep Date.....:** 12/18/07  
**Prep Batch #...:** 7353354

**Analysis Date..:** 12/18/07  
**Dilution Factor:** 1

PARAMETER	REPORTING			
	RESULT	LIMIT	UNITS	METHOD
Acetone	ND	10	ug/L	SW846 8260B
Benzene	ND	1.0	ug/L	SW846 8260B
Bromodichloromethane	ND	1.0	ug/L	SW846 8260B
Bromoform	ND	1.0	ug/L	SW846 8260B
Bromomethane	ND	1.0	ug/L	SW846 8260B
2-Butanone	ND	10	ug/L	SW846 8260B
Carbon disulfide	ND	1.0	ug/L	SW846 8260B
Carbon tetrachloride	ND	1.0	ug/L	SW846 8260B
Chlorobenzene	ND	1.0	ug/L	SW846 8260B
Chloroethane	ND	1.0	ug/L	SW846 8260B
Chloroform	ND	1.0	ug/L	SW846 8260B
Chloromethane	ND	1.0	ug/L	SW846 8260B
Dibromochloromethane	ND	1.0	ug/L	SW846 8260B
1,1-Dichloroethane	ND	1.0	ug/L	SW846 8260B
1,2-Dichloroethane	ND	1.0	ug/L	SW846 8260B
1,1-Dichloroethene	ND	1.0	ug/L	SW846 8260B
cis-1,2-Dichloroethene	ND	0.50	ug/L	SW846 8260B
trans-1,2-Dichloroethene	ND	0.50	ug/L	SW846 8260B
1,2-Dichloropropane	ND	1.0	ug/L	SW846 8260B
cis-1,3-Dichloropropene	ND	1.0	ug/L	SW846 8260B
trans-1,3-Dichloropropene	ND	1.0	ug/L	SW846 8260B
Ethylbenzene	ND	1.0	ug/L	SW846 8260B
2-Hexanone	ND	10	ug/L	SW846 8260B
Methylene chloride	ND	1.0	ug/L	SW846 8260B
4-Methyl-2-pentanone	ND	10	ug/L	SW846 8260B
Styrene	ND	1.0	ug/L	SW846 8260B
1,1,2,2-Tetrachloroethane	ND	1.0	ug/L	SW846 8260B
Tetrachloroethene	ND	1.0	ug/L	SW846 8260B
Toluene	ND	1.0	ug/L	SW846 8260B
1,1,1-Trichloroethane	ND	1.0	ug/L	SW846 8260B
1,1,2-Trichloroethane	ND	1.0	ug/L	SW846 8260B
Trichloroethene	ND	1.0	ug/L	SW846 8260B
Vinyl chloride	ND	1.0	ug/L	SW846 8260B
Xylenes (total)	ND	1.0	ug/L	SW846 8260B

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS	
Dibromofluoromethane	94	(73	- 122)
1,2-Dichloroethane-d4	102	(61	- 128)
Toluene-d8	92	(76	- 110)
4-Bromofluorobenzene	94	(74	- 116)

(Continued on next page)

METHOD BLANK REPORT

GC/MS Volatiles

Client Lot #...: 7L06288

Work Order #...: KEFNV1AA

Matrix.....: WATER

NOTE (S) :

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

METHOD BLANK REPORT

GC Volatiles

Client Lot #...: 7L06288  
MB Lot-Sample #: A7L100000-082

Analysis Date..: 12/07/07  
Dilution Factor: 1

Work Order #...: KDR2K1AA

Matrix.....: WATER

Prep Date.....: 12/07/07  
Prep Batch #: 7344082

PARAMETER	REPORTING		
	RESULT	LIMIT	UNITS
TPH (as Gasoline)	ND	100	ug/L
<hr/>			
SURROGATE	PERCENT	RECOVERY	
Trifluorotoluene	RECOVERY	LIMITS	
	95	(10 - 150)	

NOTE (S) :

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

METHOD BLANK REPORT

GC Volatiles

Client Lot #...: 7L06288      Work Order #...: KD09G1AA      Matrix.....: WATER  
MB Lot-Sample #: A7L120000-409  
Analysis Date..: 12/11/07      Prep Date.....: 12/11/07  
Dilution Factor: 1      Prep Batch #: 7346409

PARAMETER	RESULT	REPORTING		METHOD
		LIMIT	UNITS	
TPH (as Gasoline)	ND	100	ug/L	SW846 8015B
SURROGATE	PERCENT	RECOVERY	LIMITS	
Trifluorotoluene	83		(10 - 150)	

NOTE (S) :

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

METHOD BLANK REPORT

GC Volatiles

Client Lot #...: 7L06288  
MB Lot-Sample #: A7L120000-440

Work Order #...: KD1HK1AA Matrix.....: WATER

Analysis Date..: 12/11/07  
Dilution Factor: 1

Prep Date.....: 12/11/07  
Prep Batch #...: 7346440

PARAMETER	RESULT	REPORTING		
		LIMIT	UNITS	METHOD
TPH (as Gasoline)	ND	100	ug/L	SW846 8015B
<u>SURROGATE</u>		<u>PERCENT</u> <u>RECOVERY</u>		
Trifluorotoluene		RECOVERY	LIMITS	
		97	(10 - 150)	

NOTE(S) :

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

## LABORATORY CONTROL SAMPLE EVALUATION REPORT

## GC/MS Volatiles

Client Lot #...: 7L06288      Work Order #...: KD1PG1AC-LCS      Matrix.....: WATER  
 LCS Lot-Sample#: A7L120000-476      KD1PG1AD-LCSD  
 Prep Date.....: 12/11/07      Analysis Date...: 12/11/07  
 Prep Batch #...: 7346476  
 Dilution Factor: 1

<u>PARAMETER</u>	<u>PERCENT</u> <u>RECOVERY</u>	<u>RECOVERY</u> <u>LIMITS</u>	<u>RPD</u>	<u>LIMITS</u>	<u>METHOD</u>
Benzene	95 98	(80 - 116) (80 - 116)	3.2	(0-20)	SW846 8260B SW846 8260B
Chlorobenzene	94 101	(76 - 117) (76 - 117)	7.6	(0-20)	SW846 8260B SW846 8260B
1,1-Dichloroethene	85 91	(63 - 130) (63 - 130)	7.0	(0-20)	SW846 8260B SW846 8260B
Toluene	104 108	(74 - 119) (74 - 119)	3.6	(0-20)	SW846 8260B SW846 8260B
Trichloroethene	101 103	(75 - 122) (75 - 122)	1.7	(0-20)	SW846 8260B SW846 8260B

<u>SURROGATE</u>	<u>PERCENT</u> <u>RECOVERY</u>	<u>RECOVERY</u> <u>LIMITS</u>
Dibromofluoromethane	94 93	(73 - 122) (73 - 122)
1,2-Dichloroethane-d4	94 97	(61 - 128) (61 - 128)
Toluene-d8	102 104	(76 - 110) (76 - 110)
4-Bromofluorobenzene	102 101	(74 - 116) (74 - 116)

NOTE (S) :

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

Bold print denotes control parameters

## LABORATORY CONTROL SAMPLE DATA REPORT

## GC/MS Volatiles

Client Lot #....: 7L06288      Work Order #....: KD1PG1AC-LCS      Matrix.....: WATER  
 LCS Lot-Sample#: A7L120000-476      KD1PG1AD-LCSD  
 Prep Date.....: 12/11/07      Analysis Date...: 12/11/07  
 Prep Batch #....: 7346476  
 Dilution Factor: 1

PARAMETER	SPIKE	MEASURED		PERCENT		METHOD
	AMOUNT	AMOUNT	UNITS	RECOVERY	RPD	
Benzene	10	9.5	ug/L	95		SW846 8260B
	10	9.8	ug/L	98	3.2	SW846 8260B
Chlorobenzene	10	9.4	ug/L	94		SW846 8260B
	10	10	ug/L	101	7.6	SW846 8260B
1,1-Dichloroethene	10	8.5	ug/L	85		SW846 8260B
	10	9.1	ug/L	91	7.0	SW846 8260B
Toluene	10	10	ug/L	104		SW846 8260B
	10	11	ug/L	108	3.6	SW846 8260B
Trichloroethene	10	10	ug/L	101		SW846 8260B
	10	10	ug/L	103	1.7	SW846 8260B
<hr/>		<hr/>		PERCENT	RECOVERY	
<hr/>		<hr/>		RECOVERY	LIMITS	
SURROGATE						
Dibromofluoromethane				94	(73 - 122)	
				93	(73 - 122)	
1,2-Dichloroethane-d4				94	(61 - 128)	
				97	(61 - 128)	
Toluene-d8				102	(76 - 110)	
				104	(76 - 110)	
4-Bromofluorobenzene				102	(74 - 116)	
				101	(74 - 116)	

NOTE (S) :

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

Bold print denotes control parameters

## LABORATORY CONTROL SAMPLE EVALUATION REPORT

## GC/MS Volatiles

**Client Lot #....:** 7L06288      **Work Order #....:** KEAMW1AC-LCS      **Matrix.....:** WATER  
**LCS Lot-Sample#:** A7L170000-272      KEAMW1AD-LCSD  
**Prep Date.....:** 12/15/07      **Analysis Date...:** 12/15/07  
**Prep Batch #....:** 7351272  
**Dilution Factor:** 1

<u>PARAMETER</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>	<u>RPD</u>	<u>LIMITS</u>	<u>METHOD</u>
Benzene	90	(80 - 116)			SW846 8260B
	92	(80 - 116)	2.5	(0-20)	SW846 8260B
Chlorobenzene	93	(76 - 117)			SW846 8260B
	94	(76 - 117)	1.6	(0-20)	SW846 8260B
1,1-Dichloroethene	86	(63 - 130)			SW846 8260B
	89	(63 - 130)	3.4	(0-20)	SW846 8260B
Toluene	94	(74 - 119)			SW846 8260B
	95	(74 - 119)	0.64	(0-20)	SW846 8260B
Trichloroethene	90	(75 - 122)			SW846 8260B
	90	(75 - 122)	0.14	(0-20)	SW846 8260B

<u>SURROGATE</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>
Dibromofluoromethane	89	(73 - 122)
	90	(73 - 122)
1,2-Dichloroethane-d4	93	(61 - 128)
	89	(61 - 128)
Toluene-d8	94	(76 - 110)
	94	(76 - 110)
4-Bromofluorobenzene	92	(74 - 116)
	93	(74 - 116)

NOTE (S) :

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

Bold print denotes control parameters

## LABORATORY CONTROL SAMPLE DATA REPORT

## GC/MS Volatiles

Client Lot #...: 7L06288      Work Order #...: KEAMW1AC-LCS      Matrix.....: WATER  
 LCS Lot-Sample#: A7L170000-272      KEAMW1AD-LCSD  
 Prep Date.....: 12/15/07      Analysis Date...: 12/15/07  
 Prep Batch #...: 7351272  
 Dilution Factor: 1

PARAMETER	SPIKE	MEASURED	PERCENT	RPD	METHOD
	AMOUNT	AMOUNT	RECOVERY		
Benzene	10	9.0	ug/L	90	SW846 8260B
	10	9.2	ug/L	92	SW846 8260B
Chlorobenzene	10	9.3	ug/L	93	SW846 8260B
	10	9.4	ug/L	94	SW846 8260B
1,1-Dichloroethene	10	8.6	ug/L	86	SW846 8260B
	10	8.9	ug/L	89	SW846 8260B
Toluene	10	9.4	ug/L	94	SW846 8260B
	10	9.5	ug/L	95	SW846 8260B
Trichloroethene	10	9.0	ug/L	90	SW846 8260B
	10	9.0	ug/L	90	SW846 8260B

SURROGATE	PERCENT	RECOVERY
	RECOVERY	LIMITS
Dibromofluoromethane	89	(73 - 122)
	90	(73 - 122)
1,2-Dichloroethane-d4	93	(61 - 128)
	89	(61 - 128)
Toluene-d8	94	(76 - 110)
	94	(76 - 110)
4-Bromofluorobenzene	92	(74 - 116)
	93	(74 - 116)

NOTE (S) :

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

Bold print denotes control parameters

## LABORATORY CONTROL SAMPLE EVALUATION REPORT

## GC/MS Volatiles

Client Lot #....: 7L06288      Work Order #....: KEFNV1AC-LCS      Matrix.....: WATER  
 LCS Lot-Sample#: A7L190000-354      KEFNV1AD-LCSD  
 Prep Date.....: 12/18/07      Analysis Date...: 12/18/07  
 Prep Batch #....: 7353354  
 Dilution Factor: 1

<u>PARAMETER</u>	<u>PERCENT</u>	<u>RECOVERY</u>	<u>RPD</u>	<u>LIMITS</u>	<u>METHOD</u>	
	<u>RECOVERY</u>	<u>LIMITS</u>		<u>RPD</u>	<u>LIMITS</u>	<u>METHOD</u>
Benzene	100	(80 - 116)			SW846 8260B	
	101	(80 - 116)	0.31	(0-20)	SW846 8260B	
Chlorobenzene	100	(76 - 117)			SW846 8260B	
	101	(76 - 117)	1.1	(0-20)	SW846 8260B	
1,1-Dichloroethene	103	(63 - 130)			SW846 8260B	
	103	(63 - 130)	0.21	(0-20)	SW846 8260B	
Toluene	103	(74 - 119)			SW846 8260B	
	104	(74 - 119)	0.75	(0-20)	SW846 8260B	
Trichloroethene	95	(75 - 122)			SW846 8260B	
	96	(75 - 122)	0.94	(0-20)	SW846 8260B	

<u>SURROGATE</u>	<u>PERCENT</u>	<u>RECOVERY</u>	<u>LIMITS</u>
	<u>RECOVERY</u>	<u>LIMITS</u>	
Dibromofluoromethane	93	(73 - 122)	
	91	(73 - 122)	
1,2-Dichloroethane-d4	106	(61 - 128)	
	106	(61 - 128)	
Toluene-d8	94	(76 - 110)	
	94	(76 - 110)	
4-Bromofluorobenzene	99	(74 - 116)	
	98	(74 - 116)	

NOTE (S) :

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

Bold print denotes control parameters

## LABORATORY CONTROL SAMPLE DATA REPORT

## GC/MS Volatiles

**Client Lot #...**: 7L06288      **Work Order #...**: KEFNV1AC-LCS    **Matrix.....**: WATER  
**LCS Lot-Sample#**: A7L190000-354      **KEFNV1AD-LCSD**  
**Prep Date.....**: 12/18/07      **Analysis Date...**: 12/18/07  
**Prep Batch #...**: 7353354  
**Dilution Factor:** 1

<u>PARAMETER</u>	<u>SPIKE</u>	<u>MEASURED</u>	<u>UNITS</u>	<u>PERCENT</u>	<u>RPD</u>	<u>METHOD</u>
	<u>AMOUNT</u>	<u>AMOUNT</u>	<u>ug/L</u>	<u>RECOVERY</u>		
<b>Benzene</b>	10	10	ug/L	100	0.31	<b>SW846 8260B</b>
	10	10	ug/L	101		<b>SW846 8260B</b>
<b>Chlorobenzene</b>	10	10	ug/L	100	1.1	<b>SW846 8260B</b>
	10	10	ug/L	101		<b>SW846 8260B</b>
<b>1,1-Dichloroethene</b>	10	10	ug/L	103	0.21	<b>SW846 8260B</b>
	10	10	ug/L	103		<b>SW846 8260B</b>
<b>Toluene</b>	10	10	ug/L	103	0.75	<b>SW846 8260B</b>
	10	10	ug/L	104		<b>SW846 8260B</b>
<b>Trichloroethene</b>	10	9.5	ug/L	95	0.94	<b>SW846 8260B</b>
	10	9.6	ug/L	96		<b>SW846 8260B</b>

<u>SURROGATE</u>	<u>PERCENT</u>	<u>RECOVERY</u>	<u>LIMITS</u>
	<u>RECOVERY</u>	<u>LIMITS</u>	<u>RECOVERY</u>
<b>Dibromofluoromethane</b>	93	(73 - 122)	(73 - 122)
	91	(73 - 122)	
<b>1,2-Dichloroethane-d4</b>	106	(61 - 128)	(61 - 128)
	106	(61 - 128)	
<b>Toluene-d8</b>	94	(76 - 110)	(76 - 110)
	94	(76 - 110)	
<b>4-Bromofluorobenzene</b>	99	(74 - 116)	(74 - 116)
	98	(74 - 116)	

**NOTE (S) :**

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

**Bold print denotes control parameters**

**LABORATORY CONTROL SAMPLE EVALUATION REPORT**

GC Volatiles

PARAMETER	PERCENT RECOVERY	RECOVERY LIMITS	RPD	RPD LIMITS	METHOD
TPH (as Gasoline)	118	(72 - 140)			SW846 8015B
	118	(72 - 140)	0.28	(0-20)	SW846 8015B

<u>SURROGATE</u>	PERCENT RECOVERY	RECOVERY LIMITS
Trifluorotoluene	99	(10 - 150)
	96	(10 - 150)

**NOTE (S) :**

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

**Bold print denotes control parameters**

## LABORATORY CONTROL SAMPLE DATA REPORT

## GC Volatiles

**Client Lot #....:** 7L06288      **Work Order #....:** KDR2K1AC-LCS      **Matrix.....:** WATER  
**LCS Lot-Sample#:** A7L100000-082      **KDR2K1AD-LCSD**  
**Prep Date.....:** 12/07/07      **Analysis Date...:** 12/07/07  
**Prep Batch #....:** 7344082  
**Dilution Factor:** 1

<b>PARAMETER</b>	<b>SPIKE</b>	<b>MEASURED</b>		<b>PERCENT</b>	<b>RPD</b>	<b>METHOD</b>
	<b>AMOUNT</b>	<b>AMOUNT</b>	<b>UNITS</b>	<b>RECOVERY</b>		
<b>TPH (as Gasoline)</b>	200	240	ug/L	118	0.28	SW846 8015B
	200	240	ug/L	118		SW846 8015B

<b>SURROGATE</b>	<b>PERCENT</b>	<b>RECOVERY</b>	<b>LIMITS</b>
	<b>RECOVERY</b>	<b>LIMITS</b>	
Trifluorotoluene	99	(10 - 150)	
	96	(10 - 150)	

**NOTE (S) :**

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

Bold print denotes control parameters

**LABORATORY CONTROL SAMPLE EVALUATION REPORT**

## GC Volatiles

PARAMETER	PERCENT	RECOVERY	RPD	METHOD
	RECOVERY	LIMITS	RPD	
TPH (as Gasoline)	107	(72 - 140)		SW846 8015B
	111	(72 - 140)	3.7 (0-20)	SW846 8015B
SURROGATE	PERCENT	RECOVERY		
	RECOVERY	LIMITS		
Trifluorotoluene	84	(10 - 150)		
	82	(10 - 150)		

**NOTE (S) :**

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

**Bold print denotes control parameters**

## LABORATORY CONTROL SAMPLE DATA REPORT

## GC Volatiles

**Client Lot #....:** 7L06288      **Work Order #....:** KD09G1AC-LCS    **Matrix.....:** WATER  
**LCS Lot-Sample#:** A7L120000-409      KD09G1AD-LCSD  
**Prep Date.....:** 12/11/07      **Analysis Date..:** 12/11/07  
**Prep Batch #....:** 7346409  
**Dilution Factor:** 1

<u>PARAMETER</u>	<u>SPIKE</u>	<u>MEASURED</u>	<u>PERCENT</u>	<u>RPD</u>	<u>METHOD</u>
	<u>AMOUNT</u>	<u>AMOUNT</u>	<u>UNITS</u>		
<b>TPH (as Gasoline)</b>	200	210	ug/L	107	<b>SW846 8015B</b>
	200	220	ug/L	111	<b>SW846 8015B</b>
<u>SURROGATE</u>			<u>PERCENT</u>	<u>RECOVERY</u>	
Trifluorotoluene			<u>RECOVERY</u>	<u>LIMITS</u>	
			84	(10 - 150)	
			82	(10 - 150)	

NOTE (S) :

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

Bold print denotes control parameters

## LABORATORY CONTROL SAMPLE EVALUATION REPORT

## GC Volatiles

Client Lot #....: 7L06288      Work Order #....: KD1HK1AC-LCS      Matrix.....: WATER  
 LCS Lot-Sample#: A7L120000-440      KD1HK1AD-LCSD  
 Prep Date.....: 12/11/07      Analysis Date...: 12/11/07  
 Prep Batch #....: 7346440  
 Dilution Factor: 1

<u>PARAMETER</u>	<u>PERCENT</u>	<u>RECOVERY</u>	<u>RPD</u>	<u>LIMITS</u>	<u>METHOD</u>
	<u>RECOVERY</u>	<u>LIMITS</u>			
TPH (as Gasoline)	109	(72 - 140)			SW846 8015B
	118	(72 - 140)	7.8	(0-20)	SW846 8015B

<u>SURROGATE</u>	<u>PERCENT</u>	<u>RECOVERY</u>
	<u>RECOVERY</u>	<u>LIMITS</u>
Trifluorotoluene	98	(10 - 150)
	96	(10 - 150)

NOTE (S) :

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

Bold print denotes control parameters

## LABORATORY CONTROL SAMPLE DATA REPORT

## GC Volatiles

Client Lot #...: 7L06288      Work Order #...: KD1HK1AC-LCS      Matrix.....: WATER  
 LCS Lot-Sample#: A7L120000-440      KD1HK1AD-LCSD  
 Prep Date.....: 12/11/07      Analysis Date...: 12/11/07  
 Prep Batch #...: 7346440  
 Dilution Factor: 1

PARAMETER	SPIKE	MEASURED	PERCENT	RPD	METHOD
	AMOUNT	AMOUNT	UNITS		
TPH (as Gasoline)	200	220	ug/L	109	SW846 8015B
	200	240	ug/L	118	7.8 SW846 8015B
SURROGATE			PERCENT	RECOVERY	
Trifluorotoluene			RECOVERY	LIMITS	
			98	(10 - 150)	
			96	(10 - 150)	

NOTE(S) :

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

Bold print denotes control parameters

MATRIX SPIKE SAMPLE EVALUATION REPORT

GC/MS Volatiles

Client Lot #....: 7L06288      Work Order #....: KDLLN1AC-MS      Matrix.....: WG  
 MS Lot-Sample #: A7L060288-009      KDLLN1AD-MSD  
 Date Sampled...: 12/04/07 16:00 Date Received...: 12/06/07  
 Prep Date.....: 12/12/07      Analysis Date...: 12/12/07  
 Prep Batch #....: 7346476  
 Dilution Factor: 1

PARAMETER	PERCENT RECOVERY	RECOVERY LIMITS	RPD	LIMITS	METHOD
Benzene	94	(78 - 118)			SW846 8260B
	94	(78 - 118)	0.12	(0-20)	SW846 8260B
Chlorobenzene	93	(76 - 117)			SW846 8260B
	92	(76 - 117)	0.19	(0-20)	SW846 8260B
1,1-Dichloroethene	87	(62 - 130)			SW846 8260B
	88	(62 - 130)	1.4	(0-20)	SW846 8260B
Toluene	98	(70 - 119)			SW846 8260B
	98	(70 - 119)	0.42	(0-20)	SW846 8260B
Trichloroethene	91	(62 - 130)			SW846 8260B
	90	(62 - 130)	0.99	(0-20)	SW846 8260B

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS
Dibromofluoromethane	95	(73 - 122)
	97	(73 - 122)
1,2-Dichloroethane-d4	93	(61 - 128)
	93	(61 - 128)
Toluene-d8	101	(76 - 110)
	100	(76 - 110)
4-Bromofluorobenzene	103	(74 - 116)
	101	(74 - 116)

**NOTE(S) :**

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

**Bold print denotes control parameters**

MATRIX SPIKE SAMPLE DATA REPORT

GC/MS Volatiles

Client Lot #...: 7L06288      Work Order #...: KDLLN1AC-MS      Matrix.....: WG  
 MS Lot-Sample #: A7L060288-009      KDLLN1AD-MSD  
 Date Sampled...: 12/04/07 16:00 Date Received...: 12/06/07  
 Prep Date.....: 12/12/07      Analysis Date...: 12/12/07  
 Prep Batch #...: 7346476  
 Dilution Factor: 1

PARAMETER	SAMPLE	SPIKE	MEASRD	UNITS	PERCNT		
	AMOUNT	AMT	AMOUNT		RECVRY	RPD	METHOD
<b>Benzene</b>	ND	10	9.4	ug/L	94		SW846 8260B
	ND	10	9.4	ug/L	94	0.12	SW846 8260B
<b>Chlorobenzene</b>	ND	10	9.3	ug/L	93		SW846 8260B
	ND	10	9.2	ug/L	92	0.19	SW846 8260B
<b>1,1-Dichloroethene</b>	ND	10	8.7	ug/L	87		SW846 8260B
	ND	10	8.8	ug/L	88	1.4	SW846 8260B
<b>Toluene</b>	ND	10	9.8	ug/L	98		SW846 8260B
	ND	10	9.8	ug/L	98	0.42	SW846 8260B
<b>Trichloroethene</b>	ND	10	9.1	ug/L	91		SW846 8260B
	ND	10	9.0	ug/L	90	0.99	SW846 8260B

SURROGATE	PERCENT	RECOVERY	RECOVERY
	RECOVERY	LIMITS	
Dibromofluoromethane	95	(73 - 122)	
	97	(73 - 122)	
1,2-Dichloroethane-d4	93	(61 - 128)	
	93	(61 - 128)	
Toluene-d8	101	(76 - 110)	
	100	(76 - 110)	
4-Bromofluorobenzene	103	(74 - 116)	
	101	(74 - 116)	

**NOTE (S) :**

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

Bold print denotes control parameters

MATRIX SPIKE SAMPLE EVALUATION REPORT

GC/MS Volatiles

Client Lot #....: 7L06288      Work Order #....: KDQQA1AD-MS      Matrix.....: WG  
 MS Lot-Sample #: A7L070423-011      KDQQA1AE-MSD  
 Date Sampled....: 12/06/07 09:30 Date Received...: 12/07/07  
 Prep Date.....: 12/15/07      Analysis Date...: 12/15/07  
 Prep Batch #....: 7351272  
 Dilution Factor: 1

PARAMETER	PERCENT RECOVERY	RECOVERY LIMITS	RPD	LIMITS	METHOD
Benzene	91	(78 - 118)	2.4	(0-20)	SW846 8260B
	93	(78 - 118)			SW846 8260B
Chlorobenzene	92	(76 - 117)	0.87	(0-20)	SW846 8260B
	92	(76 - 117)			SW846 8260B
1,1-Dichloroethene	85	(62 - 130)	1.8	(0-20)	SW846 8260B
	87	(62 - 130)			SW846 8260B
Toluene	95	(70 - 119)	0.41	(0-20)	SW846 8260B
	95	(70 - 119)			SW846 8260B
Trichloroethene	86	(62 - 130)	6.2	(0-20)	SW846 8260B
	91	(62 - 130)			SW846 8260B

SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS
Dibromofluoromethane	90	(73 - 122)
	91	(73 - 122)
1,2-Dichloroethane-d4	98	(61 - 128)
	91	(61 - 128)
Toluene-d8	96	(76 - 110)
	94	(76 - 110)
4-Bromofluorobenzene	94	(74 - 116)
	96	(74 - 116)

NOTE (S) :

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

Bold print denotes control parameters

**MATRIX SPIKE SAMPLE DATA REPORT**

**GC/MS Volatiles**

**Client Lot #....:** 7L06288      **Work Order #....:** KDQQA1AD-MS      **Matrix.....:** WG  
**MS Lot-Sample #:** A7L070423-011      **KDQQA1AE-MSD**  
**Date Sampled....:** 12/06/07 09:30      **Date Received...:** 12/07/07  
**Prep Date.....:** 12/15/07      **Analysis Date...:** 12/15/07  
**Prep Batch #....:** 7351272  
**Dilution Factor:** 1

<b>PARAMETER</b>	<b>SAMPLE</b>	<b>SPIKE</b>	<b>MEASRD</b>	<b>UNITS</b>	<b>PERCNT</b>		<b>METHOD</b>
	<b>AMOUNT</b>	<b>AMT</b>	<b>AMOUNT</b>		<b>RECVRY</b>	<b>RPD</b>	
Benzene	ND	10	9.1	ug/L	91		SW846 8260B
	ND	10	9.3	ug/L	93	2.4	SW846 8260B
Chlorobenzene	ND	10	9.2	ug/L	92		SW846 8260B
	ND	10	9.2	ug/L	92	0.87	SW846 8260B
1,1-Dichloroethene	ND	10	8.5	ug/L	85		SW846 8260B
	ND	10	8.7	ug/L	87	1.8	SW846 8260B
Toluene	ND	10	9.5	ug/L	95		SW846 8260B
	ND	10	9.5	ug/L	95	0.41	SW846 8260B
Trichloroethene	ND	10	8.6	ug/L	86		SW846 8260B
	ND	10	9.1	ug/L	91	6.2	SW846 8260B

<b>SURROGATE</b>	<b>PERCENT</b>		<b>RECOVERY</b>
	<b>RECOVERY</b>	<b>LIMITS</b>	
Dibromofluoromethane	90	(73 - 122)	
	91	(73 - 122)	
1,2-Dichloroethane-d4	98	(61 - 128)	
	91	(61 - 128)	
Toluene-d8	96	(76 - 110)	
	94	(76 - 110)	
4-Bromofluorobenzene	94	(74 - 116)	
	96	(74 - 116)	

**NOTE(S) :**

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

Bold print denotes control parameters

MATRIX SPIKE SAMPLE EVALUATION REPORT

GC/MS Volatiles

Client Lot #...: 7L06288      Work Order #...: KDP471AE-MS      Matrix.....: WATER  
 MS Lot-Sample #: A7L070368-009      KDP471AF-MSD  
 Date Sampled...: 12/06/07      Date Received...: 12/07/07  
 Prep Date.....: 12/18/07      Analysis Date...: 12/18/07  
 Prep Batch #...: 7353354  
 Dilution Factor: 1

<u>PARAMETER</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>	<u>RPD</u>	<u>LIMITS</u>	<u>METHOD</u>
Benzene	99	(78 - 118)	1.2	(0-20)	SW846 8260B
	100	(78 - 118)			SW846 8260B
Chlorobenzene	98	(76 - 117)	0.55	(0-20)	SW846 8260B
	98	(76 - 117)			SW846 8260B
1,1-Dichloroethene	102	(62 - 130)	2.4	(0-20)	SW846 8260B
	104	(62 - 130)			SW846 8260B
Toluene	102	(70 - 119)	0.39	(0-20)	SW846 8260B
	102	(70 - 119)			SW846 8260B
Trichloroethene	93	(62 - 130)	2.1	(0-20)	SW846 8260B
	95	(62 - 130)			SW846 8260B

<u>SURROGATE</u>	<u>PERCENT RECOVERY</u>	<u>RECOVERY LIMITS</u>
Dibromofluoromethane	94	(73 - 122)
	95	(73 - 122)
1,2-Dichloroethane-d4	102	(61 - 128)
	109	(61 - 128)
Toluene-d8	95	(76 - 110)
	94	(76 - 110)
4-Bromofluorobenzene	98	(74 - 116)
	99	(74 - 116)

NOTE (S) :

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

Bold print denotes control parameters

## MATRIX SPIKE SAMPLE DATA REPORT

## GC/MS Volatiles

Client Lot #...: 7L06288      Work Order #...: KDP471AE-MS      Matrix.....: WATER  
 MS Lot-Sample #: A7L070368-009      KDP471AF-MSD  
 Date Sampled...: 12/06/07      Date Received...: 12/07/07  
 Prep Date.....: 12/18/07      Analysis Date...: 12/18/07  
 Prep Batch #...: 7353354  
 Dilution Factor: 1

PARAMETER	SAMPLE	SPIKE	MEASRD	PERCNT			METHOD
	AMOUNT	AMT	AMOUNT	UNITS	RECVRY	RPD	
Benzene	ND	10	9.9	ug/L	99		SW846 8260B
	ND	10	10	ug/L	100	1.2	SW846 8260B
Chlorobenzene	ND	10	9.8	ug/L	98		SW846 8260B
	ND	10	9.8	ug/L	98	0.55	SW846 8260B
1,1-Dichloroethene	ND	10	10	ug/L	102		SW846 8260B
	ND	10	10	ug/L	104	2.4	SW846 8260B
Toluene	ND	10	10	ug/L	102		SW846 8260B
	ND	10	10	ug/L	102	0.39	SW846 8260B
Trichloroethene	ND	10	9.3	ug/L	93		SW846 8260B
	ND	10	9.5	ug/L	95	2.1	SW846 8260B

SURROGATE	PERCENT		RECOVERY
	RECOVERY	LIMITS	
Dibromofluoromethane	94		(73 - 122)
	95		(73 - 122)
1,2-Dichloroethane-d4	102		(61 - 128)
	109		(61 - 128)
Toluene-d8	95		(76 - 110)
	94		(76 - 110)
4-Bromofluorobenzene	98		(74 - 116)
	99		(74 - 116)

## NOTE (S) :

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

Bold print denotes control parameters

MATRIX SPIKE SAMPLE EVALUATION REPORT

GC Volatiles

Client Lot #...: 7L06288      Work Order #...: KDLLN1AF-MS      Matrix.....: WG  
MS Lot-Sample #: A7L060288-009      KDLLN1AG-MSD  
Date Sampled...: 12/04/07 16:00 Date Received...: 12/06/07  
Prep Date.....: 12/08/07      Analysis Date..: 12/08/07  
Prep Batch #...: 7344082  
Dilution Factor: 1

PARAMETER	PERCENT	RECOVERY	RPD	LIMITS	METHOD
	RECOVERY	LIMITS			
TPH (as Gasoline)	88	(63 - 143)			SW846 8015B
	95	(63 - 143)	6.2	(0-20)	SW846 8015B
SURROGATE	PERCENT	RECOVERY		RECOVERY	
Trifluorotoluene		RECOVERY		LIMITS	
	99			(10 - 150)	
	98			(10 - 150)	

NOTE (S) :

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

Bold print denotes control parameters

MATRIX SPIKE SAMPLE DATA REPORT

GC Volatiles

Client Lot #....: 7L06288      Work Order #....: KDLLN1AF-MS      Matrix.....: WG  
MS Lot-Sample #: A7L060288-009      KDLLN1AG-MSD  
Date Sampled....: 12/04/07 16:00 Date Received...: 12/06/07  
Prep Date.....: 12/08/07      Analysis Date...: 12/08/07  
Prep Batch #....: 7344082  
Dilution Factor: 1

PARAMETER	SAMPLE	SPIKE	MEASRD	PERCNT			METHOD
	AMOUNT	AMT	AMOUNT	UNITS	RECVRY	RPD	
TPH (as Gasoline)	33	200	210	ug/L	88		SW846 8015B
	33	200	220	ug/L	95	6.2	SW846 8015B

SURROGATE	PERCENT		RECOVERY
	RECOVERY	LIMITS	
Trifluorotoluene	99	(10 - 150)	
	98	(10 - 150)	

NOTE (S) :

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

Bold print denotes control parameters

MATRIX SPIKE SAMPLE EVALUATION REPORT

GC Volatiles

Client Lot #...: 7L06288      Work Order #...: KDQPJ1AD-MS      Matrix.....: WG  
MS Lot-Sample #: A7L070423-001      KDQPJ1AE-MSD  
Date Sampled...: 12/05/07 11:30 Date Received..: 12/07/07  
Prep Date.....: 12/12/07      Analysis Date..: 12/12/07  
Prep Batch #...: 7346409  
Dilution Factor: 1

PARAMETER	PERCENT RECOVERY	RECOVERY LIMITS	RPD	RECOVERY LIMITS	METHOD
<b>TPH (as Gasoline)</b>	88	(63 - 143)			<b>SW846 8015B</b>
	82	(63 - 143)	5.4	(0-20)	<b>SW846 8015B</b>
SURROGATE	PERCENT RECOVERY	RECOVERY LIMITS			
Trifluorotoluene	84	(10 - 150)			
	78	(10 - 150)			

**NOTE (S) :**

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

Bold print denotes control parameters

MATRIX SPIKE SAMPLE DATA REPORT

GC Volatiles

Client Lot #....: 7L06288      Work Order #....: KDQPJ1AD-MS      Matrix.....: WG  
 MS Lot-Sample #: A7L070423-001      KDQPJ1AE-MSD  
 Date Sampled...: 12/05/07 11:30 Date Received...: 12/07/07  
 Prep Date.....: 12/12/07      Analysis Date...: 12/12/07  
 Prep Batch #....: 7346409  
 Dilution Factor: 1

PARAMETER	SAMPLE	SPIKE	MEASRD	UNITS	PERCNT		METHOD
	AMOUNT	AMT	AMOUNT		RECVRY	RPD	
TPH (as Gasoline)	33	200	210	ug/L	88	5.4	SW846 8015B
	33	200	200	ug/L	82	5.4	SW846 8015B

SURROGATE	PERCENT	RECOVERY
	RECOVERY	LIMITS
Trifluorotoluene	84	(10 - 150)
	78	(10 - 150)

NOTE (S) :

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

Bold print denotes control parameters

MATRIX SPIKE SAMPLE EVALUATION REPORT

GC Volatiles

Client Lot #....: 7L06288      Work Order #....: KDQQA1AF-MS      Matrix.....: WG  
MS Lot-Sample #: A7L070423-011                                    KDQQA1AG-MSD  
Date Sampled...: 12/06/07 09:30 Date Received...: 12/07/07  
Prep Date.....: 12/12/07      Analysis Date...: 12/12/07  
Prep Batch #:...: 7346440  
Dilution Factor: 1

PARAMETER	PERCENT	RECOVERY	RPD	PERCENT	METHOD
	RECOVERY	LIMITS		RECOVERY	
TPH (as Gasoline)	97	(63 - 143)			SW846 8015B
	90	(63 - 143)	6.8	(0-20)	SW846 8015B
SURROGATE	PERCENT	RECOVERY		RECOVERY	
Trifluorotoluene	103			(10 - 150)	
	100			(10 - 150)	

**NOTE (S) :**

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

Bold print denotes control parameters

MATRIX SPIKE SAMPLE DATA REPORT

GC Volatiles

Client Lot #....: 7L06288      Work Order #....: KDQQA1AF-MS      Matrix.....: WG  
 MS Lot-Sample #: A7L070423-011      KDQQA1AG-MSD  
 Date Sampled....: 12/06/07 09:30 Date Received...: 12/07/07  
 Prep Date.....: 12/12/07      Analysis Date...: 12/12/07  
 Prep Batch #....: 7346440  
 Dilution Factor: 1

PARAMETER	SAMPLE	SPIKE	MEASRD	PERCNT			METHOD
	AMOUNT	AMT	AMOUNT	UNITS	RECVRY	RPD	
TPH (as Gasoline)	33	200	230	ug/L	97	6.8	SW846 8015B
	33	200	210	ug/L	90		SW846 8015B

<u>SURROGATE</u>	PERCENT	RECOVERY
	<u>RECOVERY</u>	<u>LIMITS</u>
Trifluorotoluene	103	(10 - 150)
	100	(10 - 150)

NOTE (S) :

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

Bold print denotes control parameters

# CHAIN-OF-CUSTODY / Analytical Request Document

The Chain-of-Custody is a LEGAL DOCUMENT. All relevant fields must be completed accurately.

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## Required Client Information:

Company: **CRA** Report To: **Stacy Scarsella**  
 Address: **2055 Niagara Falls Blvd Suite #3** Invoice To:  
**Alasca Falls NY 14304** P.O.:  
 Phone: **716-297-6150** Project Name: **G.M. Fairbank**  
 Fax: **716-297-2265** Project Number: **12559-05-202040**  
 E-mail: **scarsella@CRAworld.com**

PAGE **1** OF **2**

		Laboratory:		Laboratory Location:		QA/QC Requirements:	
		<b>S &amp; T NY, Inc.</b>		<b>North Canton, OH</b>			



# CHAIN-OF-CUSTODY / Analytical Request Document

The Chain-of-Custody is a LEGAL DOCUMENT. All relevant fields must be completed accurately.

**ENCORE**

**Required Client Information:**

Company: <b>CRA</b>	Report To: <b>Susan Scrooch!</b>
Address: <b>Loss Niagara Falls, NY 14841</b>	Copy To: <b>Phil Harry</b>
Fax: <b>716-297-2265</b>	Invoice To: <b>P.O. Box 12559-05-202040</b>
E-mail: <b>Scroochi@CRAworld.com</b>	

PAGE **2** OF **2**

ID # **Nº 09006**

SSOW Ref. Code: **E095007**

Sample Identification:	<b>Trip Blank - A</b>	<b>WA</b>	Matrix Code	Date Collected	Time Collected	# Containers	Preservative	Analysis and Method									
								WG	Groundwater	WB	Borehole Water	WS	Surface Water	SO	Soil	SG	Sediment
1.																	
2.																	
3.																	
4.																	
5.																	
6.																	
7.																	
8.																	
9.																	
10.																	
11.																	
12.																	
13.																	
14.																	
15.																	

TOTAL NUMBER OF CONTAINERS **1 + 85 = 86**

SHIPMENT METHOD	NO. OF COOLERS	RELINQUISHED BY / AFFILIATION	DATE	TIME	RECEIVED BY / AFFILIATION	DATE	TIME	North America
<b>Fed - X</b>	<b>1</b>	<b>John P. Meyer / CRA</b>	<b>12/5/07</b>	<b>1930</b>	<b>A.S.C.</b>	<b>12/6/07</b>	<b>10:00</b>	<b>TestAmerica</b>

**Sample Condition:**

Temp in C	
Received on ice	Y/N
Sealed Cooler	Y/N
Samples intact	Y/N

Additional Comments:

**TestAmerica Cooler Receipt Form/Narrative  
North Canton Facility**

Lot Number: A7L060288

Client LRA

Project \_\_\_\_\_ Quote # 68181

Cooler Received on 12-6-07

Opened on 12-6-07

FedEx  Client Drop Off  UPS

DHL  FAS  TestAmerica Courier   
Stetson  US Cargo

Other \_\_\_\_\_ (Signature)

TestAmerica Cooler # 461

Foam Box  Client Cooler  Other \_\_\_\_\_

1. Were custody seals on the outside of the cooler? Yes  No  Intact? Yes  No  NA   
If YES, Quantity \_\_\_\_\_

Were custody seals on the outside of cooler signed and dated? Yes  No  NA   
Were custody seals on the bottles? Yes  No

If YES, are there any exceptions \_\_\_\_\_

2. Shipper's packing slip attached to this form? Yes  No

Relinquished by client? Yes  No

3. Did custody papers accompany the sample(s)? Yes  No

Yes  No   
Yes  No

4. Did you sign the custody papers in the appropriate place? Yes  No

Yes  No   
Other \_\_\_\_\_

5. Packing material used: Bubble Wrap  Foam  None

6. Cooler temperature upon receipt 2.7 °C (see back of form for multiple coolers/temps)

METHOD: IR  Other

COOLANT: Wet Ice  Blue Ice  Dry Ice  Water  None

Yes  No

7. Did all bottles arrive in good condition (Unbroken)? Yes  No

Yes  No

8. Could all bottle labels and/or tags be reconciled with the COC? Yes  No

Yes  No  NA

9. Were samples at the correct pH upon receipt? Yes  No

Yes  No  NA

10. Were correct bottles used for the tests indicated? Yes  No

Yes  No  NA

11. Were air bubbles >6 mm in any VOA vials? Yes  No

Yes  No  NA

12. Sufficient quantity received to perform indicated analyses? Yes  No

Yes  No

13. Was a Trip Blank present in the cooler? Yes  No

Were VOAs on the COC? Yes  No

Contacted PM \_\_\_\_\_ Date \_\_\_\_\_ by \_\_\_\_\_ via Voice Mail  Verbal  Other

Concerning \_\_\_\_\_

**14. CHAIN OF CUSTODY**

The following discrepancies occurred:	

**15. SAMPLE CONDITION**

Sample(s) \_\_\_\_\_ were received after the recommended holding time had expired.

Sample(s) \_\_\_\_\_ were received in a broken container.

**16. SAMPLE PRESERVATION**

Sample(s) \_\_\_\_\_ were further preserved in sample receiving to meet recommended pH level(s). Nitric Acid Lot #071707-HNO<sub>3</sub> - Sulfuric Acid Lot # 092006-H<sub>2</sub>SO<sub>4</sub>; Sodium Hydroxide Lot # 122805 -NaOH; Hydrochloric Acid Lot # 092006-HCl; Sodium Hydroxide and Zinc Acetate Lot # 050205-CH<sub>3</sub>COO<sub>2</sub>ZN/NaOH

What time was preservative added to sample(s)?

Sample(s) \_\_\_\_\_ were received with bubble > 6 mm in diameter (Notify PM)

Client ID	pH	Date	Initials

**STL Cooler Receipt Form/Narrative**  
**North Canton Facility**

<u>Client ID</u>	<u>pH</u>	<u>Date</u>	<u>Initials</u>
<u>Cooler</u>	<u>Temp</u>	<u>Method</u>	<u>Coolant</u>

Discrepancies Cont.

# CHAIN-OF-CUSTODY / Analytical Request Document

The Chain-of-Custody is a LEGAL DOCUMENT. All relevant fields must be completed accurately.

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## Required Client Information:

Company: **CRA** Report To: **Susan Schroeder**  
 Address: **2055 Niagara Falls Blvd Suite 43** Copy To: **PKH Harvey**  
**Blvd Niagara Falls, NY 14207** Invoice To:  
**P.O.: Project Name: G-M Fairfax**  
 Phone: **716 297 6150** Project Number: **12554-05-202046**  
 Fax: **716 297 2265** E-mail: **rscrach@CRAworld.com**

PAGE **1** OF **2**

ID# **N# 00763**

SSOW Ref. Code: **E045007**

Laboratory: **STL/TA**  
 Laboratory Location: **North Canton OH**

Laboratory Contact: **TAT: Standard**

Requested Due Date: **Q/A/QC Requirements: Trip Blank, Rinse, Mil/mid**

Sample Identification:	Matrix Code	Date Collected	Time Collected	# Containers	Preservative						Analysis and Method
					Unpreserved	HCl	H2SO4	HNO3	NaOH	Other	
1. <b>GW-120507-JH-016</b>	<b>W6</b>	<b>12/5/07</b>	<b>1330</b>	<b>5</b>	<b>V</b>						<b>TCL-VOC</b>
2. <b>017</b>		<b>1305</b>	<b>5</b>	<b>V</b>							<b>TPH-G</b>
3. <b>018</b>		<b>1350</b>	<b>5</b>	<b>V</b>							
4. <b>019</b>		<b>1420</b>	<b>5</b>	<b>V</b>							
5. <b>020</b>		<b>1450</b>	<b>5</b>	<b>V</b>							
6. <b>021</b>		<b>1505</b>	<b>5</b>	<b>V</b>							
7. <b>GW-120507-JH-022</b>	<b>W6</b>	<b>1520</b>	<b>5</b>	<b>V</b>							
8. <b>GW-120507-JH-023</b>	<b>W6</b>	<b>1605</b>	<b>5</b>	<b>V</b>							
9. <b>024</b>		<b>1645</b>	<b>5</b>	<b>V</b>							
10. <b>GW-120507-JH-025</b>	<b>W6</b>	<b>1710</b>	<b>915</b>	<b>V</b>							
11. <b>026</b>		<b>1730</b>	<b>15</b>	<b>V</b>							
12. <b>027</b>		<b>1830</b>	<b>5</b>	<b>V</b>							
13. <b>028</b>		<b>1945</b>	<b>5</b>	<b>V</b>							
14. <b>029</b>		<b>1115</b>	<b>5</b>	<b>V</b>							
15. <b>C30</b>		<b>1200</b>	<b>5</b>	<b>V</b>							

TOTAL NUMBER OF CONTAINERS **85**

SHIPMENT METHOD	NO. OF COOLERS	RELINQUISHED BY / AFFILIATION	DATE	TIME	RECEIVED BY / AFFILIATION	DATE	TIME
<b>Fed Ex</b>	<b>2</b>	<b>Airbill Niagara CRA</b>			<b>Kathy M. O'Brien</b>	<b>12/5/07</b>	<b>12:30pm</b>

## Sample Condition

Temp in C	
Received on Ice	Y/N
Sealed Cooler	Y/N
Samples intact	Y/N

Additional Comments:

Distribution: WHITE - Fully Executed Copy    YELLOW - Receiving Laboratory Copy    PINK - Sampler Copy

REV. 0 (6/04)

Sampler Name: \_\_\_\_\_  
 Sampler Signature: \_\_\_\_\_ Date: \_\_\_\_\_



# CHAIN-OF-CUSTODY / Analytical Request Document

The Chain-of-Custody is a LEGAL DOCUMENT. All relevant fields must be completed accurately.

PAGE 2 OF 2

## Required Client Information:

Company: **CRA**  
 Address: **2055 Niagara Falls Blvd Suite #3**  
**Loyola Falls NY 14204**  
 Phone: **716-247-6150**  
 Fax: **716-247-2265**  
 E-mail: **SScroach@crainwork.com**

## Sample Identification:

Sample ID	Valid Matrix Codes: WG Groundwater WB Barometric Water WS Surface Water SO Soil SE Sediment See Back for Additional Codes	Matrix Code	Date Collected	Time Collected	# Containers	Preservative					Analysis and Method	Remarks/Lab ID
						Unpreserved	HCl	H2SO4	HNO3	NaOH	Other:	
1. <b>G-W-120607-2H-031</b>	<b>WG</b>	<b>12/6/07</b>	<b>12:30</b>	<b>5</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>TCL-VOC</b>	
2. <b>G-W-120607-2H-032</b>	<b>WG</b>	<b>12/6/07</b>	<b>13:40</b>	<b>5</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>TPH-G</b>	
3. <b>G-W-120607-2H-033</b>	<b>WS</b>	<b>12/6/07</b>	<b>14:00</b>	<b>5</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>		
4. <b>G-W-120607-2H-034</b>	<b>WS</b>	<b>12/6/07</b>	<b>14:30</b>	<b>5</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>		
5. <b>G-W-120607-2H-035</b>	<b>WS</b>	<b>12/6/07</b>	<b>15:10</b>	<b>5</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>		
6. <b>G-W-120607-2H-036</b>	<b>WS</b>	<b>12/6/07</b>	<b>15:30</b>	<b>5</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>		
7. <b>G-W-120607-2H-037</b>	<b>WS</b>	<b>12/6/07</b>	<b>16:55</b>	<b>5</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>		
8. <b>G-W-120607-2H-038</b>	<b>WS</b>	<b>12/6/07</b>	<b>17:30</b>	<b>5</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>		
9. <b>G-W-120607-2H-039</b>	<b>WS</b>	<b>12/6/07</b>	<b>18:30</b>	<b>5</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>		
10. <b>G-W-120607-2H-040</b>	<b>WS</b>	<b>12/6/07</b>	<b>19:30</b>	<b>5</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>	<b>✓</b>		
11. <b>G-W-120607-2H-041</b>												
12. <b>G-W-120607-2H-042</b>												
13. <b>G-W-120607-2H-043</b>												
14. <b>G-W-120607-2H-044</b>												
15. <b>G-W-120607-2H-045</b>												

TOTAL NUMBER OF CONTAINERS **142 + 65 = 127**

SHIPMENT METHOD	NO. OF COOLERS	RELINQUISHED BY / AFFILIATION	DATE	TIME	RECEIVED BY / AFFILIATION	DATE	TIME
<b>F-2d-X</b>	<b>2</b>	<b>Jean Springer CRA</b>	<b>12/6/07</b>	<b>16:30</b>	<b>Mitchell M. Oster</b>	<b>12/7/07</b>	<b>12:30pm</b>

## Sample Condition

Temp in °C	
Received on ice	Y/N
Sealed Cooler	Y/N
Samples intact	Y/N

## Additional Comments:

Distribution: WHITE - Fully Executed Copy    YELLOW - Receiving Laboratory Copy    PINK - Sampler Copy

ID# **Nº 09007**  
 SSOW Ref. Code: **E045007**



**TestAmerica Cooler Receipt Form/Narrative  
North Canton Facility**

**TestAmerica**

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***END OF REPORT***

