Annual Summary Report for 2013 Groundwater Monitored Natural Attenuation Program

Bush Industries, Inc. 312 Fair Oak Street Little Valley, New York

Submitted to:

Bush Industries, Inc. Jamestown, New York

Submitted by:

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Executive Summary

GEI Consultants, Inc., P.C. (GEI) has been retained by Bush Industries, Inc. (Bush Industries) to conduct the 2013 Monitored Natural Attenuation (MNA) Program for groundwater at the property located at 312 Fair Oak Street, Little Valley, New York. The work was conducted pursuant to and in accordance with the Amended and Supplemental Order (File No.: 96-07 R9-4314-96-06) agreed to between Bush Industries and the New York State Department of Environmental Conservation (NYSDEC).

The subject property is located within the Little Valley Superfund Site (LVSS). The LVSS is currently being addressed by the United States Environmental Protection Agency (USEPA). The Record of Decision (ROD) for the LVSS specifies MNA as the remedy for trichloroethene (TCE) contaminated groundwater measured throughout the LVSS. The USEPA MNA remedy includes groundwater sampling on properties located throughout the LVSS including 312 Fair Oak Street. Bush Industries has agreed to conduct the MNA sampling on this property in accordance with the Amended and Supplemental Order. This report presents the validated results of the annual MNA sampling event conducted on the property by GEI in September 2013.

The results of the 2013 MNA sampling event for the property indicate that natural attenuation processes are occurring. The presence of daughter products and methane in groundwater samples reflect the reductive dechlorination occurring in groundwater at the property.

1. Introduction

1.1 Background and Site Description

GEI Consultants, Inc., P.C. (GEI) has been retained by Bush Industries, Inc. (Bush Industries) to conduct the 2013 Monitored Natural Attenuation (MNA) Program for groundwater at the property located at 312 Fair Oak Street, Little Valley, New York. The work was conducted pursuant to and in accordance with the Amended and Supplemental Order (File No.: 96-07 R9-4314-96-06) agreed to between Bush Industries and the New York State Department of Environmental Conservation (NYSDEC).

The subject property is located within the Little Valley Superfund Site (LVSS). The LVSS is currently being addressed by the United States Environmental Protection Agency (USEPA). The Record of Decision (ROD) for the LVSS specifies MNA as the remedy for TCE contaminated groundwater measured throughout the LVSS.

A topographic map of the Site and surrounding area prepared from a 7.5 minute series U.S. Geological Survey map is presented in Figure 1. The Site is situated on a 9.4 acre lot, and contains three contiguous buildings (see Figure 2). The USEPA MNA remedy includes groundwater sampling on properties located throughout the LVSS, including the property at 312 Fair Oak Street. Bush Industries has agreed to conduct the MNA sampling on this property in accordance with the Amended and Supplemental Order.

As NYSDEC was notified by letter dated September 15, 2008, Bush Industries entered into a contract to sell its land and improvements at 312 Fair Oak Street, Little Valley, N.Y. That transaction was completed on November 12, 2008. Bush Industries retained all rights-of-entry and authorization for Bush Industries (and NYSDEC) to continue to perform its obligations under the Amended and Supplemental Order. Also, deed restrictions have been placed upon the property prohibiting the use of groundwater. The current owner of the property is H2K Ventures, with addresses of 297 Howard Avenue, Jamestown, N.Y., and 312 Fair Oak Street, Little Valley, N.Y.

1.2 Previous Site Investigations

Bush Industries has conducted an extensive investigation of groundwater conditions at the 312 Fair Oak Street Site in concert with NYSDEC. Results are documented in the report entitled Groundwater Evaluation Report, prepared by Conestoga-Rovers & Associates (CRA) and dated February 21, 2000. The findings presented in the Groundwater Evaluation Report are summarized as follows:

- 1. The highest concentrations of TCE and its degradation products remain in the interior of the Site. There is a residual low level presence of TCE and its degradation products in the interior of the Site with concentrations in groundwater dropping precipitously along the downgradient flow path.
- 2. Concentrations of TCE at the downgradient perimeter of the Site are approximately equal to or below the New York State Groundwater criterion.
- 3. This distribution trend (rapidly declining concentrations with distance from the interior of the Site) indicates that natural attenuation processes occur limiting constituent migration and the Site does not pose a significant threat to downgradient groundwater quality.

The Groundwater Evaluation Report was approved by NYSDEC in March 2000. In May 2000, Bush Industries submitted the Remediation Report prepared by Geomatrix Consultants. The Remediation Report recommended implementation of an annual MNA sampling program at the Site. That Remediation Report was approved by NYSDEC in July 2007, along with EPA's concurrence.

1.3 MNA Program Objectives

The objectives of the natural attenuation monitoring are to:

- 1. Perform annual monitored natural attenuation (MNA) sampling events
- 2. Evaluate historic and new analytical data to monitor natural attenuation at the Site

2. Work Performed

2.1 MNA Scope of Work

The MNA monitoring work to be performed at the 312 Fair Oak Street Site is specified in the following documents:

Final Remedial Action Work Plan for the Little Valley Superfund Site Contract Number:68-W-98-214
Prepared by Tetra Tech EC, Inc.
Dated October 2006

Quality Assurance Project Plan Addendum for the Little Valley Superfund Site Contract Number:68-W-98-214
Prepared by Tetra Tech EC, Inc.
Dated September 2006

Work Plan for Natural Attenuation Monitoring, Bush Industries, Inc.
Prepared for Bush Industries, Inc.
Prepared by Geomatrix Consultants
Dated July 2007

The latter document prepared by Geomatrix governs the specific sampling program for the Site and is referred to herein as the Work Plan. In order to facilitate direct comparison of the Site analytical results with results from other wells within the LVSS sampled by USEPA, the sampling methods, analytical methods and QA/QC protocols specified by USEPA for the LVSS remediation are utilized for the Bush Industries MNA monitoring and are incorporated into the Work Plan.

In accordance with the Work Plan, the MNA Program for groundwater at the 312 Fair Oak Street Site includes the following:

- 1. Annual groundwater sampling events for the following wells: MW-D1, MW-D2, MW-2, MW-3, MW-5 and MW-6. Monitoring well locations are shown on Figure 2.
- 2. Sampling of wells using low flow methodology in accordance with the Work Plan
- 3. Analyses of samples for the following MNA analyses: Volatile Organic Chemicals (VOCs), alkalinity, sulfate, sulfide, nitrate, chloride, total organic carbon, ferrous iron, ethane, ethene and methane. The analytical program and methodology is summarized in Table 1 (except deviations as noted in Section 2.2, below).

- 4. Data validation.
- 5. Data evaluation and reporting.

These tasks are described in detail in the Work Plan.

2.2 2013 MNA Groundwater Sampling Event

GEI personnel conducted the annual MNA sampling event for the Site on September 19, 2013. Water level measurement, equipment decontamination, and low flow purge methods were in accordance with the Work Plan. Purge records are included in Table 2.

Deviations from the Work Plan during the 2013 sampling event are listed below:

- Monitoring wells MW-D1 contained less than 12 inches of water and therefore no samples could be obtained. EPA personnel were immediately notified that this monitoring well could not be sampled.
- With the prior concurrence of NYSDEC (by e-mail from Linda Ross, NYSDEC, dated September 18, 2008), VOCs were analyzed using SW-846 Third Edition Methods with USEPA Contract Laboratory Program (CLP) deliverables.

Groundwater samples were analyzed in accordance with Table 1 (except as noted above) by Test America Buffalo Laboratory.

The data validation and usability are discussed in Section 3.1. Results are presented in Section 3.2.

3. Sampling Event Results

3.1 Data Validation and Usability

The analytical results and data packages for the September 2013 sampling event reported by the laboratory were validated by MECX, LP of Aurora, Colorado. Data validation was performed in accordance with the Work Plan based on an evaluation of method specific QC information (holding times, calibration records, laboratory and field blanks, duplicate precision, and surrogate and matrix spike recoveries), the most current version of the USEPA Region 2 Data Validation SOPs (www.epa.gov/region02/desa/hsw/sops.htm), the most current version of the EPA National Functional Guidelines (www.epa.gov/superfund/programs/clp/guidance.htm) and the best professional judgment of the validator.

The Data Validation Report is included in its entirety in Appendix A. Results were deemed usable with appropriate qualifiers added (see Appendix A). No significant data quality issues were identified.

3.2 Groundwater Results

3.2.1 Hydraulic Head Measurements

Groundwater hydraulic head measurements were obtained on September 19, 2013 and are presented in Table 3. Figure 3 presents a water table elevation map prepared from these measurements.

Groundwater flow is indicated to be toward the northeast and the flow direction is consistent with prior measurement events.

3.2.2 Analytical Results

The validated analytical results are summarized in Table 4. Table 5 presents comparison criteria for detected constituents in groundwater used by USEPA for the LVSS. TCE and/or its reductive dechlorination products (cis-1,2-dichloroethene and vinyl chloride) were detected at or above the comparison criteria in 3 of the 5 wells sampled.

The highest TCE concentration was measured in the sample from well MW-D2 (reported concentration of 88 ug/L). Well MW-D2 is located in the central portion of the property. The reductive dechlorination product cis-1,2-dichloroethene was present in samples from 3 wells (28 ug/L in MW-2, 17 ug/L in MW-D2 and 14 ug/L in MW-6). The reductive dechlorination product vinyl chloride was detected in two wells (1.3 ug/L in MW-2 and 1.0 ug/L in MW-6).

No VOCs were detected in the sample from monitoring well MW-5.

Figure 4 presents an isoconcentration contour map for total VOCs measured during September 2013.

MNA parameter results are discussed in the following section.

4. Contaminant Trends and Progress of MNA

4.1 Contaminant Trends

Table 6 presents historical sampling results for the six wells in the Bush Industries MNA sampling program. Figures 5 through 10 present time versus concentration plots depicting the historical trend of TCE and daughter products in the Bush Industries MNA monitoring wells. As shown on these figures, all 2013 sampling event results for TCE and its reductive dechlorination products are within the general ranges of historical values. Results for all wells were approximately the same as measured during the 2012 event.

4.2 Reductive Dechlorination

The data obtained during the September 2013 groundwater sampling event were reviewed to assess the potential for degradation of VOCs at the Site via reductive dechlorination. EPA's Technical Protocol (EPA, 1998) was used as a basis for much of the following assessment.

Oxygen

Anaerobic bacteria generally cannot function at dissolved oxygen (DO) concentrations above 0.5 mg/L, and reductive dechlorination will not occur. As indicated in Table 2, stable field measured DO concentrations at the Site ranged from less than 1 mg/L to 4.48 mg/L. The lowest DO concentrations were measured in well MW-6. Reductive dechlorination products were detected in this well.

Nitrate

After dissolved oxygen has been depleted, nitrate may be used as an electron acceptor for the biodegradation of organic compounds via denitrification. Areas of depressed nitrate concentrations within a groundwater plume may indicate biodegradation via nitrate reduction, while the presence of nitrate in groundwater can indicate a fairly aerobic environment. Nitrate concentrations in the contaminant plume should be less than 1 mg/L for reductive dechlorination to occur. Nitrate concentrations were below 1 mg/L were measured in all wells sampled except for downgradient well MW-3.

Ferrous Iron

After nitrate, iron (III) may be used as an electron acceptor during anaerobic biodegradation, reducing the analyte to iron (II). Ferrous iron [iron (II)] concentrations were not detected above 0.1 ug/L in any wells.

Sulfate/Sulfide

After dissolved oxygen and nitrate depletion, sulfate may be used as an electron acceptor for

anaerobic biodegradation (EPA, 1998). This "sulfate reduction" process produces sulfide, and concentrations of sulfide greater than 1 mg/L indicate a possible reductive pathway. Sulfate concentrations ranged up to 17.5 mg/L. Sulfide was not detected in any well during the 2013 event.

Methane/Ethane/Ethene

EPA, 1998 states that methanogenesis (the reduction of carbon dioxide to methane) generally occurs after oxygen, nitrate, and sulfate have been depleted. Therefore, the presence of methane in groundwater is indicative of strongly reducing conditions. Samples from two wells, MW-2 and MW-6, contained detectable concentrations of methane in the 2013 event (0.075 mg/L and 0.10 mg/L, respectively). Reductive dechlorination products were detected in both of these wells.

Alkalinity

Zones of microbial activity are typically identified by an increase in alkalinity, resulting from increased concentrations of carbon dioxide produced by the metabolism of microorganisms. According to EPA, 1998, a two-fold increase in alkalinity values over background numbers suggests biodegradation may be occurring. Historically, the minimum value for alkalinity has occurred in well MW-5, which is considered upgradient of the TCE presence at the Site (historic range from approximately 60 mg/L to 70 mg/L as shown on Table 6). Alkalinity was measured at 68.5 mg/L in the 2013 sample from MW-5, and this value is used as "background" for comparison. The sample from the monitoring well MW-2 had alkalinity levels greater than twice the background concentration, as did the furthest downgradient well (MW-3).

Oxidation-Reduction Potential

The oxidation-reduction potential of groundwater is a relative measure of electron activity, and can influence rates of biodegradation. At less than 50 millivolts (mV), the reductive pathway is possible, and becomes more likely below -100 mV (EPA, 1998). Negative redox potentials were not measured in any of the wells sampled during the 2013 event.

pH and Temperature

Metabolic activity of bacteria is affected by the pH and temperature of the groundwater. The optimal values for these parameters for reductive biodegradation is a pH between 6 and 8 and a temperature greater than 20°C. All of the wells had pHs in this optimum range. Stable values of water temperature during the 2013 sampling event were between 11.1°C and 14.1°C.

Chloride

Chloride is released as a breakdown product during the biodegradation of chlorinated compounds. Chloride ions do not typically enter into oxidation-reduction reactions, form no important solute complexes, do not form salts of low solubility, are not significantly adsorbed on mineral surfaces, and play few vital biochemical roles (EPA, 1998). As a result, significant

increases in chloride concentrations relative to background (i.e., two times) may indicate the biodegradation of chlorinated compounds. Road salting also serves as a common, localized source of chloride to aquifer systems. The result from well MW-5 (14.2 mg/L), which as indicated above is considered upgradient of the TCE presence at the Site, was used as "background" for comparison of the chloride values. The well sampled downgradient from MW-5 well on the property (MW-2) had a chloride concentration of 16.7 mg/L. The furthest downgradient well, MW-3 had a chloride concentration of 20.3 mg/L.

Total Organic Carbon

The presence of natural or anthropogenic organic carbon can facilitate dechlorination, by acting as a carbon and energy source for aerobic microorganisms (which during aerobic respiration decrease dissolved oxygen levels, creating a reducing environment and increasing the potential for anaerobic bacteria to function). A TOC concentration of 20 mg/L is most favorable to dechlorination. TOC concentrations ranged from 0.44 J (estimated) in well MW-3 to 1.3 mg/L in well MW-6 for the 2013 event.

Daughter Products

Transformation of TCE via reductive dechlorination produces daughter products including 1,1-dichloroethene, 1,2-dichloroethene (cis- and/or trans-), and vinyl chloride. As described in Section 3.2, these daughter products were detected, suggesting that reductive dechlorination has occurred at the property.

4.3 Progress of MNA at the Site

The presence of daughter products and methane in groundwater samples reflect the reductive dechlorination occurring in groundwater at the property.

The 2013 results indicate concentrations of TCE and daughter products in groundwater are within the general historical ranges. Additional annual sampling data will be used to assess any long term trends in the MNA monitoring wells.

The next annual report will be due 90 days from completion of the 2014 yearly groundwater sampling, per the Work Plan.

Tables		

TABLE 1

SAMPLE COLLECTION AND ANALYSIS PROTOCOLS
312 Fair Oak Street, Little Valley, New York

Sample Type	Matrix	Sampling Device	No. of Samples ⁽¹⁾⁽²⁾	Parameter	Sample Container ⁽³⁾⁽⁴⁾	Sample Preservation	Analytical Method ⁽⁵⁾	PQL	Holding Time ⁽⁶⁾
Groundwater	Water	Positive Displacement Submersible Pump	6	pH; temperature; specific conductivity DO; ORP; turbidity [Field Screening]	NA	NA	Direct Field Measurement Following SOP	NA	Analyze Immediately
			6	Low Concentration TCL Volatile Organic Compounds [CLP Lab]	(4) 40 mL VOA vials w/Teflon lined septum	1:1 HCl to pH<2; Cool to 4°C	SOM01.1	Compound specific (0.5 - 20 μg/L)	10 days
			6	Total Organic Carbon [DESA Lab]	(1) L amber glass	H ₂ SO ₄ to pH<2; Cool to 4°C	SW-846 Method 9060	1 mg/L	28 days*
			6	Alkalinity [DESA Lab]	(1) 1 L polyethelyene	Cool to 4°C	MCAWW Method 310.1	1 mg/L	14 days*
			6	Sulfate [DESA Lab]	(1) 1 L polyethelyene	Cool to 4°C	EPA 300.1	1 mg/L	28 days*
			6	Sulfide [DESA Lab]	(1) 1 L polyethelyene	NaOH to pH >12; 4 drops of zinc acetate per liter; Cool to 4°C	MCAWW Method 376.1	1 mg/L	7 days*
			6	Nitrate [DESA Lab]	(1) 1 L polyethelyene	Cool to 4°C	EPA 300.1	0.05 mg/L	48 hours*
			6	Chloride [DESA Lab]	(1) 1 L polyethelyene	Cool to 4°C	EPA 300.1	1 mg/L	28 days*
			6	Ferrous Iron [Sub Lab]	(1) 100 mL amber glass	2mL HCI; Cool to 4°C	Std. Methods Method 3500Fe-D	10 μg/L	24 hours*
			6	Ethane [Sub Lab]	(5) 40-mL VOA vials w/Teflon lined septum	Cool to 4°C	GC/FID (SW-846 Method 3810)	5 μg/L	7 days*
			6	Ethene [Sub Lab]	(5) 40-mL VOA vials w/Teflon lined septum	Cool to 4°C	GC/FID (SW-846 Method 3810)	5 μg/L	7 days*
			6	Methane [Sub Lab]	(5) 40-mL VOA vials w/Teflon lined septum	Cool to 4°C	GC/FID (SW-846 Method 3810)	5 μg/L	7 days*

C:\Users\kmcintosh\Documents\Bush Industries\2012 MNA\Tables\Table 1

TABLE 1 SAMPLE COLLECTION AND ANALYSIS PROTOCOLS 312 Fair Oak Street, Little Valley, New York

Sample Type	Matrix	Sampling Device	No. of Samples ⁽¹⁾⁽²⁾	Parameter	Sample Container ⁽³⁾⁽⁴⁾	Sample Preservation	Analytical Method (5)	PQL	Holding Time ⁽⁶⁾
Field Blank	Water	Collected Rinsate Passed Over/Through Sampling Equipment	1	Low Concentration TCL Volatile Organic Compounds [CLP Lab]	(4) 40-mL VOA vials w/Teflon lined septum	1:1 HCl to pH<2; Cool to 4°C	SOM01.1	Compound specific (0.5 - 20 µg/L)	10 days
Trip Blank	Water	Direct Fill of Sample Bottles	1	Low Concentration TCL Volatile Organic Compounds [CLP Lab]	(4) 40-mL VOA vials w/Teflon lined septum	1:1 HCl to pH<2; Cool to 4°C	SOM01.1	Compound specific (0.5 - 20 µg/L)	10 days
			6	Ethane [Sub Lab]	(5) 40-mL VOA vials w/Teflon lined septum	Cool to 4°C	GC/FID (SW-846 Method 3810)	5 μg/L	7 days*
			6	Ethene [Sub Lab]	(5) 40-mL VOA vials w/Teflon lined septum	Cool to 4°C	GC/FID (SW-846 Method 3810)	5 μg/L	7 days*
			6	Methane [Sub Lab]	(5) 40-mL VOA vials w/Teflon lined septum	Cool to 4°C	GC/FID (SW-846 Method 3810)	5 μg/L	7 days*

NOTES:

- 1. The number in parentheses in the "No. of Samples" column denotes the number of duplicate samples.
- 2. The number of field, trip and DI water blanks is estimated based on the approximate number of days in the field for each type of sampling during the MNA Program events.
- 3. The number in parentheses in the "Sample Container" column denotes the number of containers needed. Additional volume must be sent for laboratory QA/QC sample analyses.
- 4. All bottles will comply with OSWER Directive 9240.0-05A: "Specifications and Guidance for Obtaining Contaminant-Free Sample Containers", EPA 540/R-93/051, December 1992.
- Method References:

SOM01.1 = USEPA Contract Laboratory Program Statement of Work for Multi-Media, Multi-Concentration Organics (May 2005 or latest revision).

MCAWW = Methods for Chemical Analysis of Water and Wastes, March 1983.

Std. Methods = Standard Methods for the Examination of Water and Wastewater, 20th Edition (January 2000).

SW-846 = Test Methods for Evaluating Solid Waste, Physical/Chemical Methods (November 1986, revised through November 2000 via Updates I through IVB).

EPA300.1 = Determination of Inorganic Anions in Drinking Water by Ion Chromatography, Revision I (27 April 1999).

EPA/600/R-98128 = Technical Protocol for Evaluating Natural Attenuation of Chlorinated Solvents in Groundwater (September 1998).

6. All holding times listed are from Verified Time of Sample Receipt (VTSR) unless noted otherwise (* denotes from time of sample collection).

7. Acronyms/Abbreviations used:

CLP = Contract Laboratory Program DESA = Division of Environmental Science and Assessment

DO = Dissolved Oxygen ORP = Oxidation-Reduction Potential PQL = Practical Quantitation Limit Sub Lab = Non-RAS Subcontract Laboratory

TCL = Target Compound List VOA = Voalitle Organic Analysis

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TABLE 2 MONITORING WELL PURGE SUMMARY 312 Fair Oak Street Little Valley, New York

Time	Cumulative Volume (L)	Temperature (degrees C) pH		Specific Conductance (us/cm)	Dissolved Oxygen (mg/L)	Redox Potential (mV)
MW-2						
12:25	Begin Purge					
12:35	3.5	11.14	6.93	0.350	0.92	147
12:40	5.3	11.11	6.97	0.351	0.82	150
12:45	7.0	11.22	6.97	0.352	0.80	151
12:50	8.7	11.23	6.98	0.353	0.78	152
MW-3						
9:22	Begin Purge					
9:30	2.4	11.27	6.60	0.323	4.66	215
9:35	3.9	11.23	6.59	0.321	4.62	146
9:40	5.4	11.13	6.63	0.318	4.53	213
9:45	6.9	11.13	6.65	0.317	4.50	212
9:50	8.4	11.14	6.67	0.317	4.48	211
MW-5			J			ı
15:05	Begin Purge					
15:15	3.5	14.12	6.14	0.164	0.58	162
15:20	5.3	14.13	6.24	0.165	0.59	165
15:25	7	14.12	6.29	0.167	0.66	167
15:30	8.7	14.13	6.30	0.167	0.66	165
MW-6	<u> </u>				·	
17:45	Begin Purge					
17:55	3.5	13.54	6.20	0.188	1.370	80
18:00	5.3	13.54	6.20	0.189	1.260	77
18:05	7.0	13.51	6.21	0.197	0.630	60
18:10	8.7	13.50	6.21	0.197	0.62	56
MW-D1	0.7	10.00	0.21	0.157	0.02	
<u></u>	Less t	han 1 foot water	in well, no s	ample collected		
MW-D2						
16:25	Begin Purge					
16:30	1.8	9.77	6.95	0.315	3.50	196
16:35	3.5	9.68	6.95	0.309	3.34	194
16:40	5.3	9.58	6.98	0.281	2.76	190
16:45	7.0	9.55	7.00	0.276	2.63	188
16:50	8.7	9.56	6.99	0.276	2.62	187

TABLE 3 GROUNDWATER ELEVATION SUMMARY 312 Fair Oak Street Little Valley, New York

Well ID	Measuring Point Elevation (fasl)	DTW (ft.) 9/19/13	Groundwater Elevation (fasl)
MW-2	1590.18	39.76	1550.42
MW-3	1591.37	55.20	1536.17
MW-5	1590.44	6.98	1583.46
MW-6	1584.99	3.71	1581.28
MW-D1	1590.31	50.52	1539.79
MW-D2	1584.17	40.10	1544.07

Notes:

DTW- depth to water fasl- feet above sea level

TABLE 4 VALIDATED GROUNDWATER ANALYTICAL SUMMARY 312 Fair Oak Street Little Valley, New York

Sample ID:	LVRA07-MNAGW- MW2	LVRA07-MNAGW- MW3	LVRA07-MNAGW- MW5	LVRA07-MNAGW- MW6	LVRA07- MNAGW-MWD2	LVRA03- MNAGW- DUP ⁽¹⁾
Date Sampled:	09/19/13	09/19/13	09/19/13	09/19/13	09/19/13	09/19/13
Volatile Organic Compounds (ug/L)	•		•			•
1,1,1-Trichloroethane	1U	1U	1U	1U	1U	1U
1,1,2,2-Tetrachloroethane	1U	1U	1U	1U	1U	1U
1,1,2-Trichloro-1,2,2,-trifluoroethane	1U	1U	1U	1U	1U	1U
1,1,2-Trichloroethane	1U	1U	1U	1U	1U	1U
1,1-Dichloroethane	1U	1U	1U	1U	1U	1U
1,1-Dichloroethane	0.57 J	1U	1U	1U	1U	1U
7	1U	1U	1U	1U	1U	1U
1,2-Dibromo-3-Chloropropane	1U	1U	1U	1U	1U	1U
1,2-Dibromoethane	1U	1U	1U	1U	1U	1U
1,2-Dichlorobenzene						
1,2-Dichloroethane	1U	1U	1U	1U	1U	1U
1,2,4-Trichlorobenzene	1U	1U	1U	1U	1U	1U
1,2-Dichloropropane	1U	1U	1U	1U	1U	1U
1,4-Dichlorobenzene	1U	1U	1U	1U	1U	1U
2-Hexanone	5U	5U	5U	5U	5U	5U
2-Butanone	10U	10U	10U	10U	10U	10U
4-Methyl-2-pentanone	5U	5U	5U	5U	5U	5U
Acetone	10U	10U	10U	10U	10U	10U
Benzene	1U	1U	1U	1U	1U	1U
1,3-Dichlorobenzene	1U	1U	1U	1U	1U	1U
Bromodichloromethane	1U	1U	1U	1U	1U	1U
Bromoform	1U	1U	1U	1U	1U	1U
Bromomethane	1U	1U	1U	1U	1U	1U
Carbon Disulfide	1U	1U	1U	1U	1U	1U
Carbon Tetrachloride	1U	1U	1U	1U	1U	1U
Chlorobenzene	1U	1U	1U	1U	1U	1U
Dibromochloromethane	1U	1U	1U	1U	1U	1U
Chloroethane	1U	1U	1U	1U	1U	1U
Chloroform	1U	1U	1U	1U	1U	1U
Chloromethane	1U	1U	1U	1U	1U	1U
cis-1,2-Dichloroethene	28	1U	1U	14	17	17
cis-1,3-Dichlroropropene	1U	1U	1U	1U	1U	1U
Ethylbenzene	1U	1U	1U	1U	1U	1U
Methylene Chloride	1U	1U	1U	1U	1U	1U
Styrene	1U	1U	1U	1U	1U	1U
Tetrachloroethene	1U	1U	1U	1U	1U	1U
Toluene	1U	1U	1U	1U	1U	1U
trans-1,2,-Dichloroethene	1U	1U	1U	1U	1U	1U
trans-1,3-Dichloropropene	1U	1U	1U	1U	1U	1U
Trichloroethene	61	3.9	1U	2.2	88	87
Trichlorofluoromethane	1U	1U	1U	1U	1U	1U
Vinyl Chloride	1.3	1U	1U	1.0	1U	1U
Total Xylenes	2U	2U	2U	2U	2U	2U
Cyclohexane	1U	1U	1U	1U	1U	1U
Dichlorodifluoromethane	1U	1U	1U	1U	1U	1U
Isopropylbenzene	1U	1U	1U	1U	1U	1U
Methyl acetate	1U	1U	1U	1U	1U	1U
Methyl-tert-butyl ether	1U	1U	1U	1U	1U	1U
Methylcyclohexane	1U	1U	1U	1U	1U	1U

Monitored Natural Attenuation Paran	Monitored Natural Attenuation Parameters (mg/L)												
Chloride	16.7	20.3	14.2	10.9	20.2	20.2							
Ethane	0.00049 J	0.0075 U											
Ethene	0.007U	0.007 U											
Ferrous Iron	0.10 U	0.10U	0.10U	0.08 J	0.10U	0.10U							
Methane	0.075	0.004 U	0.004 U	0.10	0.004 U	0.004 U							
Nitrate	0.052	1.2	0.051	0.034 J	0.22	0.22							
Sulfate	17.5	10.6	5.4	10.6	15.4	14.2							
Sulfide	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U	1.0 U							
Total Alkalinity	201	165	68.5	92.2	141	142							
Total Organic Carbon	0.82 J	0.44 J	0.69 J	1.3	1.0U	1.0U							

Notes

- $U = \ Compound \ not \ detected \ above \ specified \ laboratory \ detection \ limit$
- J= Laboratory estimated concentration
- (1) Duplicate sample collected at LVRA07-MNAGW-MW-D2 location

TABLE 5
Comparison Criteria for Detected Constituents in Groundwater

BASIS FOR CRITERIA	HUMAN HEALTH	STATE
	EPA Maximum Contaminant Level	NYSDEC Water Quality Values [Class GA]
Volatile Organics (ug/L)		
1,1,2-Trichloroethane	200	5
1,1-Dichloroethene	7	5
1,2,3-Trichlorobenzene	NC	5
1,2,4-Trichlorobenzene	70	5
1,2-Dichlorobenzene	600	3
1,2-Dichloroethane	5	0.6
1,2-Dichloroethene (total)	70 🗆	5
cis-1,2-Dichloroethene	70	5
trans-1,2-Dichloroethene	100	5
1,2-Dichloropropane	5	1
1,3-Dichlorobenzene	NC	3
1,4-Dichlorobenzene	75	3
2-Hexanone	NC	50
Acetone	NC	50
Benzene	5	1
Carbon disulfide	NC	60
Chlorobenzene	100	5
Chloroethane	NC	5
Cyclohexane	NC	NC
Ethylbenzene	700	5
Methyl chloride (Chloromethane)	NC	5
Methyl ethyl ketone (2-Butanone)	NC	50
Methyl isobutyl ketone (4-Methyl-2-pentanone)	NC	NC
Methylcyclohexane	NC	NC
Styrene	100	5
Tetrachloroethene	5	5
Toluene	1000	5
Trichloroethene	5	5
m/p-Xylene	10000	5
Xylenes (total)	10000	5

TABLE 6 Historical Summary of Detected Groundwater Constituents in MNA Wells 312 Fair Oak Street

								BIA	AMW-2							
	05/05/1999	05/05/1999	12/14/1999	12/14/1999	01/10/2001	12/11/2003	10/31/2006	10/31/2006	09/25/2007	09/25/2008	09/22/2009	09/21/2010	09/28/2011	09/26/2012	09/26/2012	09/19/2013
		Duplicate		Duplicate				Duplicate							Duplicate	
Volatile Organics (ug/L)																
1,1,2-Trichloroethane					-		-						-			
1,1-Dichloroethene	1 J		0.7 J	0.7 J		0.63	0.8	0.89	0.73	0.6	0.58 J	0.51 J		0.62 J	0.29 J	0.57 J
1,4-Dichlorobenzene	NA	NA	NA	NA	-		0.16 J	0.12 J					1			
Benzene	0.7 J		0.4 J	0.4 J	-	0.32 J	ł		0.29 J				ł			
2-Butanone					-		1						1			
Chloroethane	0.8 J				-		0.19 J	0.23 J					ł			
1,2-Dichloroethene (total)	54	51	40	42	NA											
cis-1,2-Dichloroethene	NA	NA	NA	NA	44	40 D	45 D	46 D	54 D	42	29	32	31	31	23	28
trans-1,2-Dichloroethene	NA	NA	NA	NA		0.28 J	0.51	0.49 J	0.47 J							
Ethylbenzene							0.25 J									
Isopropylbenzene	NA	NA	NA	NA	NA		0.14 J									
Toluene																
Trichloroethene	230	190	84	87	110	36 D	58 D	58 D	69 J	75	77	75	75	79	51	61
Vinyl Chloride	4 J	2 J	1 J	1 J	NA	4.8	4	4.8	4.2	3	0.77 J	2.1	1.2	1.3		1.3
m/p-Xylene	NA	NA	NA	NA	NA	NA	0.1 J									
MNA/Water Quality																
Parameters (mg/L)																
Alkalinity	NA	NA	NA	NA	NA	180	190	180	176	194	173	194	189	201	201	201
Chloride	NA	NA	NA	NA	NA	19	26	26	28.4	32.2	25.2	24.1	21.9	19.1	19.1	16.7
Ferrous Iron	NA	NA	NA	NA	NA		0.17	0.14						0.09J	0.09J	0.10 U
Methane	NA	NA	NA	NA	NA	0.54 JD	0.046 J	0.11 J	0.026	0.020	0.009	0.052	0.035	0.0059	0.0059	0.0750
Nitrate	NA	NA	NA	NA	NA		ł						ł	0.027 J	0.027 J	0.052
Sulfate	NA	NA	NA	NA	NA	16	17	17	20.5	21.2	16.5	17	16	16.3	16.3	17.5
Sulfide	NA	NA	NA	NA	NA	NA	0.02	0.018					1			
TOC	NA	NA	NA	NA	NA	2.6			1.6		1.6 J	0.9J	1.1	0.97 J	0.97 J	0.82 J

TABLE 6 Historical Summary of Detected Groundwater Constituents in MNA Wells 312 Fair Oak Street

						BIAN	MW-3					
	01/09/2001	12/10/2003	10/30/2006	09/25/2007	09/25/2008	09/25/2008 Duplicate	09/22/2009	09/22/2009 Duplicate	09/21/2010	09/28/2011	09/28/2011 Duplicate	09/19/2013
Volatile Organics (ug/L)						Duplicate		Бирпсис			Duplicate	
1.1.2-Trichloroethane												
1,1-Dichloroethene												
1,4-Dichlorobenzene												
Benzene			0.12 J									
2-Butanone												
Chloroethane			0.091 J									
1,2-Dichloroethene (total)	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
cis-1,2-Dichloroethene	3	2.2	0.36 J	0.86	0.7	0.8			1			
trans-1,2-Dichloroethene												
Ethylbenzene												
Isopropylbenzene	NA											
Toluene												
Trichloroethene	8	6.3	2.2	7.9 J	5	6	4.2	3.7	11	3.3	3.3	3.9
Vinyl Chloride												
m/p-Xylene	NA	NA										
MNA/Water Quality Parameters (mg/L)												
Alkalinity	NA	160	260	155	167	168	171	173	155	168	167	165
Chloride	NA	44	78	64.4	46.0	46.3	31.8	32.1	42.3	26.2	26.1	20.3
Ferrous Iron	NA			0.18								
Methane	NA	0.07 J N		-		-						
Nitrate	NA	1.2	1.9	1.5	1.4	1.3	1.43	1.46	1.04	1.4	1.3	1.2
Sulfate	NA	12	27	23.8	13.8	13.2	11.5	11.0	14.1	10.9	10.8	10.6
Sulfide	NA	NA	0.018									
TOC	NA		26	1.4					0.4J	1.0U	1.0U	0.44 J

TABLE 6 Historical Summary of Detected Groundwater Constituents in MNA Wells 312 Fair Oak Street

		BIAMW-5											
	12/13/1999	01/04/2001	10/30/2006	09/25/2007	09/25/2008	09/22/2009	12/15/2010	09/28/2011	09/26/2012	09/19/2013			
Volatile Organics (ug/L)	1												
1,1,2-Trichloroethane													
1,1-Dichloroethene													
1,4-Dichlorobenzene	NA												
Benzene			0.23 J										
2-Butanone						3.6 J							
Chloroethane			0.13 J										
1,2-Dichloroethene (total)		NA											
cis-1,2-Dichloroethene	NA												
trans-1,2-Dichloroethene													
Ethylbenzene			0.13 J										
Isopropylbenzene	NA	NA											
Toluene						5.5							
Trichloroethene													
Vinyl Chloride													
m/p-Xylene	NA	NA											
MNA/Water Quality Parameters (mg/L)													
Alkalinity	NA	NA	70	65	65.4	61.8	60.0	66.0	64.8	68.5			
Chloride	NA	NA	11	38.4	23.3	12	9.49	11.6	17.2	14.2			
Ferrous Iron	NA	NA	0.18						0.077J				
Methane	NA	NA		0.0061		0.00031 J							
Nitrate	NA	NA	0.73				1.07	0.31	0.14	0.051			
Sulfate	NA	NA	6.7	7.4	6.4	5.31	6.92	5.4	6.9	5.4			
Sulfide	NA	NA											
TOC	NA	NA		1.3		1.1 J	1.2	0.79J	0.86 J	0.69 J			

TABLE 6 Historical Summary of Detected Groundwater Constituents in MNA Wells 312 Fair Oak Street

	BIAMW-6										
	12/13/1999	01/10/2001	10/30/2006	09/25/2007	09/25/2008	09/22/2009	09/21/2010	09/28/2011	09/26/2012	09/19/2013	
Volatile Organics (ug/L)										<u> </u>	
1,1,2-Trichloroethane											
1,1-Dichloroethene				0.66							
1,4-Dichlorobenzene	NA										
Benzene											
2-Butanone											
Chloroethane			0.11 J								
1,2-Dichloroethene (total)	30	NA									
cis-1,2-Dichloroethene	NA	44	35 D	120	39	26	27	13	5.8	14	
trans-1,2-Dichloroethene	NA		0.48 J	0.31 J							
Ethylbenzene											
Isopropylbenzene	NA	NA									
Toluene											
Trichloroethene	17	37	19	1.6 J	3	3.7	2.1	1.1	0.60 J	2.2	
Vinyl Chloride	4 J			9.5 J	5	2.5	2.6	0.99J		1.0	
m/p-Xylene	NA	NA									
MNA/Water Quality Parameters (mg/L)											
Alkalinity	NA	NA	88	75	86.1	92	89.8	61.5	58.7	92.2	
Chloride	NA	NA	13	32.9	17.8	11.3	13.3	6.5	6.8	10.9	
Ferrous Iron	NA	NA								0.08 J	
Methane	NA	NA	0.082 J	0.098	0.064	0.098	0.037	0.020	0.0013J	0.10	
Nitrate	NA	NA					0.054			0.034 J	
Sulfate	NA	NA	11	19.4	10.1	10.9	10.6	9.4	9.0	10.6	
Sulfide	NA	NA									
TOC	NA	NA		1.7		2.5 J	1.6	2.6	1.7	1.3	

TABLE 6 Historical Summary of Detected Groundwater Constituents in MNA Wells 312 Fair Oak Street

	BIAMW-D1									
	12/13/1999	01/10/2001	12/10/2003	10/31/2006	09/22/2009	09/21/2010				
Volatile Organics (ug/L)										
1,1,2-Trichloroethane										
1,1-Dichloroethene										
1,4-Dichlorobenzene	NA									
Benzene										
2-Butanone										
Chloroethane										
1,2-Dichloroethene (total)	4 J	NA	NA	NA	NA	NA				
cis-1,2-Dichloroethene	NA	8	4.8	0.42 J	0.92 J					
trans-1,2-Dichloroethene	NA			0.55						
Ethylbenzene										
Isopropylbenzene	NA	NA								
Toluene										
Trichloroethene	9 J	18	12	1.8	6.7	2.2				
Vinyl Chloride				0.16 J						
m/p-Xylene	NA	NA	NA							
MNA/Water Quality										
Parameters (mg/L)		T		1	1					
Alkalinity	NA	NA	190	200	151	165				
Chloride	NA	NA	42	55	23.8	50.4				
Ferrous Iron	NA	NA								
Methane	NA	NA	0.06 J N							
Nitrate	NA	NA	1.4	2.7	1.6	2.67				
Sulfate	NA	NA	13	11	11.7	11.5				
Sulfide	NA	NA	NA							
TOC	NA	NA			1.0 J	0.8J				

TABLE 6
Historical Summary of Detected Groundwater Constituents in MNA Wells
312 Fair Oak Street

	BIAMW-D2											
	12/14/1999	01/10/2001	01/10/2001	12/11/2003	10/30/2006	09/25/2007	09/25/2007	09/25/2008	09/22/2009	09/21/2010	09/26/2012	09/19/2013
			Duplicate				Duplicate					
Volatile Organics (ug/L)												-
1,1,2-Trichloroethane					0.084 J							
1,1-Dichloroethene	0.4 J			0.81	0.54	0.44 J	0.47 J		0.71 J	0.71 J		
1,4-Dichlorobenzene	NA											
Benzene												
2-Butanone												
Chloroethane					0.11 J							
1,2-Dichloroethene (total)	16	NA										
cis-1,2-Dichloroethene	NA	36	29	18 D	26 D	33	33	25	32	16	19	17
trans-1,2-Dichloroethene	NA				0.71	0.31 J	0.23 J					
Ethylbenzene												
Isopropylbenzene	NA	NA	NA									
Toluene												
Trichloroethene	58	140	110	78 D	93 D	110 J	110 J	93	140	72	98	88
Vinyl Chloride												
m/p-Xylene	NA	NA	NA	NA								
MNA/Water Quality												-
Parameters (mg/L)												
Alkalinity	NA	NA	NA	130	140	116	116	133	154	126	138	141
Chloride	NA	NA	NA	22	31	37.8	37.8	33.4	27.3	28.1	23.9	20.2
Ferrous Iron	NA	NA	NA			0.23						
Methane	NA	NA	NA	0.07 JN								
Nitrate	NA	NA	NA	0.29	0.34	0.23	0.22	0.24	0.416	0.189	0.20	0.22
Sulfate	NA	NA	NA	15	13	19.8	19.1	16.8	17	13.2	13.3	15.4
Sulfide	NA	NA	NA	NA	0.027							
TOC	NA	NA	NA	2.4		1.8			0.9 J	1.0U	1.0U	1.0U

Notes:

-- Not detected

J Estimated concentration.

D Value derived from dilution analysis.N Evidence exists for constituent presence.

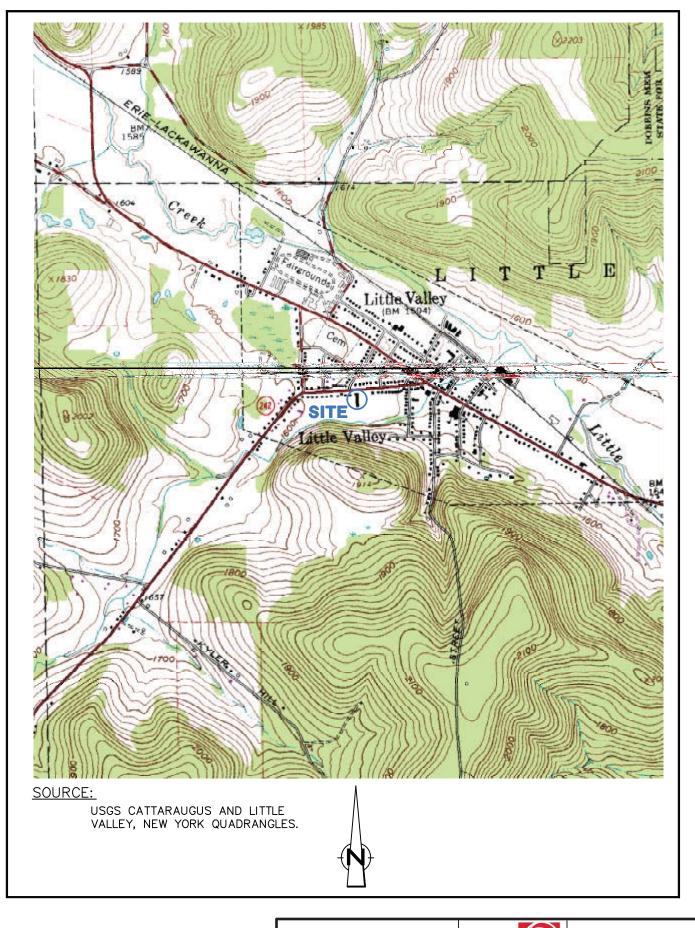
NA Not analyzed.

Above human health-based values.

Above state values.

Above both of the above values.

Figures		



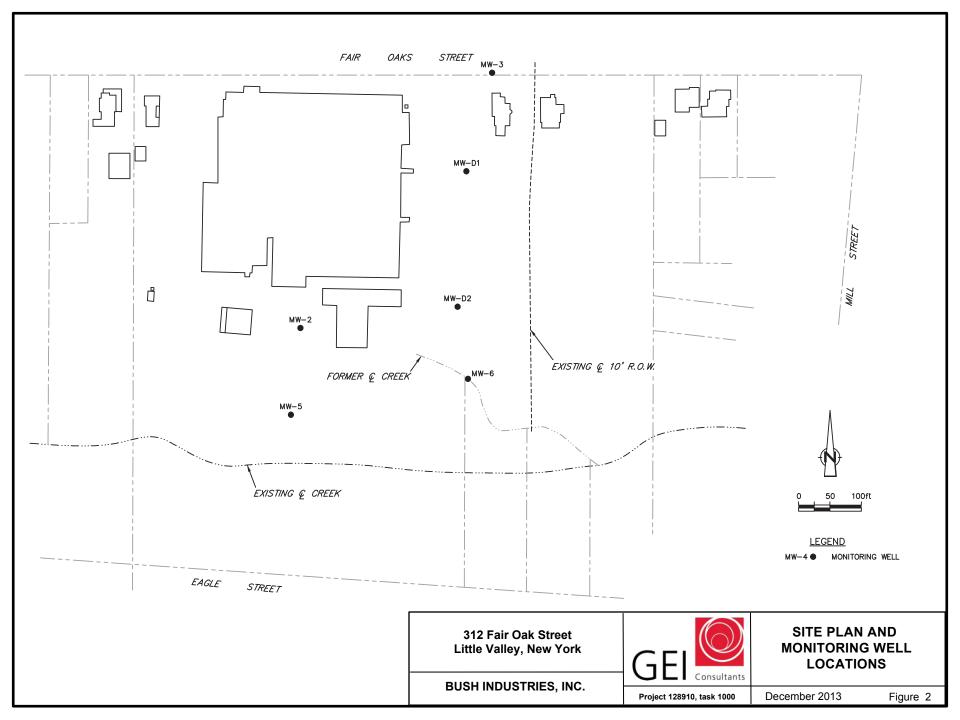
312 Fair Oak Street Little Valley, New York

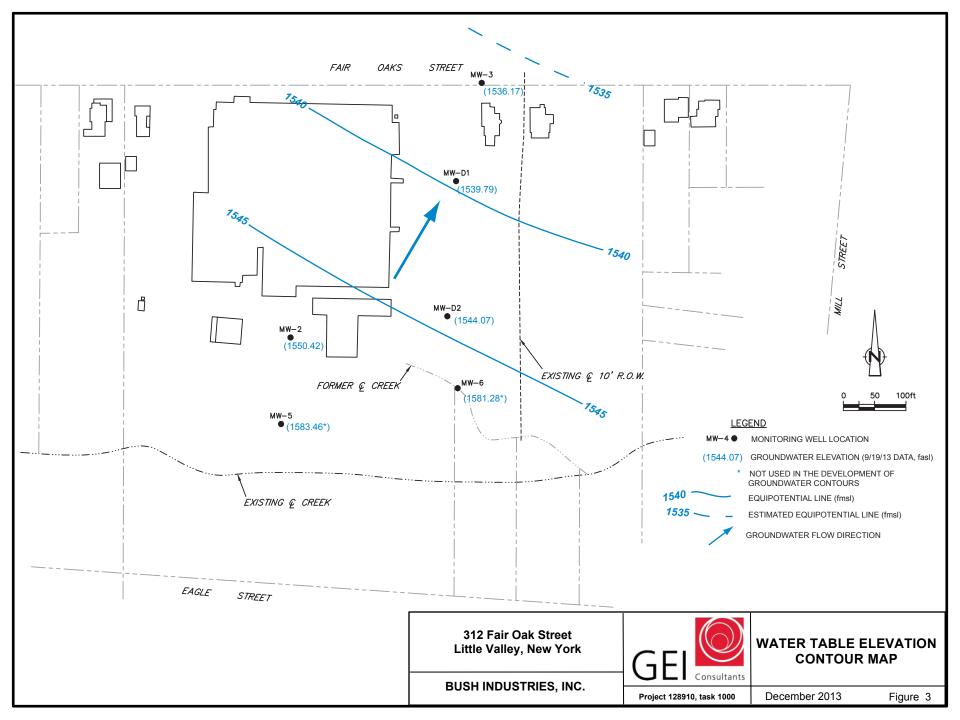
BUSH INDUSTRIES, INC.

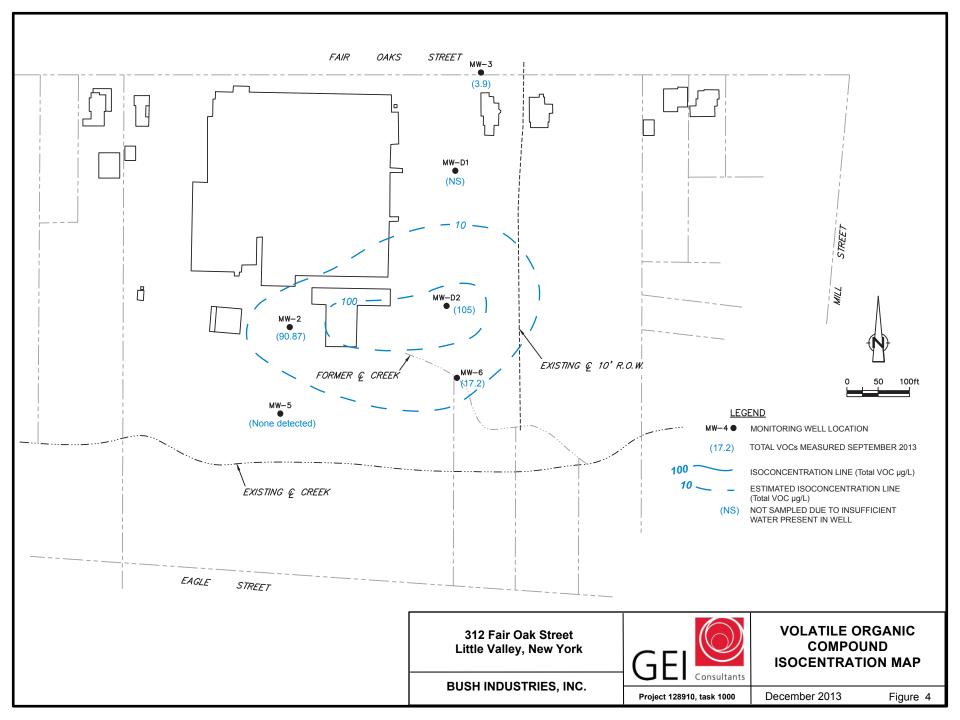


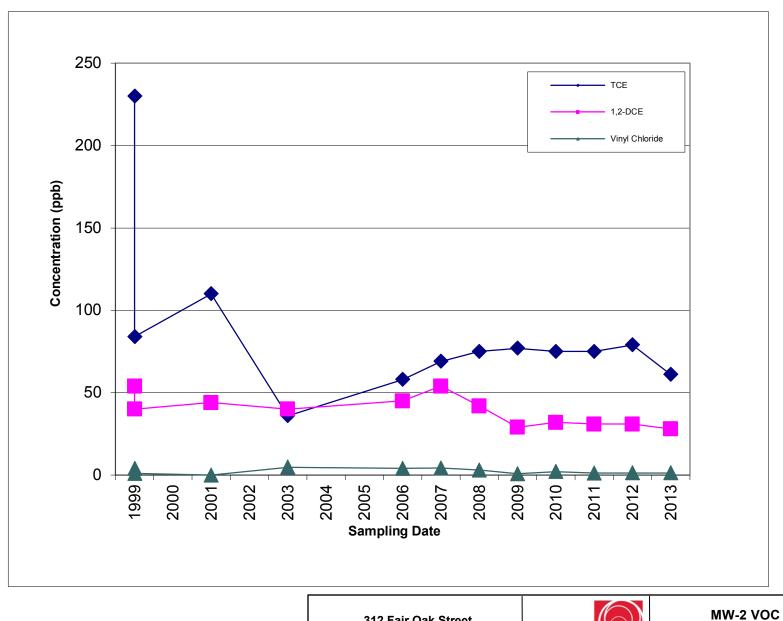
SITE LOCATION

Project 128910, task 1000 December









312 Fair Oak Street Little Valley, New York

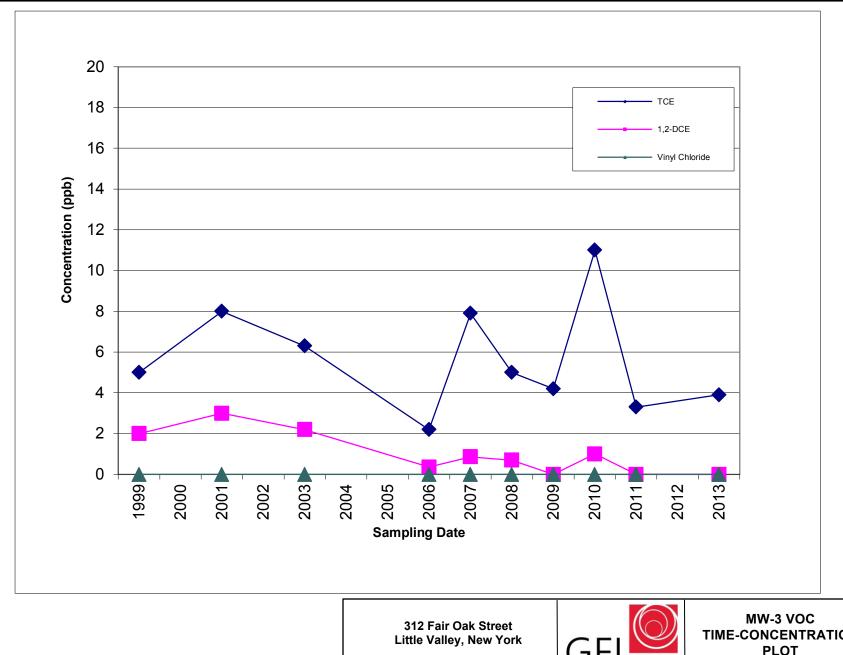
BUSH INDUSTRIES, INC.



MW-2 VOC
TIME-CONCENTRATION
PLOT

Project 128910, task 1000

December 2013

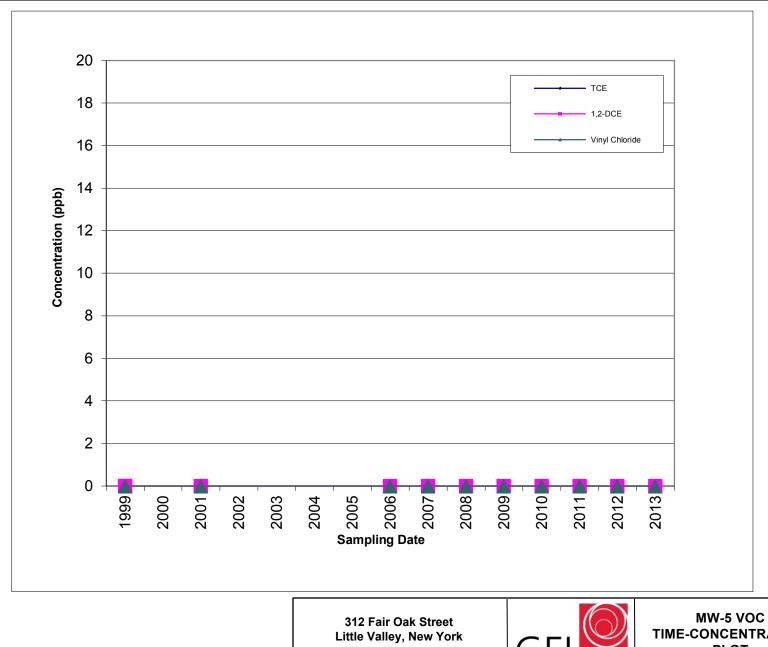


BUSH INDUSTRIES, INC.



TIME-CONCENTRATION PLOT

December 2013 Project 128910, task 1000



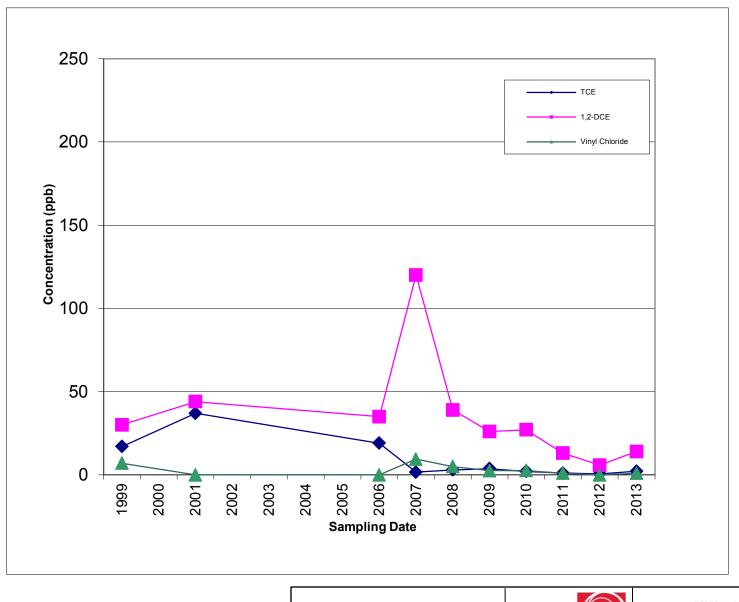
BUSH INDUSTRIES, INC.



TIME-CONCENTRATION PLOT

Project 128910, task 1000

December 2013



312 Fair Oak Street Little Valley, New York

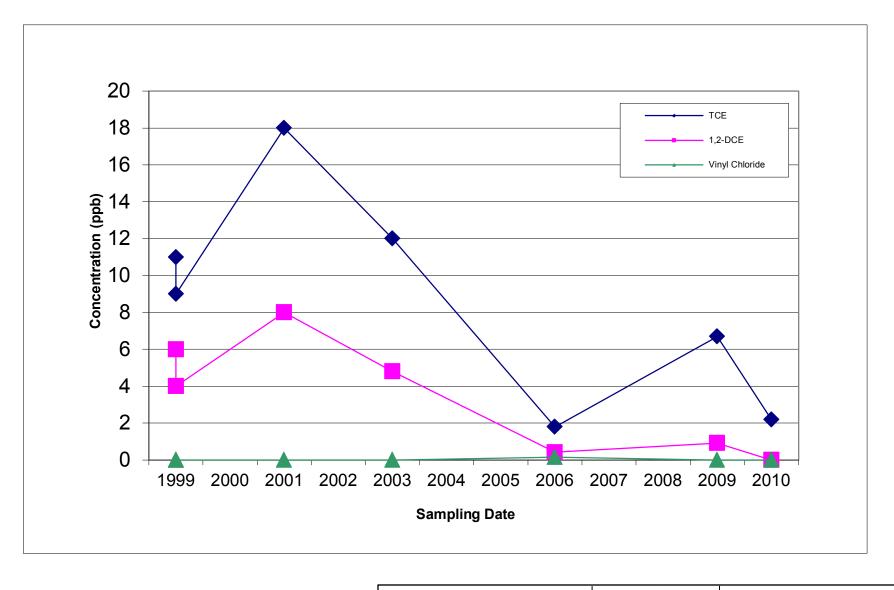
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MW-6 VOC TIME-CONCENTRATION PLOT

Project 128910, task 1000

December 2013



312 Fair Oak Street Little Valley, New York

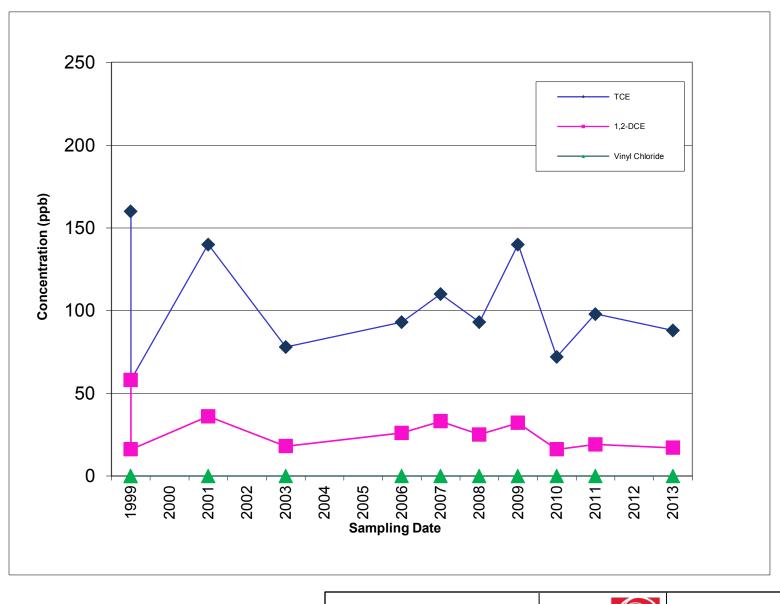
BUSH INDUSTRIES, INC.



MW-D1 VOC TIME-CONCENTRATION PLOT

Project 128910, task 1000 D

December 2013



312 Fair Oak Street Little Valley, New York

BUSH INDUSTRIES, INC.



MW-D2 VOC TIME-CONCENTRATION PLOT

Project 128910, task 1000

December 2013

Figure 10

Appendix A		
Data Validation Report		



DATA USABILITY SUMMARY REPORT for

Bush Industries

Analyses: Volatiles, Dissolved Gases, Ferrous Iron, Alkalinity, Chloride, Sulfate, Nitrate, Sulfide, Total Organic Carbon

SAMPLE DELIVERY GROUP 480-46182-1

PREPARED FOR:

GEI Consultants, Inc.

Reviewed by:

Reviewed by:

Approved by:

Prepared by

MEC^X 12269 East Vassar Drive Aurora, CO 80014



I. INTRODUCTION

Task Order Title: Bush Industries
Contract Task Order: 1427.001D.00
Sample Delivery Group: 480-46182-1
Project Manager: Kelly McIntosh

Matrix: Water QC Level: III

No. of Samples: 8
Reanalyses/Dilutions: 0

No. of Reanalyses/Dilutions: 0

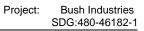
Laboratory: TestAmerica-Buffalo

Table 1. Sample Identification

Client ID	Laboratory ID	Matrix	Sample Date	Method
DUP	480-46182-7	Water	9/19/2013	300.0, 353.2, 2320B, 3500FE, 4500-SF, 8260B, 9060, RSK175
LVRA07-MNAGW-MW-2	480-46182-3	Water	9/19/2013	300.0, 353.2, 2320B, 3500FE, 4500-SF, 8260B, 9060, RSK175
LVRA07-MNAGW-MW-3	480-46182-1	Water	9/19/2013	300.0, 353.2, 2320B, 3500FE, 4500-SF, 8260B, 9060, RSK175
LVRA07-MNAGW-MW-5	480-46182-6	Water	9/19/2013	300.0, 353.2, 2320B, 3500FE, 4500-SF, 8260B, 9060, RSK175
LVRA07-MNAGW-MW-6	480-46182-5	Water	9/19/2013	300.0, 353.2, 2320B, 3500FE, 4500-SF, 8260B, 9060, RSK175
LVRA07-MNAGW-MW-D2	480-46182-2	Water	9/19/2013	300.0, 353.2, 2320B, 3500FE, 4500-SF, 8260B, 9060, RSK175
RINSE BLANK	480-46182-4	Water	9/19/2013	8260B, RSK175
TRIP BLANK	480-46182-8	Water	9/19/2013	8260B

II. Sample Management

No anomalies were observed regarding sample management. The samples in this SDG were received at the laboratory within the temperature limits of 4°C ±2°C. Changes to the COC were made by obliterating or overwriting the entry. The changes were not initialed or dated. The COC and sample receipt checklist noted the samples for the dissolved metals method were not field-filtered. The COC was appropriately signed and dated by field and/or laboratory personnel. Custody seals were present and intact on the cooler.





Data Qualifier Reference Table

Qualifie	r Organics	Inorganics
U	The analyte was analyzed for, but was not detected above the reported sample quantitation limit. The associated value is the quantitation limit or the estimated detection limit for dioxins or PCB congeners.	The material was analyzed for, but was not detected above the level of the associated value. The associated value is either the sample quantitation limit or the sample detection limit. The associated value is the sample detection limit or the quantitation limit for perchlorate only.
J	The analyte was positively identified; the associated numerical value is the approximate concentration of the analyte in the sample.	The associated value is an estimated quantity.
N	The analysis indicates the presence of an analyte for which there is presumptive evidence to make a "tentative identification."	Not applicable.
NJ	The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.	Not applicable.
UJ	The analyte was not deemed above the reported sample quantitation limit. However, the reported quantitation limit is approximate and may or may not represent the actual limit of quantitation necessary to accurately and precisely measure the analyte in the sample.	The material was analyzed for, but was not detected. The associated value is an estimate and may be inaccurate or imprecise.
R	The data are unusable. The sample results are rejected due to serious deficiencies in the ability to analyze the sample and to meet quality control criteria. The presence or absence of the analyte cannot be verified.	The data are unusable. The sample results are rejected due to serious deficiencies in the ability to analyze the sample and to meet quality control criteria. The presence or absence of the analyte cannot be verified.



Qualification Code Reference Table

Qualifier	Organics	Inorganics
Н	Holding times were exceeded.	Holding times were exceeded.
S	Surrogate recovery was outside QC limits.	The sequence or number of standards used for the calibration was incorrect
С	Calibration %RSD or %D was noncompliant.	Correlation coefficient is <0.995.
R	Calibration RRF was <0.05.	%R for calibration is not within control limits.
В	Presumed contamination as indicated by the preparation (method) blank results.	Presumed contamination as indicated by the preparation (method) or calibration blank results.
L	Laboratory Blank Spike/Blank Spike Duplicate %R was not within control limits.	Laboratory Control Sample %R was not within control limits.
Q	MS/MSD recovery was poor or RPD high.	MS recovery was poor.
Е	Not applicable.	Duplicates showed poor agreement.
I	Internal standard performance was unsatisfactory.	ICP ICS results were unsatisfactory.
Α	Not applicable.	ICP Serial Dilution %D were not within control limits.
M	Tuning (BFB or DFTPP) was noncompliant.	Not applicable.
Т	Presumed contamination as indicated by the trip blank results.	Not applicable.
+	False positive – reported compound was not present.	Not applicable.
-	False negative – compound was present but not reported.	Not applicable.
F	Presumed contamination as indicated by the FB or ER results.	Presumed contamination as indicated by the FB or ER results.
\$	Reported result or other information was incorrect.	Reported result or other information was incorrect.
?	TIC identity or reported retention time has been changed.	Not applicable.



found.

Qualification Code Reference Table Cont.

D	The analysis with this flag should not be used because another more technically sound analysis is available.	The analysis with this flag should not be used because another more technically sound analysis is available.
Р	Instrument performance for pesticides was poor.	Post Digestion Spike recovery was not within control limits.
*11, *111	Unusual problems found with the data that have been described in Section II, "Sample Management," or Section III, "Method Analyses." The	Unusual problems found with the data that have been described in Section II, "Sample Management," or Section III, "Method Analyses."

number following the asterisk (*) will

indicate the report section where a

description of the problem can be

4 Revision 0

The number following the asterisk

(*) will indicate the report section

can be found.

where a description of the problem



III. Method Analyses

A. EPA Method 8260B - Volatile Organic Compounds (VOCs)

Reviewed By: L. Calvin

Date Reviewed: December 10, 2013

The samples listed in Table 1 for this analysis were validated based on the guidelines outlined in the MEC^X Data Validation Procedure for Volatile Organics (DVP-2, Rev. 0), EPA Method 8260B, CLP Organics Data Review and Preliminary Review (9/2006), and the USEPA Hazardous Waste Support Branch Validating Volatile Organic Compounds by Gas Chromatography/Mass Spectrometry SW-846 Method 8260B (8/2008).

- Holding Times: The preserved water samples were analyzed within 14 days of collection.
- GC/MS Tuning: BFB tune ion abundance summaries were not provided, and were therefore not evaluated.
- Calibration: Initial and continuing calibration summaries were not provided; therefore, initial calibration average RRFs and %RSDs, and ICV and continuing calibration RRFs and %Ds were not evaluated.
- Blanks: The method blank had no target compound detects above the MDL.
- Blank Spikes and Laboratory Control Samples: A subset of thirteen target compounds was spiked in the LCS. All recoveries were within laboratory-established QC limits.
- Surrogate Recovery: The surrogate recoveries were within laboratory-established QC limits.
- Matrix Spike/Matrix Spike Duplicate: MS/MSD analyses were performed on sample LVRA07-MNAGW-MW-2. Chlorobenzene and tetrachloroethene were recovered marginally above the QC limits in both the MS and MSD; however, neither compound was detected in the parent sample; therefore, qualification was not required. Remaining recoveries and all RPDs were within laboratory-established QC limits.
- Field QC Samples: Field QC samples were evaluated, and if necessary, qualified based on method blanks and other laboratory QC results affecting the usability of the field QC data. Any remaining detects were used to evaluate the associated site samples. Following are findings associated with field QC samples:
 - Trip Blanks: Sample Trip Blank was the trip blank identified for the samples in this SDG. The trip blank had no detects reported above the MDL.



- Field Blanks and Equipment Rinsates: Sample Rinse Blank was identified as the equipment rinsate associated with the samples in this SDG. Acetone was detected in this sample but was not detected in the associated site samples. The equipment rinsate had no detects reported above the MDL.
- Field Duplicates: Samples LVRA07-MNAGW-MW-D2 and DUP were identified as the field duplicate pair in this SDG. There were common detects above the reporting limit for cis-1,2-dichloroethene and trichloroethene with calculated RPDs of 0% and 1.1%, respectively. The pair was considered to be in good agreement.
- Internal Standards Performance: The internal standard area counts and retention times for the samples were within the control limits established by the continuing calibration standards: -50%/+100% for internal standard areas and ±30 seconds for retention times.
- Compound Identification: Compound identification was not verified at this level of validation. The laboratory analyzed for volatiles by EPA Method 8260B.
- Compound Quantification: Compound quantitation was not verified at this level of validation. The reporting limits were supported by the laboratory MDLs. Any detect between the MDL and the reporting limit was qualified as estimated, "J," in the samples of this SDG. Reported nondetects are valid to the reporting limit.
- Tentatively Identified Compounds: TICs were not reported by the laboratory for this SDG.
- System Performance: System performance was not evaluated at this level of validation.

B. Method RSK-175-Methane, Ethane, Ethene

Reviewed By: L. Calvin

Date Reviewed: December 10, 2013

The samples listed in Table 1 for this analysis were validated based on the guidelines outlined in MEC^{x} Data Validation Procedure for Volatile Organics (DVP-2, Rev. 0), MEC^{x} Data Validation Procedure for Volatile Organics (DVP-2, Rev. 0), MEC^{x} Data Validation Procedure for Volatile Organics (DVP-2, Rev. 0), MEC^{x} Data Review and Preliminary Review (9/2006), and SW-846 ME-846 ME-847 ME-847 ME-847 ME-848 ME-849 ME

- Holding Times: The samples in the SDG were analyzed within 14 days of collection.
- GC/MS Tuning: Not applicable to this analysis.
- Calibration: Calibration criteria were met. Initial and continuing calibration summaries were not provided; therefore, initial calibration r² values, and ICV and CCV %Ds were not evaluated.



- Blanks: The method blank had no detects above the MDL.
- Blank Spikes and Laboratory Control Samples: Recoveries and RPDs were within the laboratory established QC limits.
- Surrogate Recovery: Surrogates were not utilized in this method.
- Matrix Spike/Matrix Spike Duplicate: MS/MSD analyses were performed on sample LVRA07-MNAGW-MW-2. The parent sample concentration of methane exceeded four times the spiked amount; therefore, methane recoveries were not evaluated. The remaining recoveries and RPDs were within the laboratory established QC limits.
- Compound Identification: Compound identification was not verified at this level of validation. The laboratory analyzed for methane, ethane, and ethene by EPA Method RSK-175.
- Compound Quantification and Reported Detection Limits: Compound quantification was not verified at this level of validation. The reporting limits were supported by the laboratory MDL. Any detect between the MDL and the reporting limit was qualified as estimated, "J," in the samples of this SDG. Reported nondetects are valid to the reporting limit.
- System Performance: System performance was not evaluated at this level of validation.
- Field QC Samples: Field QC samples were evaluated, and if necessary, qualified based on method blanks and other laboratory QC results affecting the usability of the field QC data. Any remaining detects were used to evaluate the associated site samples. Following are findings associated with field QC samples:
 - Trip Blanks: This SDG had no identified trip blank for this analysis.
 - Field Blanks and Equipment Rinsates: Sample Rinse Blank was identified as the equipment rinsate associated with the samples in this SDG. The equipment rinsate had no reported detects above the MDL.
 - DUP were identified as the field duplicate pair in this SDG. Neither sample had reported detects above the MDL. The pair was considered to be in good agreement.



C. VARIOUS EPA METHODS—General Minerals

Reviewed By: P. Meeks

Date Reviewed: December 10, 2013

The samples listed in Table 1 for this analysis were validated based on the guidelines outlined in the MEC^x Data Validation Procedure for General Minerals (DVP-6, Rev. 0), EPA Methods 300.0, 353.2, 2320B, 3500FE, 4500-SF, and 9060, and the Validation of Metals for the Contract Laboratory Program based on SOW ILMO5.3, SOP Revision 13 (9/2006).

- Holding Times: The analytical holding times, 28 days from collection for chloride, sulfate
 and TOC, 14 days from collection for alkalinity, seven days from collection for sulfide, and
 48 hours from collection for nitrate, were met. As per the method, the analytical holding
 time for ferrous iron is noted as "in field". As the ferrous iron analyses were performed
 within 24 hours of receipt at the laboratory, no qualifications were required.
- Calibration: Initial calibration, ICV and CCVs were not assessed as the laboratory did not provide summary or raw data.
- Blanks: Method blanks had no detects. CCBs were not assessed as the laboratory did not provide CCB summary data.
- Blank Spikes and Laboratory Control Samples: Recoveries and RPDs were within laboratory-established QC limits.
- Laboratory Duplicates: A laboratory duplicate analysis was performed on LVRA07-MNAGW-MW-2 for sulfide. The RPD was within the laboratory-established QC limit.
- Matrix Spike/Matrix Spike Duplicate: MS/MSD analyses were performed on LVRA07-MNAGW-MW-2 for all analytes. All recoveries and RPDs were within the laboratoryestablished control limits.
- Sample Result Verification: Compound identification was not verified at this level of validation. The sample result summaries were compared to the raw data and no transcription errors were noted.
- Field QC Samples: Field QC samples were evaluated, and if necessary, qualified based on method blanks and other laboratory QC results affecting the usability of the field QC data. Any remaining detects were used to evaluate the associated site samples.
 Following are findings associated with field QC samples:
 - Field Blanks and Equipment Rinsates: This SDG had no identified field blank or equipment rinsate samples for these analyses.



 Field Duplicates: Field Duplicates: Samples LVRA07-MNAGW-MW-D2 and DUP were identified as the field duplicate pair in this SDG. The samples were considered to be in good agreement as all detects were in common and all RPDs were less than 30%.

Validated Sample Result Forms: 480-46182-1

	<u> </u>					
Analysis Metho						
Sample Name:	DUP			Matrix:	Water Resu	ılt Type: FS
Lab Sample Name:	480-46182-7	Sample l	Date: 09/1	9/2013	V	alidation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier
Chloride	16887-00-6	20.2	0.28	MG/L		
Sulfate	14808-79-8	14.2	0.35	MG/L		
Sample Name:	LVRA07-MNA	GW-MW-2		Matrix:	Water Resu	ılt Type: FS
Lab Sample Name:	480-46182-3	Sample l	Date: 09/1	9/2013	v	alidation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier
Chloride	16887-00-6	16.7	0.28	MG/L		
Sulfate	14808-79-8	17.5	0.35	MG/L		
Sample Name:	LVRA07-MNA	GW-MW-3		Matrix:	Water Resu	ult Type: FS
Lab Sample Name:	480-46182-1	Sample l	Date: 09/1	9/2013	v	alidation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier
Chloride	16887-00-6	20.3	0.28	MG/L		
Sulfate	14808-79-8	10.6	0.35	MG/L		
Sample Name:	LVRA07-MNA	GW-MW-5		Matrix:	Water Resu	ılt Type: FS
Lab Sample Name:	480-46182-6	Sample l	Date: 09/1	9/2013	V	alidation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier
Chloride	16887-00-6	14.2	0.28	MG/L		
Sulfate	14808-79-8	5.4	0.35	MG/L		
Sample Name:	LVRA07-MNA	GW-MW-6		Matrix:	Water Resu	ılt Type: FS
Lab Sample Name:	480-46182-5	Sample l	Date: 09/1	9/2013	V	alidation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier
Chloride	16887-00-6	10.9	0.28	MG/L		
Sulfate	14808-79-8	10.6	0.35	MG/L		
)2.	Matrix:	Water Resu	ılt Type: FS
Sample Name:	LVRA07-MNA	GW-MW-L	_			
Sample Name: Lab Sample Name:	LVRA07-MNA 480-46182-2		Date: 09/1	9/2013	V	alidation Level: III
-				9/2013 Result Units	V Lab Qualifier	Validation Level: III Validation Validation Notes Qualifier
Lab Sample Name:	480-46182-2	Sample l	Date: 09/1	Result	Lab	Validation Validation Notes

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Analysis Method 353.2

Sample Name:	DUP			Matrix:	Water Result Type: FS
Lab Sample Name:	480-46182-7	Sample 1	Date: 09/1	9/2013	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Validation Validation Notes Qualifier Qualifier
Nitrate	14797-55-8	0.22	0.020	MG/L A	
Sample Name:	LVRA07-MNA	GW-MW-2		Matrix:	Water Result Type: FS
Lab Sample Name:	480-46182-3	Sample l	Date: 09/1	9/2013	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Validation Validation Notes Qualifier Qualifier
Nitrate	14797-55-8	0.052	0.020	MG/L A	
Sample Name:	LVRA07-MNA	GW-MW-3		Matrix:	Water Result Type: FS
Lab Sample Name:	480-46182-1	Sample 1	Date: 09/1	9/2013	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Validation Validation Notes Qualifier Qualifier
Nitrate	14797-55-8	1.2	0.020	MG/L A	
Sample Name:	LVRA07-MNA	GW-MW-5		Matrix:	Water Result Type: FS
Lab Sample Name:	480-46182-6	Sample 1	Date: 09/1	9/2013	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Validation Validation Notes Qualifier Qualifier
Nitrate	14797-55-8	0.051	0.020	MG/L A	
Sample Name:	LVRA07-MNA	GW-MW-6		Matrix:	Water Result Type: FS
Lab Sample Name:	480-46182-5	Sample l	Date: 09/1	9/2013	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Validation Validation Notes Qualifier Qualifier
Nitrate	14797-55-8	0.034	0.020	MG/L A	J J
Sample Name:	LVRA07-MNA	GW-MW-D	2	Matrix:	Water Result Type: FS
Lab Sample Name:	480-46182-2	Sample l	Date: 09/1	9/2013	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Validation Validation Notes Qualifier Qualifier
Nitrate	14797-55-8	0.22	0.020	MG/L A	

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Sample Name: DUP Matrix: Water Result Type: FS

Lab Sample Name: 480-46182-7 Sample Date: 09/19/2013 Validation Level: III

Lab Sample Name: 48	30-46182-7	Sample Date: 09/19/2013			7	vel: III	
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Qualifier	Validation Notes
1,1,1-Trichloroethane	71-55-6	0.82	0.82	UG/L	U	U	
1,1,2,2-Tetrachloroethane	79-34-5	0.21	0.21	UG/L	U	U	
1,1,2-Trichloro-1,2,2- trifluoroethane	76-13-1	0.31	0.31	UG/L	U	U	
1,1,2-Trichloroethane	79-00-5	0.23	0.23	UG/L	U	U	
1,1-Dichloroethane	75-34-3	0.38	0.38	UG/L	U	U	
1,1-Dichloroethene	75-35-4	0.29	0.29	UG/L	U	U	
1,2,4-Trichlorobenzene	120-82-1	0.41	0.41	UG/L	U	U	
1,2-Dibromo-3-Chloropropane	96-12-8	0.39	0.39	UG/L	U	U	
1,2-Dibromoethane	106-93-4	0.73	0.73	UG/L	U	U	
1,2-Dichlorobenzene	95-50-1	0.79	0.79	UG/L	U	U	
1,2-Dichloroethane	107-06-2	0.21	0.21	UG/L	U	U	
1,2-Dichloropropane	78-87-5	0.72	0.72	UG/L	U	U	
1,3-Dichlorobenzene	541-73-1	0.78	0.78	UG/L	U	U	
1,4-Dichlorobenzene	106-46-7	0.84	0.84	UG/L	U	U	
2-Butanone	78-93-3	1.3	1.3	UG/L	U	U	
2-Hexanone	591-78-6	1.2	1.2	UG/L	U	U	
4-Methyl-2-pentanone	108-10-1	2.1	2.1	UG/L	U	U	
Acetone	67-64-1	3.0	3.0	UG/L	U	U	
Benzene	71-43-2	0.41	0.41	UG/L	U	U	
Bromodichloromethane	75-27-4	0.39	0.39	UG/L	U	U	
Bromoform	75-25-2	0.26	0.26	UG/L	U	U	
Bromomethane	74-83-9	0.69	0.69	UG/L	U	U	
Carbon disulfide	75-15-0	0.19	0.19	UG/L	U	U	
Carbon tetrachloride	56-23-5	0.27	0.27	UG/L	U	U	
Chlorobenzene	108-90-7	0.75	0.75	UG/L	U	U	
Chloroethane	75-00-3	0.32	0.32	UG/L	U	U	
Chloroform	67-66-3	0.34	0.34	UG/L	U	U	
Chloromethane	74-87-3	0.35	0.35	UG/L	U	U	
cis-1,2-Dichloroethene	156-59-2	17	0.81	UG/L			
cis-1,3-Dichloropropene	10061-01-5	0.36	0.36	UG/L	U	U	
Cyclohexane	110-82-7	0.18	0.18	UG/L	U	U	
Dibromochloromethane	124-48-1	0.32	0.32	UG/L	U	U	
Dichlorodifluoromethane	75-71-8	0.68	0.68	UG/L	U	U	
Ethylbenzene	100-41-4	0.74	0.74	UG/L	U	U	
Isopropylbenzene	98-82-8	0.79	0.79	UG/L	U	U	
Methyl acetate	79-20-9	0.50	0.50	UG/L	U	U	
Methyl tert-butyl ether	1634-04-4	0.16	0.16	UG/L	U	U	

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Analys	is Method	8260B
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Methylcyclohexane	108-87-2	0.16	0.16	UG/L	U	U	
Methylene Chloride	75-09-2	0.44	0.44	UG/L	U	Ŭ	
Styrene	100-42-5	0.73	0.73	UG/L	U	U	
Tetrachloroethene	127-18-4	0.36	0.36	UG/L	U	U	
Toluene	108-88-3	0.51	0.51	UG/L	U	U	
trans-1,2-Dichloroethene	156-60-5	0.90	0.90	UG/L	U	U	
trans-1,3-Dichloropropene	10061-02-6	0.37	0.37	UG/L	U	U	
Trichloroethene	79-01-6	87	0.46	UG/L			
Trichlorofluoromethane	75-69-4	0.88	0.88	UG/L	U	U	
Vinyl chloride	75-01-4	0.90	0.90	UG/L	U	U	
Xylenes, Total	1330-20-7	0.66	0.66	UG/L	U	U	

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Sample Name: LVRA07-MNAGW-MW-2 Matrix: Water Result Type: FS

Lab Sample Name: 480-46182-3 Sample Date: 09/19/2013 Validation Level: III

Lab Sample Name:	180-46182-3	Sample Date: 09/19/2013			Validation Level: III		
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Qualifier	Validation Notes
1,1,1-Trichloroethane	71-55-6	0.82	0.82	UG/L	U	U	
1,1,2,2-Tetrachloroethane	79-34-5	0.21	0.21	UG/L	U	U	
1,1,2-Trichloro-1,2,2- trifluoroethane	76-13-1	0.31	0.31	UG/L	U	U	
1,1,2-Trichloroethane	79-00-5	0.23	0.23	UG/L	U	U	
1,1-Dichloroethane	75-34-3	0.38	0.38	UG/L	U	U	
1,1-Dichloroethene	75-35-4	0.57	0.29	UG/L	J	J	
1,2,4-Trichlorobenzene	120-82-1	0.41	0.41	UG/L	U	U	
1,2-Dibromo-3-Chloropropane	96-12-8	0.39	0.39	UG/L	U	U	
1,2-Dibromoethane	106-93-4	0.73	0.73	UG/L	U	U	
1,2-Dichlorobenzene	95-50-1	0.79	0.79	UG/L	U	U	
1,2-Dichloroethane	107-06-2	0.21	0.21	UG/L	U	U	
1,2-Dichloropropane	78-87-5	0.72	0.72	UG/L	U	U	
1,3-Dichlorobenzene	541-73-1	0.78	0.78	UG/L	U	U	
1,4-Dichlorobenzene	106-46-7	0.84	0.84	UG/L	U	U	
2-Butanone	78-93-3	1.3	1.3	UG/L	U	U	
2-Hexanone	591-78-6	1.2	1.2	UG/L	U	U	
4-Methyl-2-pentanone	108-10-1	2.1	2.1	UG/L	U	U	
Acetone	67-64-1	3.0	3.0	UG/L	U	U	
Benzene	71-43-2	0.41	0.41	UG/L	U	U	
Bromodichloromethane	75-27-4	0.39	0.39	UG/L	U	U	
Bromoform	75-25-2	0.26	0.26	UG/L	U	U	
Bromomethane	74-83-9	0.69	0.69	UG/L	U	U	
Carbon disulfide	75-15-0	0.19	0.19	UG/L	U	U	
Carbon tetrachloride	56-23-5	0.27	0.27	UG/L	U	U	
Chlorobenzene	108-90-7	0.75	0.75	UG/L	U	U	
Chloroethane	75-00-3	0.32	0.32	UG/L	U	U	
Chloroform	67-66-3	0.34	0.34	UG/L	U	U	
Chloromethane	74-87-3	0.35	0.35	UG/L	U	U	
cis-1,2-Dichloroethene	156-59-2	28	0.81	UG/L			
cis-1,3-Dichloropropene	10061-01-5	0.36	0.36	UG/L	U	U	
Cyclohexane	110-82-7	0.18	0.18	UG/L	U	U	
Dibromochloromethane	124-48-1	0.32	0.32	UG/L	U	U	
Dichlorodifluoromethane	75-71-8	0.68	0.68	UG/L	U	U	
Ethylbenzene	100-41-4	0.74	0.74	UG/L	U	U	
Isopropylbenzene	98-82-8	0.79	0.79	UG/L	U	U	
Methyl acetate	79-20-9	0.50	0.50	UG/L	U	U	
Methyl tert-butyl ether	1634-04-4	0.16	0.16	UG/L	U	U	

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Analys	is Method	8260B
11.0000,0	is in contour	02000

Methylcyclohexane	108-87-2	0.16	0.16	UG/L	U	U	
Methylene Chloride	75-09-2	0.44	0.44	UG/L	U	U	
Styrene	100-42-5	0.73	0.73	UG/L	U	U	
Tetrachloroethene	127-18-4	0.36	0.36	UG/L	U	U	
Toluene	108-88-3	0.51	0.51	UG/L	U	U	
trans-1,2-Dichloroethene	156-60-5	0.90	0.90	UG/L	U	U	
trans-1,3-Dichloropropene	10061-02-6	0.37	0.37	UG/L	U	U	
Trichloroethene	79-01-6	61	0.46	UG/L			
Trichlorofluoromethane	75-69-4	0.88	0.88	UG/L	U	U	
Vinyl chloride	75-01-4	1.3	0.90	UG/L			
Xylenes, Total	1330-20-7	0.66	0.66	UG/L	U	U	

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Sample Name: LVRA07-MNAGW-MW-3 Matrix: Water Result Type: FS

Lab Sample Name: 480-46182-1 Sample Date: 09/19/2013 Validation Level: III

Lab Sample Name: 48	480-46182-1 Sample Date: 09/19/2013			9/2013	Validation Level: III			
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Qualifier	Validation Notes	
1,1,1-Trichloroethane	71-55-6	0.82	0.82	UG/L	U	U		
1,1,2,2-Tetrachloroethane	79-34-5	0.21	0.21	UG/L	U	U		
1,1,2-Trichloro-1,2,2- trifluoroethane	76-13-1	0.31	0.31	UG/L	U	U		
1,1,2-Trichloroethane	79-00-5	0.23	0.23	UG/L	U	U		
1,1-Dichloroethane	75-34-3	0.38	0.38	UG/L	U	U		
1,1-Dichloroethene	75-35-4	0.29	0.29	UG/L	U	U		
1,2,4-Trichlorobenzene	120-82-1	0.41	0.41	UG/L	U	U		
1,2-Dibromo-3-Chloropropane	96-12-8	0.39	0.39	UG/L	U	U		
1,2-Dibromoethane	106-93-4	0.73	0.73	UG/L	U	U		
1,2-Dichlorobenzene	95-50-1	0.79	0.79	UG/L	U	U		
1,2-Dichloroethane	107-06-2	0.21	0.21	UG/L	U	U		
1,2-Dichloropropane	78-87-5	0.72	0.72	UG/L	U	U		
1,3-Dichlorobenzene	541-73-1	0.78	0.78	UG/L	U	U		
1,4-Dichlorobenzene	106-46-7	0.84	0.84	UG/L	U	U		
2-Butanone	78-93-3	1.3	1.3	UG/L	U	U		
2-Hexanone	591-78-6	1.2	1.2	UG/L	U	U		
4-Methyl-2-pentanone	108-10-1	2.1	2.1	UG/L	U	U		
Acetone	67-64-1	3.0	3.0	UG/L	U	U		
Benzene	71-43-2	0.41	0.41	UG/L	U	U		
Bromodichloromethane	75-27-4	0.39	0.39	UG/L	U	U		
Bromoform	75-25-2	0.26	0.26	UG/L	U	U		
Bromomethane	74-83-9	0.69	0.69	UG/L	U	U		
Carbon disulfide	75-15-0	0.19	0.19	UG/L	U	U		
Carbon tetrachloride	56-23-5	0.27	0.27	UG/L	U	U		
Chlorobenzene	108-90-7	0.75	0.75	UG/L	U	U		
Chloroethane	75-00-3	0.32	0.32	UG/L	U	U		
Chloroform	67-66-3	0.34	0.34	UG/L	U	U		
Chloromethane	74-87-3	0.35	0.35	UG/L	U	U		
cis-1,2-Dichloroethene	156-59-2	0.81	0.81	UG/L	U	U		
cis-1,3-Dichloropropene	10061-01-5	0.36	0.36	UG/L	U	U		
Cyclohexane	110-82-7	0.18	0.18	UG/L	U	U		
Dibromochloromethane	124-48-1	0.32	0.32	UG/L	U	U		
Dichlorodifluoromethane	75-71-8	0.68	0.68	UG/L	U	U		
Ethylbenzene	100-41-4	0.74	0.74	UG/L	U	U		
Isopropylbenzene	98-82-8	0.79	0.79	UG/L	U	U		
Methyl acetate	79-20-9	0.50	0.50	UG/L	U	U		
Methyl tert-butyl ether	1634-04-4	0.16	0.16	UG/L	U	U		

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Methylcyclohexane	108-87-2	0.16	0.16	UG/L	U	U	
Methylene Chloride	75-09-2	0.44	0.44	UG/L	U	U	
Styrene	100-42-5	0.73	0.73	UG/L	U	U	
Tetrachloroethene	127-18-4	0.36	0.36	UG/L	U	U	
Toluene	108-88-3	0.51	0.51	UG/L	U	U	
trans-1,2-Dichloroethene	156-60-5	0.90	0.90	UG/L	U	U	
trans-1,3-Dichloropropene	10061-02-6	0.37	0.37	UG/L	U	U	
Trichloroethene	79-01-6	3.9	0.46	UG/L			
Trichlorofluoromethane	75-69-4	0.88	0.88	UG/L	U	U	
Vinyl chloride	75-01-4	0.90	0.90	UG/L	U	U	
Xylenes, Total	1330-20-7	0.66	0.66	UG/L	U	U	

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Sample Name: LVRA07-MNAGW-MW-5 Matrix: Water Result Type: FS

Lab Sample Name: 480-46182-6 Sample Date: 09/19/2013 Validation Level: III

Lab Sample Name: 48	30-46182-6	Sample Date: 09/19/2013			7	evel: III	
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Qualifier	Validation Notes
1,1,1-Trichloroethane	71-55-6	0.82	0.82	UG/L	U	U	
1,1,2,2-Tetrachloroethane	79-34-5	0.21	0.21	UG/L	U	U	
1,1,2-Trichloro-1,2,2-trifluoroethane	76-13-1	0.31	0.31	UG/L	U	U	
1,1,2-Trichloroethane	79-00-5	0.23	0.23	UG/L	U	U	
1,1-Dichloroethane	75-34-3	0.38	0.38	UG/L	U	U	
1,1-Dichloroethene	75-35-4	0.29	0.29	UG/L	U	U	
1,2,4-Trichlorobenzene	120-82-1	0.41	0.41	UG/L	U	U	
1,2-Dibromo-3-Chloropropane	96-12-8	0.39	0.39	UG/L	U	U	
1,2-Dibromoethane	106-93-4	0.73	0.73	UG/L	U	U	
1,2-Dichlorobenzene	95-50-1	0.79	0.79	UG/L	U	U	
1,2-Dichloroethane	107-06-2	0.21	0.21	UG/L	U	U	
1,2-Dichloropropane	78-87-5	0.72	0.72	UG/L	U	U	
1,3-Dichlorobenzene	541-73-1	0.78	0.78	UG/L	U	U	
1,4-Dichlorobenzene	106-46-7	0.84	0.84	UG/L	U	U	
2-Butanone	78-93-3	1.3	1.3	UG/L	U	U	
2-Hexanone	591-78-6	1.2	1.2	UG/L	U	U	
4-Methyl-2-pentanone	108-10-1	2.1	2.1	UG/L	U	U	
Acetone	67-64-1	3.0	3.0	UG/L	U	U	
Benzene	71-43-2	0.41	0.41	UG/L	U	U	
Bromodichloromethane	75-27-4	0.39	0.39	UG/L	U	U	
Bromoform	75-25-2	0.26	0.26	UG/L	U	U	
Bromomethane	74-83-9	0.69	0.69	UG/L	U	U	
Carbon disulfide	75-15-0	0.19	0.19	UG/L	U	U	
Carbon tetrachloride	56-23-5	0.27	0.27	UG/L	U	U	
Chlorobenzene	108-90-7	0.75	0.75	UG/L	U	U	
Chloroethane	75-00-3	0.32	0.32	UG/L	U	U	
Chloroform	67-66-3	0.34	0.34	UG/L	U	U	
Chloromethane	74-87-3	0.35	0.35	UG/L	U	U	
cis-1,2-Dichloroethene	156-59-2	0.81	0.81	UG/L	U	U	
cis-1,3-Dichloropropene	10061-01-5	0.36	0.36	UG/L	U	U	
Cyclohexane	110-82-7	0.18	0.18	UG/L	U	U	
Dibromochloromethane	124-48-1	0.32	0.32	UG/L	U	U	
Dichlorodifluoromethane	75-71-8	0.68	0.68	UG/L	U	U	
Ethylbenzene	100-41-4	0.74	0.74	UG/L	U	U	
Isopropylbenzene	98-82-8	0.79	0.79	UG/L	U	U	
Methyl acetate	79-20-9	0.50	0.50	UG/L	U	U	
Methyl tert-butyl ether	1634-04-4	0.16	0.16	UG/L	U	U	

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Methylcyclohexane	108-87-2	0.16	0.16	UG/L	U	U	
Methylene Chloride	75-09-2	0.44	0.44	UG/L	U	U	
Styrene	100-42-5	0.73	0.73	UG/L	U	U	
Tetrachloroethene	127-18-4	0.36	0.36	UG/L	U	U	
Toluene	108-88-3	0.51	0.51	UG/L	U	U	
trans-1,2-Dichloroethene	156-60-5	0.90	0.90	UG/L	U	U	
trans-1,3-Dichloropropene	10061-02-6	0.37	0.37	UG/L	U	U	
Trichloroethene	79-01-6	0.46	0.46	UG/L	U	U	
Trichlorofluoromethane	75-69-4	0.88	0.88	UG/L	U	U	
Vinyl chloride	75-01-4	0.90	0.90	UG/L	U	U	
Xylenes, Total	1330-20-7	0.66	0.66	UG/L	U	U	

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Sample Name: LVRA07-MNAGW-MW-6 Matrix: Water Result Type: FS

Lab Sample Name: 480-46182-5 Sample Date: 09/19/2013 Validation Level: III

Lab Sample Name: 4	80-46182-5	Sample Date: 09/19/2013			7	vel: III	
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Qualifier	Validation Notes
1,1,1-Trichloroethane	71-55-6	0.82	0.82	UG/L	U	U	
1,1,2,2-Tetrachloroethane	79-34-5	0.21	0.21	UG/L	U	U	
1,1,2-Trichloro-1,2,2- trifluoroethane	76-13-1	0.31	0.31	UG/L	U	U	
1,1,2-Trichloroethane	79-00-5	0.23	0.23	UG/L	U	U	
1,1-Dichloroethane	75-34-3	0.38	0.38	UG/L	U	U	
1,1-Dichloroethene	75-35-4	0.29	0.29	UG/L	U	U	
1,2,4-Trichlorobenzene	120-82-1	0.41	0.41	UG/L	U	U	
1,2-Dibromo-3-Chloropropane	96-12-8	0.39	0.39	UG/L	U	U	
1,2-Dibromoethane	106-93-4	0.73	0.73	UG/L	U	U	
1,2-Dichlorobenzene	95-50-1	0.79	0.79	UG/L	U	U	
1,2-Dichloroethane	107-06-2	0.21	0.21	UG/L	U	U	
1,2-Dichloropropane	78-87-5	0.72	0.72	UG/L	U	U	
1,3-Dichlorobenzene	541-73-1	0.78	0.78	UG/L	U	U	
1,4-Dichlorobenzene	106-46-7	0.84	0.84	UG/L	U	U	
2-Butanone	78-93-3	1.3	1.3	UG/L	U	U	
2-Hexanone	591-78-6	1.2	1.2	UG/L	U	U	
4-Methyl-2-pentanone	108-10-1	2.1	2.1	UG/L	U	U	
Acetone	67-64-1	3.0	3.0	UG/L	U	U	
Benzene	71-43-2	0.41	0.41	UG/L	U	U	
Bromodichloromethane	75-27-4	0.39	0.39	UG/L	U	U	
Bromoform	75-25-2	0.26	0.26	UG/L	U	U	
Bromomethane	74-83-9	0.69	0.69	UG/L	U	U	
Carbon disulfide	75-15-0	0.19	0.19	UG/L	U	U	
Carbon tetrachloride	56-23-5	0.27	0.27	UG/L	U	U	
Chlorobenzene	108-90-7	0.75	0.75	UG/L	U	U	
Chloroethane	75-00-3	0.32	0.32	UG/L	U	U	
Chloroform	67-66-3	0.34	0.34	UG/L	U	U	
Chloromethane	74-87-3	0.35	0.35	UG/L	U	U	
cis-1,2-Dichloroethene	156-59-2	14	0.81	UG/L			
cis-1,3-Dichloropropene	10061-01-5	0.36	0.36	UG/L	U	U	
Cyclohexane	110-82-7	0.18	0.18	UG/L	U	U	
Dibromochloromethane	124-48-1	0.32	0.32	UG/L	U	U	
Dichlorodifluoromethane	75-71-8	0.68	0.68	UG/L	U	U	
Ethylbenzene	100-41-4	0.74	0.74	UG/L	U	U	
Isopropylbenzene	98-82-8	0.79	0.79	UG/L	U	U	
Methyl acetate	79-20-9	0.50	0.50	UG/L	U	U	
Methyl tert-butyl ether	1634-04-4	0.16	0.16	UG/L	U	U	

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Methylcyclohexane	108-87-2	0.16	0.16	UG/L	U	U	
Methylene Chloride	75-09-2	0.44	0.44	UG/L	U	U	
Styrene	100-42-5	0.73	0.73	UG/L	U	U	
Tetrachloroethene	127-18-4	0.36	0.36	UG/L	U	U	
Toluene	108-88-3	0.51	0.51	UG/L	U	U	
trans-1,2-Dichloroethene	156-60-5	0.90	0.90	UG/L	U	U	
trans-1,3-Dichloropropene	10061-02-6	0.37	0.37	UG/L	U	U	
Trichloroethene	79-01-6	2.2	0.46	UG/L			
Trichlorofluoromethane	75-69-4	0.88	0.88	UG/L	U	U	
Vinyl chloride	75-01-4	1.0	0.90	UG/L			
Xylenes, Total	1330-20-7	0.66	0.66	UG/L	U	U	

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Sample Name: LVRA07-MNAGW-MW-D2 Matrix: Water Result Type: FS

Lab Sample Name: 480-46182-2 Sample Date: 09/19/2013 Validation Level: III

Lab Sample Name: 4	80-46182-2	Sample Date: 09/19/2013			7	vel: III	
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Qualifier	Validation Notes
1,1,1-Trichloroethane	71-55-6	0.82	0.82	UG/L	U	U	
1,1,2,2-Tetrachloroethane	79-34-5	0.21	0.21	UG/L	U	U	
1,1,2-Trichloro-1,2,2- trifluoroethane	76-13-1	0.31	0.31	UG/L	U	U	
1,1,2-Trichloroethane	79-00-5	0.23	0.23	UG/L	U	U	
1,1-Dichloroethane	75-34-3	0.38	0.38	UG/L	U	U	
1,1-Dichloroethene	75-35-4	0.29	0.29	UG/L	U	U	
1,2,4-Trichlorobenzene	120-82-1	0.41	0.41	UG/L	U	U	
1,2-Dibromo-3-Chloropropane	96-12-8	0.39	0.39	UG/L	U	U	
1,2-Dibromoethane	106-93-4	0.73	0.73	UG/L	U	U	
1,2-Dichlorobenzene	95-50-1	0.79	0.79	UG/L	U	U	
1,2-Dichloroethane	107-06-2	0.21	0.21	UG/L	U	U	
1,2-Dichloropropane	78-87-5	0.72	0.72	UG/L	U	U	
1,3-Dichlorobenzene	541-73-1	0.78	0.78	UG/L	U	U	
1,4-Dichlorobenzene	106-46-7	0.84	0.84	UG/L	U	U	
2-Butanone	78-93-3	1.3	1.3	UG/L	U	U	
2-Hexanone	591-78-6	1.2	1.2	UG/L	U	U	
4-Methyl-2-pentanone	108-10-1	2.1	2.1	UG/L	U	U	
Acetone	67-64-1	3.0	3.0	UG/L	U	U	
Benzene	71-43-2	0.41	0.41	UG/L	U	U	
Bromodichloromethane	75-27-4	0.39	0.39	UG/L	U	U	
Bromoform	75-25-2	0.26	0.26	UG/L	U	U	
Bromomethane	74-83-9	0.69	0.69	UG/L	U	U	
Carbon disulfide	75-15-0	0.19	0.19	UG/L	U	U	
Carbon tetrachloride	56-23-5	0.27	0.27	UG/L	U	U	
Chlorobenzene	108-90-7	0.75	0.75	UG/L	U	U	
Chloroethane	75-00-3	0.32	0.32	UG/L	U	U	
Chloroform	67-66-3	0.34	0.34	UG/L	U	U	
Chloromethane	74-87-3	0.35	0.35	UG/L	U	U	
cis-1,2-Dichloroethene	156-59-2	17	0.81	UG/L			
cis-1,3-Dichloropropene	10061-01-5	0.36	0.36	UG/L	U	U	
Cyclohexane	110-82-7	0.18	0.18	UG/L	U	U	
Dibromochloromethane	124-48-1	0.32	0.32	UG/L	U	U	
Dichlorodifluoromethane	75-71-8	0.68	0.68	UG/L	U	U	
Ethylbenzene	100-41-4	0.74	0.74	UG/L	U	U	
Isopropylbenzene	98-82-8	0.79	0.79	UG/L	U	U	
Methyl acetate	79-20-9	0.50	0.50	UG/L	U	Ū	
Methyl tert-butyl ether	1634-04-4	0.16	0.16	UG/L	U	U	

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Methylcyclohexane	108-87-2	0.16	0.16	UG/L	U	U	
Methylene Chloride	75-09-2	0.44	0.44	UG/L	U	U	
Styrene	100-42-5	0.73	0.73	UG/L	U	U	
Tetrachloroethene	127-18-4	0.36	0.36	UG/L	U	U	
Toluene	108-88-3	0.51	0.51	UG/L	U	U	
trans-1,2-Dichloroethene	156-60-5	0.90	0.90	UG/L	U	U	
trans-1,3-Dichloropropene	10061-02-6	0.37	0.37	UG/L	U	U	
Trichloroethene	79-01-6	88	0.46	UG/L			
Trichlorofluoromethane	75-69-4	0.88	0.88	UG/L	U	U	
Vinyl chloride	75-01-4	0.90	0.90	UG/L	U	U	
Xylenes, Total	1330-20-7	0.66	0.66	UG/L	U	U	

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Sample Name:	RINSE BLANK			Matrix:	Water Resi	ult Type: FS)
Lab Sample Name:	480-46182-4	Sample 1	Sample Date: 09/19/2013			Validation Le	evel: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Qualifier	Validation Notes
1,1,1-Trichloroethane	71-55-6	0.82	0.82	UG/L	U	U	
1,1,2,2-Tetrachloroethane	79-34-5	0.21	0.21	UG/L	U	U	
1,1,2-Trichloro-1,2,2- trifluoroethane	76-13-1	0.31	0.31	UG/L	U	U	
1,1,2-Trichloroethane	79-00-5	0.23	0.23	UG/L	U	U	
1,1-Dichloroethane	75-34-3	0.38	0.38	UG/L	U	U	
1,1-Dichloroethene	75-35-4	0.29	0.29	UG/L	U	U	
1,2,4-Trichlorobenzene	120-82-1	0.41	0.41	UG/L	U	U	
1,2-Dibromo-3-Chloropropan	e 96-12-8	0.39	0.39	UG/L	U	U	
,2-Dibromoethane	106-93-4	0.73	0.73	UG/L	U	U	
,2-Dichlorobenzene	95-50-1	0.79	0.79	UG/L	U	U	
,2-Dichloroethane	107-06-2	0.21	0.21	UG/L	U	U	
,2-Dichloropropane	78-87-5	0.72	0.72	UG/L	U	U	
,3-Dichlorobenzene	541-73-1	0.78	0.78	UG/L	U	U	
,4-Dichlorobenzene	106-46-7	0.84	0.84	UG/L	U	U	
-Butanone	78-93-3	1.3	1.3	UG/L	U	U	
-Hexanone	591-78-6	1.2	1.2	UG/L	U	U	
-Methyl-2-pentanone	108-10-1	2.1	2.1	UG/L	U	U	
cetone	67-64-1	3.0	3.0	UG/L	U	U	
enzene	71-43-2	0.41	0.41	UG/L	U	U	
romodichloromethane	75-27-4	0.39	0.39	UG/L	U	U	
romoform	75-25-2	0.26	0.26	UG/L	U	U	
Fromomethane	74-83-9	0.69	0.69	UG/L	U	U	
arbon disulfide	75-15-0	0.19	0.19	UG/L	U	U	
arbon tetrachloride	56-23-5	0.27	0.27	UG/L	U	U	
hlorobenzene	108-90-7	0.75	0.75	UG/L	U	U	
hloroethane	75-00-3	0.32	0.32	UG/L	U	U	
Chloroform	67-66-3	0.34	0.34	UG/L	U	U	
Chloromethane	74-87-3	0.35	0.35	UG/L	U	U	
is-1,2-Dichloroethene	156-59-2	0.81	0.81	UG/L	U	U	
is-1,3-Dichloropropene	10061-01-5	0.36	0.36	UG/L	U	U	
Cyclohexane	110-82-7	0.18	0.18	UG/L	U	U	
Dibromochloromethane	124-48-1	0.32	0.32	UG/L	U	U	
Dichlorodifluoromethane	75-71-8	0.68	0.68	UG/L	U	U	
Ethylbenzene	100-41-4	0.74	0.74	UG/L	U	U	
sopropylbenzene	98-82-8	0.79	0.79	UG/L	U	U	
Methyl acetate	79-20-9	0.50	0.50	UG/L	U	U	
Methyl tert-butyl ether	1634-04-4	0.16	0.16	UG/L	U	U	

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Methylcyclohexane	108-87-2	0.16	0.16	UG/L	U	U	
Methylene Chloride	75-09-2	0.44	0.44	UG/L	U	U	
Styrene	100-42-5	0.73	0.73	UG/L	U	U	
Tetrachloroethene	127-18-4	0.36	0.36	UG/L	U	U	
Toluene	108-88-3	0.51	0.51	UG/L	U	U	
trans-1,2-Dichloroethene	156-60-5	0.90	0.90	UG/L	U	U	
trans-1,3-Dichloropropene	10061-02-6	0.37	0.37	UG/L	U	U	
Trichloroethene	79-01-6	0.46	0.46	UG/L	U	U	
Trichlorofluoromethane	75-69-4	0.88	0.88	UG/L	U	U	
Vinyl chloride	75-01-4	0.90	0.90	UG/L	U	U	
Xylenes, Total	1330-20-7	0.66	0.66	UG/L	U	U	

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CAS No	Sample Name:	TRIP BLANK			Matrix:	Water Res	ult Type: FS		
National	Lab Sample Name:	480-46182-8	Sample Date: 09/19/2013			Validation Level: III			
1.2.2-Tetrachloroethane 79-34-5 0.21 0.21 UGL U U U U U U U U U U U U U U U U U U	Analyte:	CAS No		RL				Validation Notes	
1,2-Trichloro-1,2,2- 76-13-1 0.31 0.31 UGA. U U U U U U U U U	1,1,1-Trichloroethane	71-55-6	0.82	0.82	UG/L	U	U		
	1,1,2,2-Tetrachloroethane	79-34-5	0.21	0.21	UG/L	U	U		
1-Dichloroethane	1,1,2-Trichloro-1,2,2- rifluoroethane	76-13-1	0.31	0.31	UG/L	U	U		
1-Dichloroethene	1,1,2-Trichloroethane	79-00-5	0.23	0.23	UG/L	U	U		
2,4-Trichlorobenzene 120-82-1 0.41 0.41 UG/L U U 2-Dibromo-3-Chloropropane 96-12-8 0.39 0.39 UG/L U U 2-Dichlorobenzene 106-93-4 0.73 0.73 UG/L U U 2-Dichloropenzene 95-50-1 0.79 0.79 UG/L U U 2-Dichloropropane 78-75 0.72 0.72 UG/L U U 2-Dichloropropane 78-87-5 0.72 0.72 UG/L U U 3-Dichlorobenzene 541-73-1 0.78 0.78 UG/L U U 4-Dichlorobenzene 106-46-7 0.84 0.84 UG/L U U 4-Dichlorobenzene 106-46-7 0.84 0.84 UG/L U U 4-Dichlorobenzene 106-46-7 0.84 0.84 UG/L U U Butanone 78-93-3 1.3 1.3 UG/L U U Butano	,1-Dichloroethane	75-34-3	0.38	0.38	UG/L	U	U		
2-Dibromo-3-Chloropropane 96-12-8 0.39 0.39 UG/L U U 2-Dichlorobenzene 106-93-4 0.73 0.73 UG/L U U 2-Dichlorobenzene 95-50-1 0.79 0.79 UG/L U U 2-Dichloropenzene 107-06-2 0.21 0.21 UG/L U U 2-Dichloropenzene 78-87-5 0.72 0.72 UG/L U U 3-Dichloropenzene 541-73-1 0.78 0.78 UG/L U U 4-Dichlorobenzene 106-46-7 0.84 0.84 UG/L U U 4-Dichlorobenzene 108-10-1 2.1 2.1 UG/L U U 4-Dichlorobenzene 108-10-1 2.1 2.1 UG/L U U 4-Dichlorobenzene 108-10-1 2.1 2.1 UG/L U U 4-Dichlorobenzene 108-10-1 0.1 0.1 UU 4-Dichlorobenzene 108-90-7 0.75 0.75 UG/L UU 4-Dichlorobenzene 100-10-10-10-10-10-10-10-10-10-10-10-10-	,1-Dichloroethene	75-35-4	0.29	0.29	UG/L	U	U		
2-Dibromoethane 106-93-4 0.73 0.73 UG/L U U 2-Dichlorobenzene 95-50-1 0.79 0.79 UG/L U U 2-Dichlorobenzene 95-50-1 0.79 0.79 UG/L U U 2-Dichloropenane 107-06-2 0.21 0.21 UG/L U U 3-Dichloropenane 78-87-5 0.72 0.72 UG/L U U 3-Dichlorobenzene 541-73-1 0.78 0.78 UG/L U U 4-Dichlorobenzene 106-46-7 0.84 0.84 UG/L U U 4-Dichlorobenzene 106-46-7 0.84 0.84 UG/L U U 4-Dichlorobenzene 106-46-7 1.3 1.3 UG/L U U 4-Dichlorobenzene 108-10-1 2.1 1.2 UG/L U U 4-Dichlorobenzene 591-78-6 1.2 1.2 UG/L U U 4-Dichlorobenzene 108-10-1 2.1 1.2 UG/L U U 4-Dichlorobenzene 67-64-1 3.0 3.0 UG/L U U 4-Dichlorobenzene 108-10-1 2.1 1.2 UG/L U U 4-Dichlorobenzene 67-64-1 3.0 3.0 UG/L U U 4-Dichlorobenzene 71-43-2 0.41 0.41 UG/L U U 4-Dichlorobenzene 75-25-2 0.26 0.26 UG/L U U 4-Dichlorobenzene 75-25-2 0.26 0.26 UG/L U U 4-Dichlorobenzene 75-25-2 0.26 0.26 UG/L U U 4-Dichlorobenzene 75-15-0 0.19 0.19 UG/L U U 4-Dichlorobenzene 108-90-7 0.75 0.75 UG/L U U 4-Dichlorobenzene 108-90-7 0.75 0.35 UG/L U U 4-Dichlorobenzene 108-90-7 0.36 0.36 UG/L U U 4-Dichlorobenzene 108-90-7 0.18 0.18 UG/L U U 4-Dichlorobenzene 100-41-4 0.34 0.34 UG/L U U 4-Dichlorobenzene 100-41-4 0.74 0.74 UG/L U U 4-Dichlorobenzene 98-82-8 0.79 0.79 UG/L U U 4-Dichlorobenzene 98-82-8 0.79 0.79 UG/L U U 4-Dichlorobenzene 100-41-4 0.74 0.74 UG/L U U	,2,4-Trichlorobenzene	120-82-1	0.41	0.41	UG/L	U	U		
2-Dichlorobenzene 95-50-1 0.79 0.79 UG/L U U 2-Dichloroethane 107-06-2 0.21 0.21 UG/L U U 3-Dichloropropane 78-87-5 0.72 0.72 UG/L U U 3-Dichloropropane 541-73-1 0.78 0.78 UG/L U U 4-Dichlorobenzene 541-73-1 0.78 0.78 UG/L U U 4-Dichlorobenzene 106-46-7 0.84 0.84 UG/L U U 4-Dichlorobenzene 108-10-1 2.1 2.1 UG/L U U 4-Dichlorobenzene 108-10-1 2.1 2.1 UG/L U U 4-Dichlorobenzene 108-10-1 2.1 2.1 UG/L U U 4-Dichlorobenzene 67-64-1 3.0 3.0 UG/L U U 4-Dichlorobenzene 67-64-1 3.0 3.0 UG/L U U 4-Dichlorobenzene 75-27-4 0.39 0.39 UG/L U U 4-Dichlorobenzene 75-25-2 0.26 0.26 UG/L U U 4-Dichlorobenzene 75-25-2 0.26 0.26 UG/L U U 4-Dichlorobenzene 108-90-7 0.75 0.75 UG/L U U 4-Dichlorobenzene 108-90-7 0.75 0.35 UG/L U U 4-Dichlorobenzene 108-90-7 0.78 0.36 UG/L U U 4-Dichlorobenzene 108-90-7 0.38 0.38 UG/L U U 4-Dichlorobenzene 108-90-7 0.48 0.48 UG/L U U 4-Dichlorobenzene 108-90-7 0.48 0.48 UG/L U U 4-Dichlorobenzene 108-90-7 0.48 0.48 UG/L U U 4-Dichlorobenzene 109-90-90-90-90-90-90-90-90-90-90-90-90-9	,2-Dibromo-3-Chloropropan	e 96-12-8	0.39	0.39	UG/L	U	U		
2-Dichloroethane 107-06-2 0.21 0.21 UG/L U U U U U U U U U	,2-Dibromoethane	106-93-4	0.73	0.73	UG/L	U	U		
2-Dichloropropane 78-87-5 0.72 0.72 UG/L U U U U U U U U U	,2-Dichlorobenzene	95-50-1	0.79	0.79	UG/L	U	U		
S-Dichlorobenzene S41-73-1 0.78 0.78 UG/L U U U U U U U U U	,2-Dichloroethane	107-06-2	0.21	0.21	UG/L	U	U		
Helichlorobenzene 106-46-7 0.84 0.84 UG/L U U U U U U U U U	,2-Dichloropropane	78-87-5	0.72	0.72	UG/L	U	U		
Butanone 78-93-3 1.3 1.3 UG/L U U Hexanone 591-78-6 1.2 1.2 UG/L U U Methyl-2-pentanone 108-10-1 2.1 2.1 UG/L U U methyl-2-pentanone 108-10-1 2.1 2.1 UG/L U U methyl-2-pentanone 67-64-1 3.0 3.0 UG/L U U methyl-2-pentanone 71-43-2 0.41 0.41 UG/L U U methyl-2-pentanone 75-27-4 0.39 0.39 UG/L U U methyl-2-pentanone 75-27-4 0.39 0.39 UG/L U U methyl-2-pentanone 75-25-2 0.26 0.26 UG/L U U methyl-2-pentanone 75-25-2 0.26 0.26 UG/L U U methyl-2-pentanone 75-25-2 0.26 0.26 UG/L U U methyl-2-pentanone 75-15-0 0.19 0.19 UG/L U U methyl-2-pentanone 75-15-0 0.19 0.19 UG/L U U methyl-2-pentanone 75-25-2 0.27 0.27 UG/L U U methyl-2-pentanone 108-90-7 0.75 0.75 UG/L U U methyl-2-pentanone 75-00-3 0.32 0.32 UG/L U U methyl-2-pentanone 75-00-3 0.34 0.34 UG/L U U methyl-2-pentanone 108-10-1-5 0.36 0.36 UG/L U U methyl-2-pentanone 108-10-1-5 0.36 0.36 UG/L U U methyl-2-pentanone 108-2-7 0.18 0.18 UG/L U U methyl-2-pentanone 109-41-4 0.74 0.74 UG/L U U methyl-2-pentanone 109-41-4 0.74 0.74 UG/L U U methyl-2-penyl-benzene 98-82-8 0.79 0.79 UG/L U U methyl-benzene 100-41-4 0.74 0.74 UG/L U U methyl-benzene 98-82-8 0.79 0.79 UG/L U U methyl-acetate 79-20-9 0.50 0.50 UG/L U U methyl-acetate 79-20-9 0.50 0.50 UG/L U U methyl-acetate 79-20-9 0.50 UG/L U U methyl-a	3-Dichlorobenzene	541-73-1	0.78	0.78	UG/L	U	U		
Hexanone 591-78-6 1.2 1.2 UG/L U U Methyl-2-pentanone 108-10-1 2.1 2.1 UG/L U U methyl-2-pentanone 67-64-1 3.0 3.0 UG/L U U methyl-2-pentanone 67-64-1 3.0 UG/L U U methyl-2-pentanone 71-43-2 0.41 0.41 UG/L U U methyl-2-pentanone 71-43-2 0.41 0.41 UG/L U U methyl-2-pentanone 71-43-2 0.41 0.41 UG/L U U methyl-2-pentanone 75-27-4 0.39 0.39 UG/L U U methyl-2-pentanone 75-27-4 0.39 0.39 UG/L U U methyl-2-pentanone 75-25-2 0.26 0.26 UG/L U U methyl-2-pentanone 75-25-2 0.26 0.26 UG/L U U methyl-2-pentanone 74-83-9 0.69 0.69 UG/L U U methyl-2-pentanone 74-83-9 0.69 UG/L U U methyl-2-pentanone 74-83-9 0.69 0.69 UG/L U U methyl-2-pentanone 74-83-9 0.69 0.69 UG/L U U methyl-2-pentanone 74-83-9 0.69 0.69 UG/L U U methyl-2-pentanone 75-71-8 0.68 0.68 UG/L U U methyl-2-pentanone 75-71-8 0.69 0.69 UG/L U U methyl-2-pentanone 75-71-8 0.69 0.69 UG/L U U methyl-2-pentanone 75-71-8 0.69 0.69 UG/L U U methyl-2-pentanone 75-71-8 0.69 0.50 0.50 UG/L U U methyl-2-pentanone 75-70-9 0.50 0.50 UG/L U U methyl-2-pentanone 71-4-8-1 0.50 0.50 UG/L U U methyl-2-pentanone 71-4-8-1 0.50 0.50 UG/L U U methyl-2-pentanone 71-4-8-1 0.50 0.50 UG/L	4-Dichlorobenzene	106-46-7	0.84	0.84	UG/L	U	U		
Methyl-2-pentanone 108-10-1 2.1 2.1 UG/L U U exterior 67-64-1 3.0 3.0 UG/L U U emzene 71-43-2 0.41 0.41 UG/L U U comodichloromethane 75-27-4 0.39 0.39 UG/L U U comodichloromethane 75-27-4 0.39 0.39 UG/L U U comodichloromethane 75-27-4 0.39 0.39 UG/L U U comodichloromethane 74-83-9 0.69 0.69 UG/L U U comomethane 74-83-9 0.69 0.69 UG/L U U colorobenzene 108-90-7 0.75 0.27 UG/L U U colorobenzene 108-90-7 0.75 0.75 UG/L U U coloroform 67-66-3 0.34 0.34 UG/L U U coloroform 67-66-3	Butanone	78-93-3	1.3	1.3	UG/L	U	U		
Section Sect	Hexanone	591-78-6	1.2	1.2	UG/L	U	U		
Transpare Tran	Methyl-2-pentanone	108-10-1	2.1	2.1	UG/L	U	U		
Description Temperature	cetone	67-64-1	3.0	3.0	UG/L	U	U		
Tomonform Total	enzene	71-43-2	0.41	0.41	UG/L	U	U		
commomethane 74-83-9 0.69 0.69 UG/L U U rbon disulfide 75-15-0 0.19 0.19 UG/L U U urbon tetrachloride 56-23-5 0.27 0.27 UG/L U U dlorobenzene 108-90-7 0.75 0.75 UG/L U U dloroforme 67-66-3 0.32 0.32 UG/L U U dloroformethane 74-87-3 0.35 0.34 UG/L U U -1,2-Dichloroethene 156-59-2 0.81 0.81 UG/L U U -1,3-Dichloropropene 10061-01-5 0.36 0.36 UG/L U U -1,3-Dichloropropene 10061-01-5 0.36 0.36 UG/L U U bromochloromethane 124-48-1 0.32 0.32 UG/L U U chlorodifluoromethane 75-71-8 0.68 0.68 UG/L U U opropylb	omodichloromethane	75-27-4	0.39	0.39	UG/L	U	U		
rbon disulfide 75-15-0 0.19 0.19 UG/L U U rbon tetrachloride 56-23-5 0.27 0.27 UG/L U U rbon tetrachloride 56-23-5 0.27 0.75 UG/L U U rbon diorobenzene 108-90-7 0.75 0.75 UG/L U U rbon dioroform 75-00-3 0.32 0.32 UG/L U U rbon dioroform 67-66-3 0.34 0.34 UG/L U U rbon dioroform 74-87-3 0.35 0.35 UG/L U U rbon dioroform 156-59-2 0.81 0.81 UG/L U U rbon dioroform 10061-01-5 0.36 0.36 UG/L U U rbon dioroform 10061-01-5 0.36 0.36 UG/L U U rbon dioroform 10041-48-1 0.32 0.32 UG/L U U rbon dioroform 10041-4 0.74 0.74 UG/L U U rbon dioroformethane 10041-4 0.7	omoform	75-25-2	0.26	0.26	UG/L	U	U		
Arbon tetrachloride 56-23-5 0.27 0.27 UG/L U U U U U U U U U U U U U U U U U U U	omomethane	74-83-9	0.69	0.69	UG/L	U	U		
108-90-7 0.75 0.75 UG/L U U U U U U U U U	rbon disulfide	75-15-0	0.19	0.19	UG/L	U	U		
Description 100 10	urbon tetrachloride	56-23-5	0.27	0.27	UG/L	U	U		
Description Color	nlorobenzene	108-90-7	0.75	0.75	UG/L	U	U		
100 100	hloroethane	75-00-3	0.32	0.32	UG/L	U	U		
S-1,2-Dichloroethene 156-59-2 0.81 0.81 UG/L U U U S-1,3-Dichloropropene 10061-01-5 0.36 0.36 UG/L U U U U U U U U U U U U U U U U U U U	hloroform	67-66-3	0.34	0.34	UG/L	U	U		
S-1,3-Dichloropropene 10061-01-5 0.36 0.36 UG/L U U U U U U U U U	nloromethane	74-87-3	0.35	0.35	UG/L	U	U		
Veclohexane	s-1,2-Dichloroethene	156-59-2	0.81	0.81	UG/L	U	U		
bromochloromethane 124-48-1 0.32 0.32 UG/L U U ichlorodifluoromethane 75-71-8 0.68 0.68 UG/L U U hylbenzene 100-41-4 0.74 0.74 UG/L U U opropylbenzene 98-82-8 0.79 0.79 UG/L U U ethyl acetate 79-20-9 0.50 0.50 UG/L U U	s-1,3-Dichloropropene	10061-01-5	0.36	0.36	UG/L	U	U		
ichlorodifluoromethane 75-71-8 0.68 0.68 UG/L U U hylbenzene 100-41-4 0.74 0.74 UG/L U U opropylbenzene 98-82-8 0.79 0.79 UG/L U U ethyl acetate 79-20-9 0.50 0.50 UG/L U U	yclohexane	110-82-7	0.18	0.18	UG/L	U	U		
hylbenzene 100-41-4 0.74 0.74 UG/L U U propylbenzene 98-82-8 0.79 0.79 UG/L U U ethyl acetate 79-20-9 0.50 0.50 UG/L U U	bromochloromethane	124-48-1	0.32	0.32	UG/L	U	U		
propylbenzene 98-82-8 0.79 0.79 UG/L U U ethyl acetate 79-20-9 0.50 0.50 UG/L U U	ichlorodifluoromethane	75-71-8	0.68	0.68	UG/L	U	U		
ethyl acetate 79-20-9 0.50 0.50 UG/L U U	thylbenzene	100-41-4	0.74	0.74	UG/L	U	U		
	sopropylbenzene	98-82-8	0.79	0.79	UG/L	U	U		
ethyl tert-butyl ether 1634-04-4 0.16 0.16 UG/L U U	Methyl acetate	79-20-9	0.50	0.50	UG/L	U	U		
	Methyl tert-butyl ether	1634-04-4	0.16	0.16	UG/L	U	U		

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Analysis	Method	8260B

Methylcyclohexane	108-87-2	0.16	0.16	UG/L	U	U	
Methylene Chloride	75-09-2	0.44	0.44	UG/L	U	U	
Styrene	100-42-5	0.73	0.73	UG/L	U	U	
Tetrachloroethene	127-18-4	0.36	0.36	UG/L	U	U	
Toluene	108-88-3	0.51	0.51	UG/L	U	U	
trans-1,2-Dichloroethene	156-60-5	0.90	0.90	UG/L	U	U	
trans-1,3-Dichloropropene	10061-02-6	0.37	0.37	UG/L	U	U	
Trichloroethene	79-01-6	0.46	0.46	UG/L	U	U	
Trichlorofluoromethane	75-69-4	0.88	0.88	UG/L	U	U	
Vinyl chloride	75-01-4	0.90	0.90	UG/L	U	U	
Xylenes, Total	1330-20-7	0.66	0.66	UG/L	U	U	

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Analysis Method 9060

Sample Name:	DUP			Matrix:	Water Resu	lt Type: FS
Lab Sample Name:	480-46182-7	Sample 1	Date: 09/1	9/2013	V	alidation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier
Total Organic Carbon	7440-44-0	0.43	0.43	MG/L	U	U
Sample Name:	LVRA07-MNA	GW-MW-2		Matrix:	Water Resu	lt Type: FS
Lab Sample Name:	480-46182-3	Sample 1	Date: 09/1	9/2013	V	alidation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier
Total Organic Carbon	7440-44-0	0.82	0.43	MG/L	J	J
Sample Name:	LVRA07-MNA	GW-MW-3		Matrix:	Water Resu	lt Type: FS
Lab Sample Name:	480-46182-1	Sample 1	Date: 09/1	9/2013	V	alidation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Notes Qualifier
Total Organic Carbon	7440-44-0	0.44	0.43	MG/L	J	1
Sample Name:	LVRA07-MNA	GW-MW-5		Matrix:	Water Resu	lt Type: FS
Lab Sample Name:	480-46182-6	Sample 1	Date: 09/1	9/2013	v	alidation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Notes Qualifier
Total Organic Carbon	7440-44-0	0.69	0.43	MG/L	J	1
Sample Name:	LVRA07-MNA	GW-MW-6		Matrix:	Water Resu	lt Type: FS
Lab Sample Name:	480-46182-5	Sample 1	Date: 09/1	9/2013	v	alidation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier
Total Organic Carbon	7440-44-0	1.3	0.43	MG/L		
Sample Name:	LVRA07-MNA	GW-MW-D	02	Matrix:	Water Resu	lt Type: FS
Lab Sample Name:	480-46182-2	Sample 1	Date: 09/1	9/2013	v	alidation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Notes Qualifier
Total Organic Carbon	7440-44-0	0.43	0.43	MG/L	U	U

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Analysis Method RSK-175

Sample Name:	DUP			Matrix:	Water Res	ult Type: FS	
Lab Sample Name:	480-46182-7	Sample	Date: 09/1	9/2013	•	Validation Lev	vel: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Qualifier	Validation Notes
Ethane	74-84-0	0.49	0.49	UG/L	U	U	
Ethene	74-85-1	0.52	0.52	UG/L	U	U	
Methane	74-82-8	0.22	0.22	UG/L	U	U	
Sample Name:	LVRA07-MNA	AGW-MW-2	2	Matrix:	Water Res	ult Type: FS	
Lab Sample Name:	480-46182-3	Sample 3	Date: 09/1	9/2013	•	Validation Lev	vel: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Qualifier	Validation Notes
Ethane	74-84-0	0.49	0.49	UG/L	J	J	
Ethene	74-85-1	0.52	0.52	UG/L	U	U	
Methane	74-82-8	75	0.22	UG/L			
Sample Name:	LVRA07-MNA	AGW-MW-3	}	Matrix:	Water Res	ult Type: FS	
Lab Sample Name:	480-46182-1	Sample	Date: 09/1	9/2013	•	Validation Lev	vel: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Qualifier	Validation Notes
Ethane	74-84-0	0.49	0.49	UG/L	U	U	
Ethene	74-85-1	0.52	0.52	UG/L	U	U	
Methane	74-82-8	0.22	0.22	UG/L	U	U	
Sample Name:	LVRA07-MNA	AGW-MW-5	i	Matrix:	Water Res	ult Type: FS	
Lab Sample Name:	480-46182-6	Sample	Date: 09/1	9/2013	•	Validation Lev	vel: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Qualifier	Validation Notes
Ethane	74-84-0	0.49	0.49	UG/L	U	U	
Ethene	74-85-1	0.52	0.52	UG/L	U	U	
Methane	74-82-8	0.22	0.22	UG/L	U	U	
Sample Name:	LVRA07-MNA	AGW-MW-6	,	Matrix:	Water Res	ult Type: FS	
Lab Sample Name:	480-46182-5	Sample	Date: 09/1	9/2013	•	Validation Lev	vel: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Qualifier	Validation Notes
Ethane	74-84-0	0.49	0.49	UG/L	U	U	
	74.05.1	0.52	0.52	TIC/I	T.T.	U	
Ethene	74-85-1	0.52	0.52	UG/L	U	U	

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Analysis Method RSK-175

Sample Name:	LVRA07-MNA0	VRA07-MNAGW-MW-D2 Matrix: Water Result Type: FS					
Lab Sample Name:	480-46182-2	Sample 1	Date: 09/19/	2013	•	Validation Le	evel: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Qualifier	Validation Notes
Ethane	74-84-0	0.49	0.49	UG/L	U	U	
Ethene	74-85-1	0.52	0.52	UG/L	U	U	
Methane	74-82-8	0.22	0.22	UG/L	U	U	
Sample Name:	RINSE BLANK			Matrix:	Water Resi	ult Type: FS	3
Lab Sample Name:	480-46182-4	Sample 1	Date: 09/19/	2013	7	Validation Le	vel: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Qualifier	Validation Notes
Ethane	74-84-0	0.49	0.49	UG/L	U	U	
Ethene	74-85-1	0.52	0.52	UG/L	U	U	
Methane	74-82-8	0.22	0.22	UG/L	U	U	

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Analysis Method SM 2320B

Comple Name	DIID			Motui	Water Dogult Trues ES
Sample Name:	DUP			Matrix:	· -
Lab Sample Name:	480-46182-7	Sample I	Date: 09/1	9/2013	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Validation Validation Notes Qualifier Qualifier
Alkalinity, Total	STL00171	142	0.79	MG/L	
Sample Name:	LVRA07-MNA	GW-MW-2		Matrix:	Water Result Type: FS
Lab Sample Name:	480-46182-3	Sample I	Date: 09/1	9/2013	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Validation Validation Notes Qualifier Qualifier
Alkalinity, Total	STL00171	201	0.79	MG/L	
Sample Name:	LVRA07-MNA	GW-MW-3		Matrix:	Water Result Type: FS
Lab Sample Name:	480-46182-1	Sample I	Date: 09/1	9/2013	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Validation Validation Notes Qualifier Qualifier
Alkalinity, Total	STL00171	165	0.79	MG/L	
Sample Name:	LVRA07-MNA	GW-MW-5		Matrix:	Water Result Type: FS
Lab Sample Name:	480-46182-6	Sample I	Date: 09/1	9/2013	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Validation Validation Notes Qualifier Qualifier
Alkalinity, Total	STL00171	68.5	0.79	MG/L	
Sample Name:	LVRA07-MNA	GW-MW-6		Matrix:	Water Result Type: FS
Lab Sample Name:	480-46182-5	Sample I	Date: 09/1	9/2013	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Validation Validation Notes Qualifier Qualifier
Alkalinity, Total	STL00171	92.2	0.79	MG/L	
Sample Name:	LVRA07-MNA	GW-MW-D)2	Matrix:	Water Result Type: FS
Lab Sample Name:	480-46182-2	Sample I	Date: 09/1	9/2013	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Validation Validation Notes Qualifier Qualifier
Alkalinity, Total	STL00171	141	0.79	MG/L	

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Analysis Method SM 3500 FE D

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Sample Name:	DUP			Matrix:	water Res	ult Type: FS
Lab Sample Name:	480-46182-7	Sample 1	Date: 09/1	9/2013	•	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier
Ferrous Iron, Dissolved	15438-31-0	0.075	0.075	MG/L	U HF	U
Sample Name:	LVRA07-MNA	GW-MW-2	,	Matrix:	Water Res	ult Type: FS
Lab Sample Name:	480-46182-3	Sample	Date: 09/1	9/2013	•	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier
Ferrous Iron, Dissolved	15438-31-0	0.075	0.075	MG/L	U HF	U
Sample Name:	LVRA07-MNA	GW-MW-3		Matrix:	Water Res	ult Type: FS
Lab Sample Name:	480-46182-1	Sample	Date: 09/1	9/2013	•	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier
Ferrous Iron, Dissolved	15438-31-0	0.075	0.075	MG/L	U HF	U
Sample Name:	LVRA07-MNA	GW-MW-5		Matrix:	Water Res	ult Type: FS
Lab Sample Name:	480-46182-6	Sample	Date: 09/1	9/2013	•	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier
Ferrous Iron, Dissolved	15438-31-0	0.075	0.075	MG/L	U HF	U
Sample Name:	LVRA07-MNA	GW-MW-6		Matrix:	Water Res	ult Type: FS
Lab Sample Name:	480-46182-5	Sample	Date: 09/1	9/2013	•	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier
Ferrous Iron, Dissolved	15438-31-0	0.080	0.075	MG/L	J HF	J
Sample Name:	LVRA07-MNA	GW-MW-I	D2	Matrix:	Water Res	ult Type: FS
Lab Sample Name:	480-46182-2	Sample	Date: 09/1	9/2013	•	Validation Level: III
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier
Ferrous Iron, Dissolved	15438-31-0	0.075	0.075	MG/L	U HF	U

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Analysis Method SM 4500 S2 F

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Sample Name:	DUP Matrix: Water Result Type: FS							
Lab Sample Name:	480-46182-7	Sample Date: 09/19/201			13 Validation Level: III			
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier		
Sulfide	18496-25-8	0.67	0.67	MG/L	U	U		
Sample Name:	LVRA07-MNAGW-MW-2 Mat			Matrix:	Water Res	ult Type: FS		
Lab Sample Name:	480-46182-3 Sample Date: 09/19/2			9/2013	Validation Level: III			
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier		
Sulfide	18496-25-8	0.67	0.67	MG/L	U	U		
Sample Name:	LVRA07-MNAGW-MW-3			Matrix:	Water Res	ult Type: FS		
Lab Sample Name:	480-46182-1 Sample Date: 09/19/2			9/2013	•	Validation Level: III		
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier		
Sulfide	18496-25-8	0.67	0.67	MG/L	U	U		
Sample Name:	LVRA07-MNAGW-MW-5			Matrix:	Water Res	ult Type: FS		
Lab Sample Name:	480-46182-6	Sample Date: 09/19/2		9/2013	•	Validation Level: III		
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier		
Sulfide	18496-25-8	0.67	0.67	MG/L	U	υ		
Sample Name:	LVRA07-MNAGW-MW-6 Ma			Matrix:	Water Res	ult Type: FS		
Lab Sample Name:	480-46182-5	-5 Sample Date: 09/19/20		9/2013	•	Validation Level: III		
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier		
Sulfide	18496-25-8	0.67	0.67	MG/L	U	U		
Sample Name:	LVRA07-MNAGW-MW-D2		Matrix:	Water Res	ult Type: FS			
Lab Sample Name:	480-46182-2 Sample Date: 09/19/2		9/2013	•	Validation Level: III			
Analyte:	CAS No	Result Value	RL	Result Units	Lab Qualifier	Validation Validation Notes Qualifier		

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