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March 23, 2017
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Mr. Bill Sadler, P.G.
ADEQ – Solid Waste Management Division
5301 Northshore Drive
North Little Rock, Arkansas 72118

Subject: **NABORS Landfill**
Fourth Quarter 2016 Groundwater Monitoring and Gas Monitoring Reports

Mr. Sadler,

In accordance with Contract Number 4600033394, SCS Engineers is submitting the Fourth Quarter Groundwater Monitoring Report (GWMR) and Fourth Quarter Gas Monitoring Report summarizing monitoring activities performed in December 2016 at the NABORS Landfill.

If you have questions or comments regarding these reports, please do not hesitate to contact us at (913) 681-0030.

Sincerely,

Dillon Baird, P.E.
Project Engineer
SCS ENGINEERS

Floyd Cotter, P.E.
Project Manager
SCS ENGINEERS

Enclosure: Fourth Quarter 2016 Groundwater Monitoring Report
Fourth Quarter 2016 Gas Monitoring Report
Fourth Quarter 2016 Sanitas Database Flat File

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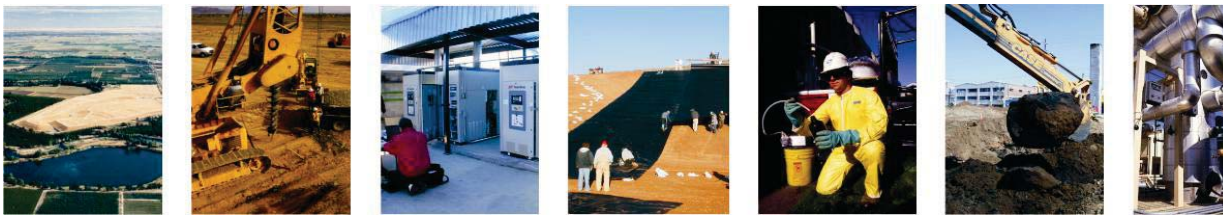

FOURTH QUARTER 2016 NABORS LANDFILL GROUNDWATER MONITORING REPORT

Presented to:
Arkansas Department of Environmental Quality

ADEQ
ARKANSAS
Department of Environmental Quality

5301 Northshore Drive
North Little Rock, AR 72118-5317
501-682-0744

Presented by:
SCS ENGINEERS
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North Little Rock, AR 72113
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March 23, 2017
File No. 27214218.01

Offices Nationwide
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1 INTRODUCTION

This report summarizes the results of the Fourth Quarter 2016 Groundwater Assessment Monitoring event at the NABORS Landfill. Field work was conducted by personnel from SCS Engineers (SCS) on December 19-21, 2016. This report was prepared under the provisions of Contract No. 4600033394 between the Arkansas Department of Environmental Quality (ADEQ) and SCS.

The NABORS Landfill originally operated under Solid Waste Disposal Permit number 0249-S, as issued to RLH, Inc. (RLH) by the Arkansas Department of Environmental Quality (ADEQ) on June 14, 1988. The solid waste permit was transferred to NABORS on August 31, 2005. The NABORS Landfill is currently under Solid Waste Disposal Permit 0249-S1-R2 issued by the ADEQ on August 10, 2006. Although the permit is considered "open", the landfill is not receiving waste. The ADEQ is currently managing the closure of the site. The analytical work for this sampling event was conducted by Environmental Science Corporation (ESC) of Nashville, Tennessee and groundwater sampling was conducted by SCS.

1.1 SITE DESCRIPTION

The Northwest Arkansas Regional Solid Waste Management District (District) owns both a Class 1 and Class 4 Landfill at the subject site. The NABORS Landfills (herein referred to as Landfill) are located on approximately 700 acres near Three Brothers (Baxter County), Arkansas. The permitted Class 1 area is located in a portion of the SW1/4 of SE1/4 of Section 26, Township 21 North, Range 14 West, and NW1/4 of NE1/4 of Section 35, Township 21 North, Range 14 West. A general geographic location map is included as **FIGURE 1** in **APPENDIX A**. All report Figures are provided in **APPENDIX A**.

1.2 SITE GROUNDWATER MONITORING SYSTEM

The Landfill groundwater monitoring system has evolved over time into a system of monitoring points including both monitoring wells and local springs. The current Assessment Monitoring Program was triggered by the statistically significant occurrence of certain parameters documented in the original Detection Monitoring System. In turn, the required contingencies for Nature and Extent characterization and other regulatory provisions were addressed through the development of the site.

More specifically, the Second Detection Monitoring System for the site that was intended to comply with the provisions of Reg.22.1202 through Reg.22.1204 was approved by ADEQ in 1998. Following the Second Half 2005 sampling event, the ADEQ was notified under the provisions of 22.1204(c) of a statistically significant increase for volatile organic compounds (VOC) at MW-1. These detections were verified during the Second Half 2006 sampling event and the facility began Assessment Monitoring at wells MW-1, MW-2, MW-3, MW-4, MW-5, MW-6, and MW-7 in June of 2006 as required by Regulation 22.1205. At that time these seven wells were the only permitted facility Detection Monitoring wells.

The Assessment Monitoring program was later expanded to include all the wells/springs listed below when the new permit was issued in August of 2006 (0249-S1-R2), and when the Nature and Extent Characterization was completed. Monitoring was conducted under this program through the Third Quarter 2012.

Assessment Monitoring continued on a quarterly basis as per 22.1205(d) based upon:

- The presence of VOCs at MW-1
- Detected concentrations of arsenic at various wells, some of which were above the Groundwater Protection Standard (GWPS)
- Vinyl chloride detections at CAO-1 which historically exceed the GWPS.

The current groundwater monitoring system for the NABORS Landfill consists of twenty-four (24) monitoring wells, thirteen (13) springs, and a Leachate sample for a total of thirty-eight (38) Assessment Monitoring points. These sampling points are listed below and are located on **FIGURE 2.**

NABORS GROUNDWATER ASSESSMENT MONITORING POINTS				
MW-1	**NAB-1**	MW-577	SP-4	NE-6
MW-2	NAB-2	MW-689D	SP-5	MW-1R
MW-3	NAB-3	MW-633D	SP-7	
MW-4	NAB-4	TSP-1	Spring A	
MW-5	NAB-7	TSP-2	Spring B	
MW-6	CAO-3	TSP-3	Leachate	
MW-7	NAB-8	TSP-4	NE-2	
CAO-1	MW-509D	Entrance Seep	NE-3	
CAO-2	Class I Draw	Class IV Draw	Spring near NE-3	
Note: NAB-1 is damaged at approximately 68 feet below top of casing and is no longer being sampled.				

2 GROUNDWATER SAMPLING

The Fourth Quarter 2016 Assessment Monitoring sampling event was conducted on December 19-21, 2016. A representative of SCS collected samples from twenty-three (23) monitoring wells (MW-1 through MW-7, MW-1R, MW-509D, MW-577, MW-633D, MW-689D, CAO-1, CAO-2, CAO-3, NAB-2, NAB-3, NAB-4, NAB-7, NAB-8, NE-2, NE-3 and NE-6) and five (5) springs (Spring A, Class I Draw Spring, Class IV Draw Spring, SP-7, and the Landfill Entrance Seep). All additional springs were dry, or had been covered by landfill construction and no longer exist. The current *Groundwater Sampling and Analysis Plan* dated December 2011 (ADEQ document #61474) was utilized for this sampling event. The procedures for obtaining groundwater samples, parameters analyzed, sample preservation and handling are discussed in the following sections.

2.1 WATER LEVEL DETERMINATION

Prior to evacuating each well for sampling, the depth to water was measured using an electronic water level probe. The measurements were taken to the nearest 0.01-foot from the top of the well casing and the information was used to calculate the volume of water in the well. Because non-dedicated equipment was used to obtain water levels, procedures were instituted to insure the samples were not contaminated. The electronic water level probe is constructed of inert materials and was de-contaminated with distilled water prior to use at each well.

2.2 WELL EVACUATION

The water in a well prior to sampling may not be representative of in-situ groundwater quality. Therefore, the groundwater technician used an electric submersible pump with dedicated sampling tubing to purge a minimum of three casing volumes at a rate that did not excessively agitate the recharge water. Since non-dedicated equipment was used to purge the wells, procedures were utilized to insure the samples were not contaminated. Clean, non-powdered, nitrile gloves were worn by the sampling personnel. Measures were taken to prevent surface soils from coming in contact with the purging equipment and lines, which could introduce contaminants to the well.

2.3 EQUIPMENT DECONTAMINATION PROCEDURES

All equipment that was used in the monitoring wells and had contact with the samples was thoroughly cleaned before use. These devices included a water level probe and a submersible pump. The water level probe was washed with potable water and phosphate-free laboratory detergent. Next, the probe was rinsed with potable water and finally, rinsed with distilled water. The water level probe was then placed in a plastic bag to reduce contact with air and transported into the field. After a water level was measured at each well, a paper towel was soaked with distilled water and as the probe was reeled up the tape and probe were wiped clean.

The submersible pump was initially flushed with potable water and phosphate-free detergent. Next, the pump was rinsed with potable water, and finally rinsed and flushed with distilled water in a portable decontamination tub prior to use in each well. The pump was transported in a clean, sealed tub to minimize contact with the air prior to use at each well.

2.4 SAMPLE EXTRACTION

The technique used to withdraw each groundwater sample from the wells was selected based on consideration of the parameters analyzed in the sample. To insure the groundwater sample is representative of the formation, it is important to minimize physically altering or chemically contaminating the sample during the withdrawal process. In order to minimize the possibility of sample contamination the groundwater technicians did the following:

- Made sure clean sampling equipment was not placed directly on the ground or other contaminated surfaces prior to insertion into the well.

- Gently lowered and retrieved sampling equipment in order to prevent undue disturbance of the water column. Monitoring wells were purged and samples were collected using a submersible pump. Wells that went dry during purging were allowed to recharge for approximately 24 hours and samples were collected using a disposable bailer.

- Transferred samples to the appropriate containers in a manner that minimized agitation and aeration.

Samples were collected and containerized in the order of volatilization sensitivity of the parameters. Spring samples are collected where able, by utilizing a grab sample method by holding the sample bottle within the spring water flow path. The list of parameters analyzed is presented in **TABLE 1**.

TABLE 1. ASSESSMENT MONITORING CONSTITUENTS (AMC)

APPENDIX 1 VOLATILES	INDICATOR PARAMETERS	INORGANICS
ACETONE	CHLORIDE	ANTIMONY
ACRYLONITRILE	pH	ARSENIC
BENZENE	SULFATE	BARIUM
BROMOCHLOROMETHANE	TDS	BERYLLIUM
BROMODICHLOROMETHANE	TOC	CADMIUM
CHROMIUM	SPECIFIC CONDUCTANCE	CHROMIUM
CARBON TETRACHLORIDE		COBALT
CARBON DISULFIDE		COPPER
CHLOROBENZENE		LEAD
CHLOROETHANE		NICKEL
CHLOROFORM		SELENIUM
DIBROMOCHLOROMETHANE		SILVER
1,2, DICHLOROBENZENE		THALLIUM
1,4, DICHLOROBENZENE		VANADIUM
TRANS- 1,4-DICHLORO-2-BUTENE		ZINC
1,1 DICHLOROETHANE		IRON
1,2 DICHLOROETHANE		MANGANESE
CIS-1,2,-DICHLOROETHYLENE		TIN
TRANS-1,2-DICHLOROETHYLENE		
1,2-DICHLOROPROPANE		
CIS-1,3-DICHLOROPROPENE		
TRANS-1,3-DICHLOROPROPENE		
ETHYLBENZENE		
2-HEXANONE		
METHYL BROMIDE		
METHYL CHLORIDE		
METHYLENE BROMIDE		
METHYLENE CHLORIDE		
METHYL ETHYL KETONE		
METHYL IODIDE		
4- METHYL-2-PENTANONE		
STYRENE		
1,1,1,2-TETRACHLOROETHANE		
1,1,2,2,-TETRACHLOROETHANE		
TETRACHLOROETHANE		
TOLUENE		
1,1,1-TRICHLOROETHANE		
1,1,2-TRICHLOROETHANE		
TRICHLOROETHYLENE		
TRICHLOROFLUOROMETHANE		
1,2,3-TRICHLOROPROPANE		
VINYL ACETATE		
VINYL CHLORIDE		
XYLENE		

The collection order for the common groundwater parameters was as follows:

- Volatile Organic Compounds (VOCs)
- Total Organic Carbon (TOC)
- Sulfate, Chloride, and Total Dissolved Solids (TDS)
- Total Metals

2.5 FIELD TESTING

Some of the parameters evaluated are physically or chemically unstable and were measured immediately after collection by the SCS representative using a flow cell. Examples of unstable elements or properties include pH and temperature. Although the turbidity and specific conductance (inverse of electrical resistance) of a substance are relatively stable, these parameters were also measured in the field. This information was recorded on *Groundwater Monitoring Sampling Records* presented in **APPENDIX B**. A summary of the field measurements for the Fourth Quarter 2016 Assessment Monitoring sampling event is presented in **TABLE 2**.

TABLE 2. FOURTH QUARTER 2016 FIELD MEASUREMENTS

Well #	Date	Time	TOC Elev. (fmsl)	GW Depth (ft.)	GW Elev. (fmsl)	pH (SU)	Temp. (°C)	Spec. Cond. (uS/cm)	Turbidity (NTU)
MW-1R	12/20/2016	1629	1067.57	69.77	997.80	6.14	16.8	1523	18.5
MW-1	12/20/2016	1558	1067.26	67.65	999.61	6.15	16.8	1618	35.7
MW-2	12/20/2016	1652	1001.21	35.00	966.21	6.60	16.5	622	2.22
MW-3	12/21/2016	0804	994.48	26.84	967.64	6.32	17.3	868	3.07
MW-4	12/21/2016	0847	1012.11	87.02	925.09	6.34	15.7	655	10.7
MW-5	12/21/2016	0911	1004.38	77.77	926.61	6.27	15.2	702	8.47
MW-6	12/21/2016	0946	1000.38	56.75	943.63	6.50	14.3	733	5.68
MW-7	12/20/2016	1438	999.66	8.64	991.02	6.25	15.4	578	8.40
CAO-1	12/21/2016	1022	1026.40	26.45	999.95	5.91	16.7	1466	5.65
CAO-2	12/21/2016	0748	998.80	28.23	970.57	6.35	16.5	489	12.7
CAO-3	12/21/2016	0720	984.20	10.62	973.58	6.60	15.6	469	4.13
NAB-2	12/20/2016	1253	993.98	79.26	914.72	6.24	16.0	744	19.8
NAB-3	12/20/2016	0819	921.49	25.18	896.31	7.75	15.3	721	13.5
NAB-4	12/20/2016	1042	1004.20	108.07	896.13	6.26	15.8	743	13.4
NAB-7	12/20/2016	1515	1012.36	22.51	989.85	6.30	16.0	636	2.71
*NAB-8	12/20/2016	0700	1039.21	78.54	960.67	4.00	5.2	624	17.8
MW-509D	12/20/2016	1415	1014.20	24.25	989.95	6.20	16.8	667	59.1
MW-577	12/20/2016	1210	982.60	42.25	940.35	6.28	16.1	674	4.79
MW-633D	12/20/2016	1130	1050.10	64.40	985.70	6.41	17.0	719	4.99
MW-689D	12/20/2016	0931	966.20	26.08	940.12	6.38	16.9	692	16.8
NE-2	12/19/2016	1150	976.98	53.10	923.88	8.00	13.5	2192	19.2
*NE-3	12/20/2016	0730	846.91	8.30	838.61	4.64	7.1	592	13.8
NE-6	12/21/2016	0820	901.42	9.35	892.07	6.32	15.5	623	3.36

*Note: NAB-8 and NE-3 went dry during purging (at approx. 3 and 4.5 gallons, respectively)

2.6 FIELD QA/QC PROCEDURES

For QA/QC purposes, a duplicate sample of MW-6 was collected and labeled Dupe. Procedures utilized for collecting the duplicate sample were identical to the sampling protocol detailed in Section 2.4 and collected at the same time as the MW-6 samples. The duplicate sample was collected to verify the consistency and precision of the sampling and testing procedures.

A field blank was also collected and labeled FB. The field blank consisted of distilled water poured into a sample container under field conditions and returned for laboratory analysis. The SCS field representative prepared the field blank for all the required monitoring parameters. The field blank was used to verify that the sample collection and handling process or ambient conditions, such as airborne materials or other factors unique to the sampling area did not affect the quality of the samples. A volatile organic analyte (VOA) trip blank was also included as part of the field QA/QC procedures. The trip blank was prepared in the laboratory utilizing de-ionized water, transported to the site, handled as a sample (yet never opened in the field), and returned to the laboratory for analysis. Trip blank results are used to verify that the sample containers were adequately prepared/handled in the laboratory, and that the groundwater samples were protected from contamination during transport.

An equipment blank, labeled EB, was prepared on site by pouring de-ionized water over the water level probe, gloves, and through a disposable bailer. Equipment blank results are used to verify that proper protocols for collection of samples and decontamination of equipment were followed.

2.7 HANDLING/TRANSPORT/CUSTODY

Samples were accompanied by a Chain-of-Custody record that includes the name of the facility, collector's signature, monitoring point identification number, date, time, type of sample, number of containers, and analyses required. Samples collected from the Landfill site were placed in sample containers provided by the Laboratory. Containers were certified clean by the supplier and transported with ice to preserve samples.

Attached to the sample container at the time of collection is the sample label. The following information is recorded on the sample label:

- Project or facility name
- Sample type
- Sample location number (well number)
- Preservation type
- Sampling date and time
- Sample collector's name or initials

Documentation for the sample collection process and other important information was recorded on the chain of custody. The standard format includes the date, time, type of sample, code for sample analysis, unique sample number, and sampling location. The entries were signed by the sample collector.

2.8 SAMPLE PRESERVATION

In accordance with the facility's *Sampling and Analysis Plan*, the samples were placed in an ice chest for preservation and cooled to approximately 4 degrees Celsius. Custody was retained by a SCS representative from the time of collection until shipment via Federal Express to Environmental Science Corp. (ESC) in Nashville, Tennessee. Laboratory analytical results and a copy of the ESC Chain-of-Custody form are included in **APPENDIX C**.

3 FOURTH QUARTER 2016 ASSESSMENT MONITORING EVENT

The sampling results summarized in this report are for the Fourth Quarter 2016 Assessment Monitoring Event. The results for this event, conducted on December 19-21, 2016 are provided in the following sections, tables, and appendices. In addition, all historical groundwater data was evaluated statistically to determine if significant differences exist between compliance and background concentrations at each monitoring point.

3.1 GROUNDWATER ELEVATION, FLOW DIRECTION & RATE

There are currently twenty-three monitoring wells located around the Landfill area. Water level elevations were measured for monitoring wells MW-1R, MW-1, MW-2, MW-3, MW-4, MW-5, MW-6, MW-7, CAO-1, CAO-2, CAO-3, NAB-2, NAB-3, NAB-4, NAB-7, NAB-8, NE-2, NE-3, NE-6, MW-509D, MW-577, MW-689 and MW-633D during the Fourth Quarter 2016 event. **TABLE 2** summarizes the results of the water level and field measurements for this event. The highest water level elevation during this event was measured in monitoring well CAO-1, located east of Area 1-2, and the lowest elevation occurred in monitoring well NE-3, located southeast of Area 1-3. A potentiometric surface map was constructed utilizing the water levels measured during the Fourth Quarter 2016 Assessment Monitoring Event and is presented as **FIGURE 2**.

As **FIGURE 2** indicates, groundwater within the uppermost aquifer was found to flow to the west/northwest in Area 1-2 and generally to the east-southeast in Area 1-3. Based on this flow pattern, monitoring wells MW-1, MW-1R, MW-7, and NAB-7 are upgradient wells, and wells MW-2, MW-3, MW-4, MW-5, MW-6, NAB-2, NAB-3, NAB-4, NAB-8, MW-509D, MW-633D, MW-577, MW-689, CAO-1, CAO-2, and CAO-3 monitor the groundwater downgradient of the landfill.

Based on the principles of Darcian flow, the average linear velocity (groundwater flow rate) during the Fourth Quarter 2016 Assessment Monitoring event was calculated utilizing the following equation:

$$V_x = (K*i) / n_e$$

where,

V_x is the average linear velocity (length/time),
K is the hydraulic conductivity (length/time),
i is the hydraulic gradient (length/length),
and n_e is the effective porosity (decimal).

The hydraulic gradient in the Area 1-2 portion of the Landfill was calculated by comparing upgradient well, MW-1R, to the most directly downgradient well, MW-5. The change in head of 71.20 feet between the two wells over a distance of approximately 1,432 feet produces a hydraulic gradient of 0.05 (ft/ft).

The hydraulic gradient in the Area 1-3 portion of the Landfill was calculated by comparing upgradient well, MW-1R, to the most directly downgradient well, NAB-2. The change in head of 83.08 feet between the two wells over a distance of approximately 693 feet produces a hydraulic gradient of 0.12 (ft/ft).

Grubbs, Garner, & Hoskyn, Inc. reported an average hydraulic conductivity of $1.0x10^{-3}$ cm/sec in the site's *Hydrogeologic and Geotechnical Report* (1987). This hydraulic conductivity for the uppermost aquifer was used to aid in the flow rate calculations. SCS utilized an estimated porosity for this report of 10 percent for dolomite bedrock (Freeze and Cherry, 1979). An effective porosity was then determined by multiplying the porosity by 0.90 (90 percent). Effective porosity is always equal to or less than the porosity and utilizing 90 percent is a conservative approach since part of the total porosity is occupied by static fluid held to the mineral surface by surface tension. The effective porosity determined was nine percent.

Area 1-2:

$$V_x = [(1.0x10^{-3} \text{ cm/sec})(0.05]/(0.09) = 5.52x10^{-4} \text{ cm/sec}$$ or **1.57 ft/day**

Area 1-3:

$$V_x = [(1.0x10^{-3} \text{ cm/sec})(0.120)]/(0.09) = 1.33x10^{-3} \text{ cm/sec}$$ or **3.78 ft/day**

3.2 STATISTICAL EVALUATION

The historical statistical database (provided on CD in **APPENDIX D**) was first reviewed for anomalies or outliers using the statistical program SANITAS™ for Groundwater. There were statistical outliers calculated for dissolved solids at MW-633D, manganese at NE-2, sulfate at NE-6, total organic carbon at NAB-3 and zinc at NE-6.

In addition to outlier analysis, parameter concentrations were plotted versus time for each of the parameters. These graphs are included in **APPENDIX E**. Graphs provide a summary of the historical data that are more easily visualized on graphs than tables. Another important application of these graphs is for detecting possible trends or drifts in the data from a given well. Furthermore, when visually comparing the plots of all the wells for a parameter, it is easy to identify the variability among the wells. This variability may be spatial or due to contamination from an off-site source.

An important consideration in any graphical presentation is whether the data is significantly influenced by seasonal changes. If this is the case, then the data should be adjusted for seasonal influences. In order to make such a determination, there should exist at least eight and preferably sixteen observations for each parameter. However, seasonal influences will likely be first suspected from visual observation of the data graphs discussed above. Based on the data generated thus far and as stated in previous reports, it is suspected that concentration levels are significantly influenced by the amount of precipitation and the timing of individual precipitation events in relation to sampling events.

The methods used to evaluate the groundwater data for statistically significant increases (SSIs) are based on statistical procedures outlined in the *Statistical Analysis of Ground-Water Monitoring Data at RCRA Facilities (Unified Guidance, March 2009)* and ASTM D6312-98 *Standard Guide for Developing Appropriate Statistical Approaches for Groundwater Detection Monitoring Programs*. The *SANITAS™ for Groundwater* program was utilized to statistically evaluate the data for the Fourth Quarter 2016 sampling event. A brief description of the procedures used in the statistical evaluation is provided on each statistical plot (See **APPENDIX E**).

3.2.1 Sen's Slope/Mann-Kendall

When used in conjunction with one another, the Mann-Kendall test for temporal trend and the Sen's slope estimate are two types of Evaluation Monitoring Statistics useful in determining the significance of an apparent trend and to estimate the magnitude of that trend. The Sen's Slope/Mann-Kendall was performed on all detected constituents from each well to determine whether a statistical trend is present in the data. The results of the Sen's Slope/Mann-Kendall plots associated with the Fourth Quarter 2016 sampling event are presented in **APPENDIX E**. A number of constituents presented a false trend due to a change in reporting limits. Only constituents with true statistical trends in detected concentrations are presented below:

Well	Significant Decreasing Trends	Significant Increasing Trends
CAO-1	manganese	arsenic
CAO-2	iron	chloride, manganese, total organic carbon
CAO-3	sulfate	
MW-1		arsenic, barium, chloride, iron, sulfate
MW-1R	manganese	cis-1,2-dichloroethene
MW-3		arsenic, barium, chloride, manganese, nickle
MW-4		chloride, zinc
MW-5	sulfate	chloride
MW-6		chloride
MW-7	chloride, sulfate	
MW-509D		sulfate
MW-577	barium, dissolved soilds, sulfate	chloride
MW-633D	sulfate	1,1-dichloroethane, zinc
MW-689D	barium, iron	chloride
NAB-2		chloride, zinc
NAB-4		barium,
NAB-8	barium,	
NE-2	chloride, dissolved solids, sulfate, total organic carbon	
NE-6	zinc	

3.2.2 Results of Assessment Monitoring Statistical Analyses

Confidence intervals are the recommended statistical strategy for assessment monitoring. The Groundwater Protection Standards (GWPS) utilized in the statistical evaluation are included in **TABLE 3.**

TABLE 3. Groundwater Protection Standards (Regulation 22.1205 (h)(l))

Compound	MCL**	RBSL***	Compound	MCL**	RBSL***
Total Arsenic	0.01 mg/l		Chloroethane (Ethyl Chloride)	--	21,000 ug/l
Total Barium	2 mg/l		Vinyl Chloride	2 ug/l	
Total Cadmium	0.005 mg/l		1,1-Dichloroethane	--	2.7 ug/l
Total Cobalt	--	0.011 mg/l	cis-1,2-Dichloroethene (-ethylene)	70 ug/l	--
Total Chromium	0.1 mg/l	--	Tetrachloroethene (-ethylene)	5 ug/l	
Total Nickel	--	0.039 mg/l	1,4-dichlorobenzene	75 ug/l	
Total Zinc	--	6 mg/l	Methylene Chloride	5 ug/l	
Total Selenium	0.05 mg/l		Toluene	1,000 ug/l	
Tin	--	12mg/l	Trans 1,2-dichloroethylene	100 ug/l	
Silver	--	0.094 mg/l	Xylene (Total)	10,000 ug/l	
Thallium	0.002 mg/l		Ethyl Benzene	700 ug/l	
Vanadium	--	0.086 mg/l	Carbon Disulfide	--	810 ug/l
Antimony	0.006 mg/l		Chlorobenzene	100 ug/l	
Beryllium	0.004 mg/l		Cyanide	200 ug/l	
Copper	1.3 mg/l		Mercury	200 ug/l	
Lead	0.015 mg/l		Trichloroethene (-ethylene)	5 ug/l	
Benzene	5 ug/l				

Available MCL's will be used as the Groundwater Protection Standard
**MCL = Maximum Contaminant Levels (current or proposed)
***RBSL = Risk-Based Screen Levels (EPA Region 6 Human Health Medium Specific Screening Level)
Residential Water (Residential Scenario: Ingestion and Inhalation)

Data evaluation consisted of the establishment of 95% Lower Confidence Limits (LCL), assuming that a minimum of four (4) background samples exist for each parameter and well detected during the Assessment Monitoring Program. If inadequate background data exists, sufficient background data will be collected to provide adequate sample size for statistical analysis. If the 95% LCL of one parameter exceeds action levels defined as Maximum Contaminant Levels (MCLs), if applicable, or a health-based alternate Groundwater Protection Standards (GWPS) as stipulated in Section 22.1205(d)(4), the site will conduct an Assessment of Corrective Measures in coordination with ADEQ.

Because the entire monitoring system is currently in the Assessment Monitoring Program, each detected volatile organic compound and metal parameter at each well was evaluated utilizing events from 2011 to present. This number of events was selected based on the interpretation that data population has been generally stable at each monitoring point over this time period. Data prior to these events varied substantially in some cases from the utilized data set. The statistical evaluation was conducted in

accordance with recommended procedures found in the Unified Guidance. Confidence Interval statistical analysis was not performed for indicator parameters.

The results of the assessment monitoring statistical evaluations indicated the concentrations of the following parameters at the following wells, statistically exceeded the established Groundwater Protection Standards.

FOURTH QUARTER 2016 GWPS EXCEEDANCES

MONITORING POINT	AMC
CAO-1	1,1-dichloroethane, arsenic, cobalt, vinyl chloride
MW-1	1,1-dichloroethane, cobalt, nickel
MW-1R	1,1-dichloroethane, arsenic, cobalt, nickel, vinyl chloride

The results of the Fourth Quarter 2016 Sampling Event indicate that AMC concentrations statistically exceed the GWPS at sample points CAO-1, MW-1, and MW-1R

The results of the confidence interval evaluation associated with the Fourth Quarter 2016 sampling event are presented in **APPENDIX E**. In accordance with Regulation 12.1205(d)(4), an *Assessment of Corrective Measures Report* (June 2015, Document 67822) was submitted to ADEQ for approval.

3.3 LABORATORY ANALYTICAL RESULTS

The analytical laboratory results for the Fourth Quarter 2016 Assessment Monitoring sampling event are summarized in **TABLE 4** and **TABLE 5**. The Tables present a comparison of parameter concentrations from the current sampling event to the applicable Primary Drinking Water Standards-Maximum Contaminant Levels (MCLs) and Secondary Drinking Water Standards (SDWS). The SDWS are set primarily for aesthetic reasons and are generally not considered health-based criteria. Constituents covered by the SDWS are those which may adversely affect the aesthetic qualities of drinking water such as taste, odor, color, and appearance and are not federally enforced.

Primary Drinking Water Standard MCL exceedances noted for well samples consisted of arsenic at MW-1, MW-1R, and CAO-1; cadmium at MW-509D; 1,1-dichloroethane at MW-1 and MW-1R; and vinyl chloride at MW-1R and CAO-1 (see TABLE 4 and TABLE 5) during the Fourth Quarter 2016 Assessment Monitoring event.

As shown in **TABLE 5**, Appendix 1 volatile organic compound (VOC) detections in well samples above the PQL consisted of the following:

- **MW-1 –** chlorobenzene, 1,1-dichloroethane, cis-1,2-dichloroethene, and vinyl chloride
- **MW-1R –** chlorobenzene, 1,4-dichlorobenzene, 1,1-dichloroethane, cis-1,2-dichloroethene, trichloroethene, and vinyl chloride
- **CAO-1 –** benzene, 1,1-dichloroethane, and vinyl chloride
- **MW-633D** - 1,1-dichloroethane

Because VOCs are not naturally occurring, detected concentrations above the PQL are considered SSIs.

3.4 QUALITY ASSURANCE/QUALITY CONTROL RESULTS

A QA/QC comparison for the Fourth Quarter 2016 Assessment Monitoring event is presented in **TABLES 4** and **5**. The duplicate sample was consistent with the representative sample during this event. The field and equipment blank had detections of toluene, while the trip blank had no VOC detections, during the Fourth Quarter 2016 Assessment Monitoring event

TABLE 4. INORGANIC GROUNDWATER QUALITY RESULTS

WELL ID	pH (SU)	TDS (mg/L)	Sulfide (mg/L)	Cyanide (mg/L)	Cl (mg/L)	SO_4 (mg/L)	TOC (mg/L)	Hg (mg/L)
MW-1	**6.14**	**972**	<0.05	<0.005	122	23.7	9.23	<0.0002
MW-1R	**6.15**	**928**	<0.05	<0.005	102	23.7	7.52	<0.0002
MW-2	6.60	381	<0.05	<0.005	3.61	14.7	0.519 J	<0.0002
MW-3	**6.32**	**525**	<0.05	<0.005	9.54	12.3	2.32	<0.0002
MW-4	**6.34**	376	<0.05	<0.005	7.31	14.9	1.09	<0.0002
MW-5	**6.27**	416	<0.05	<0.005	5.76	7.07	1.2	<0.0002
MW-6	6.50	432	<0.05	<0.005	21.0	7.76	1.75	<0.0002
Dupe (MW-6)	N/A	459	<0.05	<0.005	20.8	8.32	1.39	<0.0002
MW-7	**6.25**	338	<0.05	<0.005	1.50	6.39	1.07	<0.0002
CAO-1	**5.91**	**794**	<0.05	<0.005	95.9	0.104 J	10.6	<0.0002
CAO-2	**6.35**	433	<0.05	<0.005	66.2	0.261 J	5.77	<0.0002
CAO-3	6.60	305	<0.05	<0.005	17.9	6.01	1.65	<0.0002
NAB-2	**6.24**	443	<0.05	<0.005	10.4	13.5	1.39	<0.0002
NAB-3	7.75	425	<0.05	<0.005	9.17	8.97	94.7	0.000134 J
NAB-4	**6.26**	458	<0.05	<0.005	4.11	41.9	1.79	<0.0002
NAB-7	**6.30**	399	<0.05	<0.005	3.05	29.7	1.61	<0.0002
NAB-8	**4.00**	371	<0.05	<0.005	2.23	11.2	23.3	<0.0002
MW-509D	**6.20**	395	<0.05	<0.005	6.42	16.0	1.23	<0.0002
MW-577	**6.28**	398	<0.05	<0.005	2.48	20.2	1.14	<0.0002
MW-633D	6.41	366	<0.05	<0.005	15.4	10.6	1.53 B	<0.0002
MW-689D	**6.38**	408	<0.05	<0.005	2.49	14.2	0.579 J,B	<0.0002
NE-2	8.00	**1660**	<0.05	<0.005	24.9	737	12.4	<0.0002
NE-3	4.64	347	<0.05	<0.005	5.54	15.3	1.60	<0.0002
NE-6	6.32	398	<0.05	<0.005	2.31	22.4	0.663 J	<0.0002
Spring A	6.81	359	<0.05	<0.005	2.37	8.17	0.967 J	<0.0002
Class I Draw Sp	6.65	319	<0.05	<0.005	45	2.8 J	4.23	<0.0002
Landfill Ent Seep	6.95	335	<0.05	<0.005	12.7	4.41 J	2.38	<0.0002
SP-7	6.75	275	<0.05	<0.005	16.1	5.17	2.37	<0.0002
Field Blank	N/A	26	<0.05	<0.005	0.122 J	<5.0	0.397 J	<0.0002
LEACHATE	7.59	**1830**	<0.05	<0.005	**659**	11.2	60.0	<0.0002
EPA Standards	***6.5-8.5*****	***500*****	**---**	**---**	***250*****	***250*****	**---**	***0.002****

*Primary Drinking Water Standard-Maximum Contaminant Level (MCL)
**Secondary Drinking Water Standard (SDWS)
"J" Value= estimated concentration above the MDL but below the PQL
Values in **bold** exceed applicable Primary Drinking Water EPA Standards.
Values with a "B" suffix denotes the same analyte is found in the associated blank.

TABLE 4 (CONT'D). INORGANIC GROUNDWATER QUALITY RESULTS

WELL ID	Sb (mg/L)	As (mg/L)	Ba (mg/L)	Be (mg/L)	Cd (mg/L)	Cr (mg/L)	Cu (mg/L)	Co (mg/L)
MW-1	0.00115 J	**0.071**	0.201	<0.002	0.000654 J	0.00588	0.0104 B	0.109
MW-1R	<0.002	**0.0627**	0.173	<0.002	0.000239 J	0.00168 J	0.00493 J,B	0.0692
MW-2	<0.002	0.00466	0.0315	<0.002	<0.001	<0.002	0.000563 J,B	0.000295 J
MW-3	<0.002	0.00231	0.0812	<0.002	0.00043 J	0.000594 J	0.00108 J,B	0.0041
MW-4	<0.002	0.000505 J	0.0358	<0.002	0.000883 J	0.000895 J	0.00122 J,B	<0.002
MW-5	<0.002	0.000536 J	0.0341	<0.002	<0.001	0.00328	0.00377 J,B	<0.002
MW-6	<0.002	0.000423 J	0.0465	<0.002	0.000439 J	0.000805 J	0.00121 J,B	<0.002
Dupe (MW-6)	<0.002	0.000422 J	0.0482	<0.002	0.00047 J	<0.002	0.0016 J,B	<0.002
MW-7	<0.002	0.000687 J	0.0463	<0.002	0.000307 J	<0.002	0.0017 J,B	0.0056
CAO-1	<0.002	**0.123**	0.25	<0.002	<0.001	0.00132 J	0.000868 J,B	0.0286
CAO-2	<0.002	0.00135 J	0.0997	<0.002	0.00148	0.000821 J	0.0015 J,B	0.0111
CAO-3	<0.002	<0.002	0.0412	<0.002	0.000765 J	<0.002	0.00105 J,B	<0.002
NAB-2	0.000806 J	0.00628	0.0394	<0.002	<0.001	0.00577	0.00185 J,B	<0.002
NAB-3	<0.002	0.000561 J	0.0502	<0.002	0.000995 J	0.00121 J	0.00263 J,B	0.00125 J
NAB-4	<0.002	0.00458	0.0419	<0.002	0.00128	0.00571	0.004 J,B	0.00106 J
NAB-7	<0.002	0.00574	0.023	<0.002	<0.001	0.00744	0.00529 B	0.000774 J
NAB-8	<0.002	0.00354	0.0319	<0.002	0.000291 J	<0.002	0.000959 J,B	<0.002
MW-509D	<0.002	0.00136 J	0.034	<0.002	0.00621	0.00233	0.00262 J,B	0.000342 J
MW-577	<0.002	0.00139 J	0.0233	<0.002	<0.001	<0.002	0.000986 J,B	<0.002
MW-633D	<0.002	0.000414 J	0.041	<0.002	0.00168	0.00174 J	0.00351 J,B	<0.002
MW-689D	<0.002	0.000733 J	0.0277	<0.002	<0.001	0.00118 J	0.00256 J,B	<0.002
NE-2	<0.002	0.00199 J	0.0185	<0.002	<0.001	<0.002	0.00146 J,B	0.000603 J
NE-3	<0.002	0.000766 J	0.128	<0.002	0.000181 J	<0.002	0.000727 J,B	0.000394
NE-6	<0.002	0.00157 J	0.0328	<0.002	<0.001	0.00312	0.00189 J,B	0.000682 J
Spring A	<0.002	0.000375 J	0.0418	<0.002	0.00161	<0.002	0.00119 J,B	0.00169 J
Class I Draw Sp	<0.002	0.000799 J	0.0661	<0.002	<0.001	<0.002	0.000709 J,B	0.00108 J
Landfill Ent Seep	<0.002	0.003	0.0798	<0.002	<0.001	<0.002	<0.005	0.00241
SP-7	<0.002	0.00473	0.0727	<0.002	<0.001	<0.002	0.000792 J,B	0.00271
Field Blank	<0.002	<0.002	<0.005	<0.002	<0.001	<0.002	0.00619 B	<0.002
LEACHATE	<0.002	0.00912	**2.56**	<0.002	<0.001	0.0307	0.0122 B	0.00624
EPA Standards	***0.006****	***0.01****	***2****	***0.004****	***0.005****	***0.1****	***1.3****	**---**

*Primary Drinking Water Standard-Maximum Contaminant Level (MCL) / **Secondary Drinking Water Standard (SDWS)

"J" Value= estimated concentration above the MDL but below the PQL / Values in **bold** exceed applicable Primary Drinking Water EPA Standards.

Values with a "B" suffix denotes the same analyte is found in the associated blank.

TABLE 4 (CONT'D). INORGANIC GROUNDWATER QUALITY RESULTS

WELL ID	Fe (mg/L)	Pb (mg/L)	Mn (mg/L)	Ni (mg/L)	Se (mg/L)	Ag (mg/L)	Tl (mg/L)	Sn (mg/L)	V (mg/L)	Zn (mg/L)
MW-1	**6.3**	0.0029 B	**0.78**	0.158	<0.002	<0.002	0.000372 J	0.000693 J,B	0.00178 J	0.619
MW-1R	**23.7**	0.00241 B	**0.346**	0.0802	<0.002	<0.002	0.00164 J	0.00143 J,B	0.00045 J	1.6
MW-2	0.118	0.000636 J,B	0.0104	0.00979	<0.002	<0.002	<0.002	<0.002	0.000361 J	0.00366 J,B
MW-3	**0.473**	0.000937 J,B	**2.23**	0.0622	<0.002	<0.002	0.000661 J	<0.002	0.000426 J	0.146
MW-4	0.0411	0.000888 J,B	0.00209 J,B	0.114 J	<0.002	<0.002	<0.002	<0.002	0.0004 J	0.105
MW-5	0.0606 J	0.000765 J,B	0.00185 J,B	0.00281	<0.002	<0.002	<0.002	0.000313 J,B	0.000385 J	0.0368 B
MW-6	0.0334 J	0.000601 J,B	0.000867 J,B	0.000957 J	<0.002	<0.002	<0.002	0.000311 J,B	0.000427 J	0.0415 B
Dupe (MW-6)	0.0637 J,B	0.000635 J,B	0.00134 J,B	0.000804 J	<0.002	<0.002	<0.002	0.000326 J,B	0.000306 J	0.0416
MW-7	0.153	0.00153 J,B	**0.363**	0.0106	<0.002	<0.002	<0.002	0.000318 J,B	0.000497 J	0.0151 J,B
CAO-1	**36.6**	0.000886 J,B	**1.2**	0.0571	<0.002	<0.002	<0.002	0.000812 J,B	0.00115 J	0.0272 B
CAO-2	**0.305**	0.0013 J,B	**4.35**	0.0438	<0.002	<0.002	0.000612 J	0.000432 J,B	0.000711 J	0.0429 B
CAO-3	0.039 J	0.00106 J,B	0.0084	0.00196 J	<0.002	<0.002	<0.002	0.000506 J,B	0.000514 J	0.0213 J,B
NAB-2	**0.536**	0.00316 B	0.0175	0.00586	<0.002	<0.002	<0.002	0.000638 J,B	0.00128 J	0.0602
NAB-3	0.108	0.000842 J,B	0.0483	0.00268	<0.002	<0.002	<0.002	0.000407 J,B	0.000542 J	0.0823
NAB-4	0.247	0.00148 J,B	0.0166	0.0187	<0.002	<0.002	0.000547 J	0.000531 J,B	0.000569 J	0.656
NAB-7	**0.4**	0.00555 B	0.00701	0.0195	<0.002	<0.002	<0.002	<0.002	<0.000359 J	0.043 B
NAB-8	0.0378 J	0.000583 J,B	0.00405 J	<0.002	<0.002	<0.002	<0.002	<0.002	0.000351 J	0.00799 J,B
MW-509D	**0.882**	0.0029 B	0.0149	0.00283	<0.002	<0.002	<0.002	0.000307 J,B	0.00256 J	0.457
MW-577	0.189	0.000479 J,B	0.00484 J	0.0048	<0.002	<0.002	<0.002	<0.002	0.000297 J	<0.025
MW-633D	0.0548 J	0.000434 J,B	0.0035 J	0.00465	<0.002	<0.002	<0.002	0.000552 J,B	0.000561 J	0.283
MW-689D	0.143	0.00124 J,B	0.00803	0.00153 J	<0.002	<0.002	<0.002	0.000472 J,B	0.000508 J	0.00523 J,B
NE-2	**0.6**	0.00123 J,B	0.022	0.00719	0.000467 J	<0.002	<0.002	0.000305 J,B	0.000924 J	0.0797
NE-3	0.062 J,B	0.000515 J,B	0.0329	0.00364	<0.002	<0.002	<0.002	0.000479 J,B	0.00046 J	0.0411
NE-6	**0.345**	0.00102 J,B	0.0153	0.00782	<0.002	<0.002	<0.002	0.000323 J,B	0.000521 J	0.202
Spring A	0.118 B	0.000597 J,B	**0.327**	0.00446	<0.002	<0.002	<0.002	<0.002	0.000706 J	0.122
Class I Draw Sp	**0.567**	0.000762 J,B	**0.464**	0.00333	<0.002	<0.002	<0.002	0.000322 J,B	0.000474 J	0.00299 J,B
Landfill Ent Seep	**2.01**	0.000245 J,B	**0.914**	0.00236	<0.002	<0.002	<0.002	0.000379 J,B	0.000288 J	<0.025
SP-7	**2.05**	0.000428 J,B	**1.14**	0.00335	<0.002	<0.002	<0.002	<0.002	0.000281 J	0.00614 J,B
Field Blank	0.0466 J,B	0.000543 J,B	0.000294 J,B	<0.002	<0.002	<0.002	<0.002	0.000331 J,B	<0.005	<0.025
LEACHATE	**49.4**	0.00315 B	**1.22**	0.0499	0.000627 J	<0.002	<0.002	0.00632	0.00208 J	0.0666
EPA Standards	***0.3*****	***0.015****	***0.05*****	**---**	***0.05****	***0.1*****	***0.002****	**---**	**---**	***5*****

*Primary Drinking Water Standard-Maximum Contaminant Level (MCL) / **Secondary Drinking Water Standard (SDWS)

"J" Value= estimated concentration above the MDL but below the PQL / Values in **bold** exceed applicable Primary Drinking Water EPA Standards.

Values with a "B" suffix denotes the same analyte is found in the associated blank.

TABLE 5. VOLATILE ORGANIC GROUNDWATER QUALITY RESULTS

WELL ID	Acetone (ug/L)	Benzene (ug/L)	ChlBenz (ug/L)	ClEthane (ug/L)	1,4-DCB (ug/L)	1,1-DCA (ug/L)	1,2-DCA (ug/L)	CisDCEE (ug/L)	TranDCEE (ug/L)	TCE (ug/L)	VC (ug/L)	Xylene (ug/L)	Toluene (ug/L)	(MIBK) (ug/L)
MW-1	<10	0.572 J	1.2	<5.0	<1.0	**11.1**	0.399 J	7.76	<1.0	0.607 J	1.37	<3.0	<1.0	<10.0
MW-1R	<10	0.857 J	1.46	2.07 J	2.63	**16.2**	0.369 J	10.2	0.405 J	1.11	**3.69**	<3.0	<1.0	<10.0
MW-2	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
MW-3	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
MW-4	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
MW-5	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
MW-6	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
Dupe (MW-6)	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
MW-7	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
CAO-1	<10	4.34	0.849 J	2.0 J	<1.0	5.43	<1.0	<1.0	<1.0	<1.0		<3.0	0.443 J	<10.0
CAO-2	<10	<1.0	<1.0	<5.0	<1.0	0.488 J	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
CAO-3	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
NAB-2	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
NAB-3	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
NAB-4	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
NAB-7	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
NAB-8	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
MW-509D	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
MW-577	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
MW-633D	<10	<1.0	<1.0	<5.0	<1.0	1.97	<1.0	0.607 J	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
MW-689D	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
NE-2	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0	<10.0
NE-3	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0
NE-6	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0
Spring A	<10	<1.0	<1.0	<5.0	<1.0	2.92	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0
Class I Draw Sp	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0
Landfill Ent Seep	<10	<1.0	<1.0	2.81 J	<1.0	0.862 J	<1.0	0.477 J	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0
SP-7	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0
Field Blank	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	0.865 J
LEACHATE	23.6	<1.0	<1.0	0.565 J	1.4	0.565 J	<1.0	<1.0	<1.0	<1.0	<1.0	0.463 J	2.04 J	<1.0
Trip Blank	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	<1.0
Equipment Blank	<10	<1.0	<1.0	<5.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<3.0	0.82 J
EPA Standards	**---**	***5***	***75***	**---**	***75***	***7***	***5***	***70***	***100***	***5***	***2***	***10,000***	***1,000***	**---**

EPA standards depicted above are Primary Drinking Water Standard MCLs / "J" Value= estimated concentration above the MDL but below the PQL

Values in **bold** exceed applicable Primary Drinking Water EPA Standards / Values with a "B" suffix denotes the same analyte is found in the associated blank.

4 CONCLUSION

Based on the results of the Fourth Quarter 2016 Assessment Monitoring and analytical testing, SCS reached the following conclusions:

Groundwater Flow:

- ***FIGURE 2*** *represents a potentiometric surface map constructed from water levels measured during the Fourth Quarter 2016 Assessment Monitoring event. As indicated, groundwater within the uppermost aquifer was found to flow in a west-northwesterly direction in the Area 1-2 portion of the site. This flow relationship is consistent with the flow direction indicated by historical water level measurements. The groundwater flow direction in the Area 1-3 portion of the site is to the east-southeast. The average linear velocity in Area 1-2 is estimated at* $5.52x10^{-4}$ *cm/sec or* ***1.57 ft/day****. The average linear velocity in Area 1-3 is estimated at* $1.33x10^{-3}$ *cm/sec or* ***3.78 ft/day****.*

Analytical Results:

- *A QA/QC comparison for the Fourth Quarter 2016 Assessment Monitoring event is presented in* ***TABLES 4*** *and* ***5****. The duplicate sample was consistent with the representative sample during this event. The field and equipment blank had detections of toluene, while the trip blank had no VOC detections, during the Fourth Quarter 2016 Assessment Monitoring event*

- *Primary Drinking Water Standard MCL exceedances noted for well samples consisted of arsenic at MW-1, MW-1R, and CAO-1; cadmium at MW-509D; 1,1-dichloroethane at MW-1 and MW-1R; and vinyl chloride at MW-1R and CAO-1 (see TABLE 4 and TABLE 5) during the Fourth Quarter 2016 Assessment Monitoring event.*
- *As shown in* ***TABLE 5****, Appendix 1 volatile organic compound (VOC) detections in well samples above the PQL consisted of the following:*
    - ***MW-1 –*** *chlorobenzene, 1,1-dichloroethane, cis-1,2-dichloroethene, and vinyl chloride*
    - ***MW-1R –*** *chlorobenzene, 1,4-dichlorobenzene, 1,1-dichloroethane, cis-1,2-dichloroethene, trichloroethene, and vinyl chloride*
    - ***CAO-1 –*** *benzene, 1,1-dichloroethane, and vinyl chloride*
    - ***MW-633D*** *- 1,1-dichloroethane*

Statistical Evaluation:

- *The results of the assessment monitoring statistical evaluations indicated the concentrations of the following parameters at the following wells, statistically exceeded the established groundwater protection standards (****TABLE 3****).*

MONITORING POINT	***AMC***
CAO-1	*1,1-dichloroethane, arsenic, cobalt, vinyl chloride*
MW-1	*1,1-dichloroethane, cobalt, nickel*
MW-1R	*1,1-dichloroethane, arsenic, cobalt, nickel, vinyl chloride*

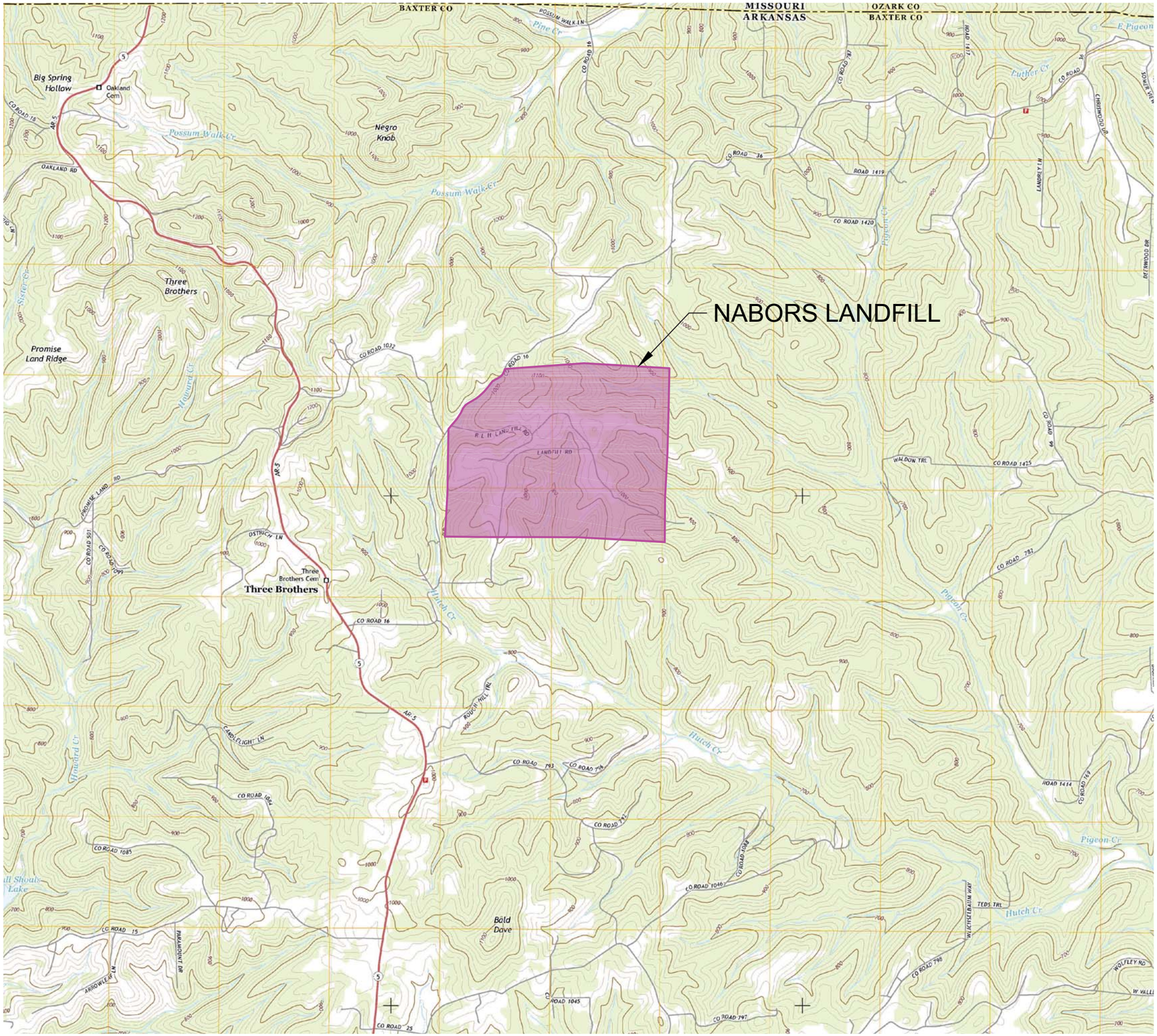
- *The results of the Fourth Quarter 2016 Sampling Event indicate that AMC concentrations statistically exceed the GWPS at sample points CAO-1, MW-1, and MW-1R*

- *In accordance with Regulation 12.1205(d)(4), an Assessment of Corrective Measures Report (June 2015, Document 67822) was submitted to ADEQ for approval.*

- *The next quarterly sampling event is scheduled for March 2017.*

APPENDIX A

FIGURES



NOTES:

1. BACKGROUND IMAGERY IS THE MIDWAY QUADRANGLE ARKANSAS-MISSOURI 7.5-MINUTE SERIES TOPO MAP RETRIEVED FROM WWW.USGS.GOV ON APRIL 27, 2016.

2. LANDFILL BOUNDARY IS APPROXIMATE.

3000 0 3000 6000

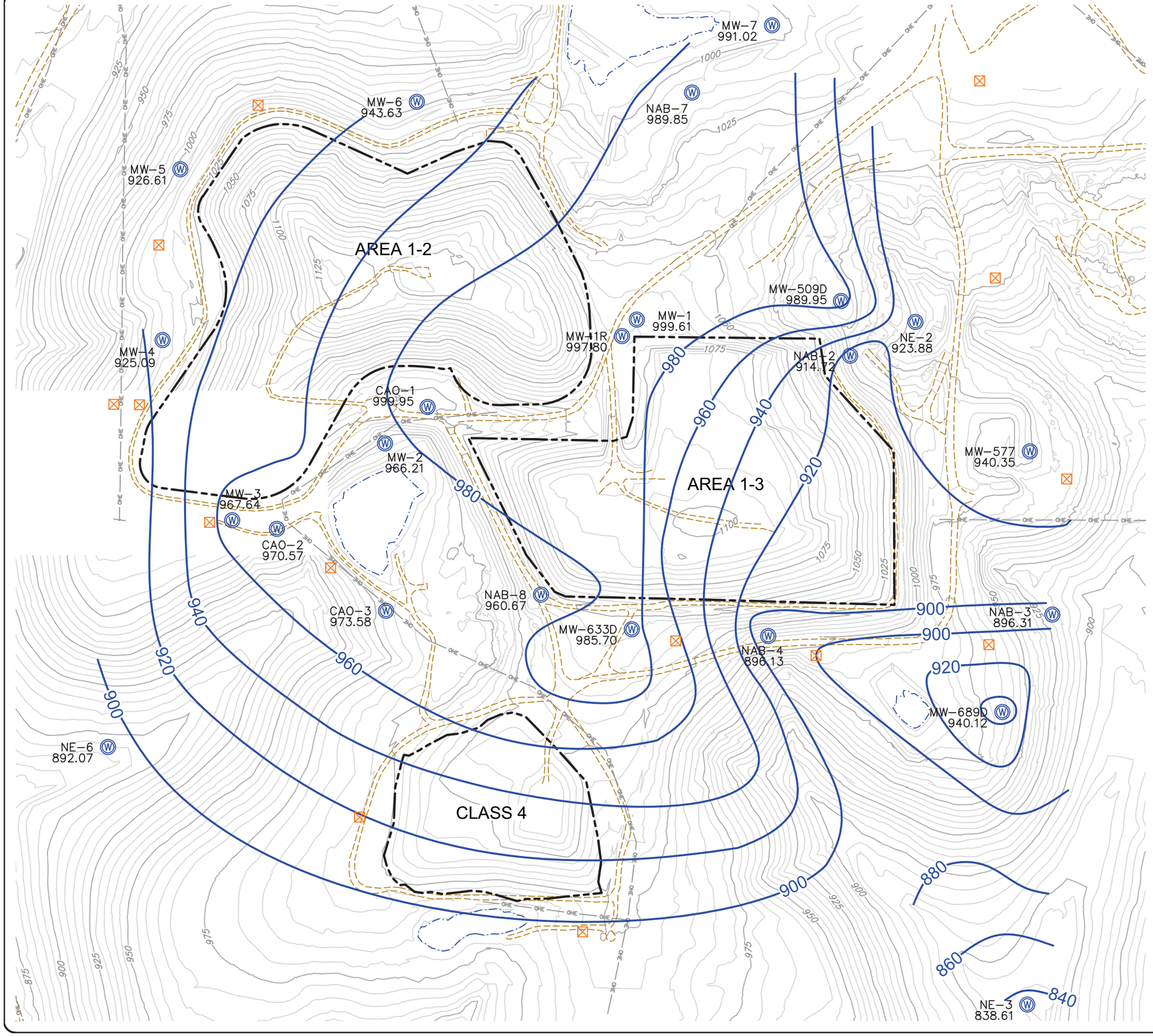
SCALE FEET

FIGURE 1 - SITE LOCATION MAP
NABORS LANDFILL
BAXTER COUNTY, ARKANSAS

SCS ENGINEERS

11219 Richardson Drive
North Little Rock, Arkansas 72113
PH: (501) 503-4781

CHK. BY: KV	DWN. BY: DMB	DSN. BY: DMB	PROJ. NO. --
PROJ. MGR: DM	DATE: 4/27/16	CADD FILE: SITE LOCATION MAP.DWG	



LEGEND:

- EXISTING 5' MINOR CONTOUR
- 1025 EXISTING 25' MAJOR CONTOUR
- 920 GROUNDWATER CONTOUR
- EXISTING ROAD
- OHE OHE OHE EXISTING OVERHEAD ELECTRIC
- DISPOSAL BOUNDARY (APPROXIMATE)
- EXISTING BODY OF WATER
- (W) EXISTING GROUNDWATER WELL
- EXISTING GAS PROBE

NOTES:

1. EXISTING TOPOGRAPHY BASED ON AERIAL SURVEY PERFORMED BY M.J. HARDEN ASSOCIATES, INC ON AUGUST 17, 2010. UPDATES TO THE TOPOGRAPHY AND SITE FEATURES FOR AREA 1-2, AREA 1-3, AND CLASS 4 DISPOSAL UNITS SURVEYED BY CONSOLIDATED LAND SERVICES, INC BETWEEN DECEMBER 2014 AND JANUARY 2015.
2. GROUNDWATER ELEVATIONS TAKEN ON DECEMBER 19-21, 2016.

300 0 300 600

SCALE FEET

FIGURE 2 - POTENTIOMETRIC MAP
4th QUARTER 2016 GROUNDWATER MONITORING
NABORS LANDFILL
BAXTER COUNTY, ARKANSAS

SCS ENGINEERS
11219 Richards Drive
North Little Rock, Arkansas 72113
PH. (501) 812-4551

CHK. BY: KAV	DWN. BY: DMB	DSN. BY: DMB	PROJ. NO. 27214218.01
PROJ. MGR: DM	DATE: 3/17/17	CADD FILE: POTENTIOMETRIC MAP.DWG	

APPENDIX B

GROUNDWATER SAMPLING RECORDS

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/20/2016 **Well No.** MW-1R

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** ok **Locked?** yes

Well Depth 78.55 **DTW (from TOC)** 69.77 **Volume H2O in well** 1.4

Other Information

Sunny, 45

5 mph wind

sub-pump

started purge @ 1612

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/20/2016	1614	1.5	6.18	15.8	2.57	28.4	1516	320
	1620	3.0	6.17	16.5	2.02	17.9	1517	30.4
	1626	4.5	6.15	16.7	1.34	9.9	1522	56.5
	1629	6.0	6.14	16.8	1.01	8.7	1523	28.5

Sampling Date & Time 12/20/2016 @ 1629

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/20/2016 **Well No.** MW-1

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** ok **Locked?** no

Well Depth 77.20 **DTW (from TOC)** 67.65 **Volume H2O in well** 1.5

Other Information

Sunny, 45

5 mph wind

sub-pump

started purge @ 1523

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/20/2016	1528	1.5	6.42	16.3	1.51	73.7	1656	18.6
	1532	3.0	6.24	16.5	0.74	66.7	1652	16.1
	1536	4.5	6.22	16.4	0.99	60.1	1648	17.7
	1542	6.0	5.84	16.6	0.99	52.9	1654	773
	1558	12.0	6.15	16.8	1.39	43.6	1618	35.7

Sampling Date & Time 12/20/2016 @ 1558

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/20/2016 **Well No.** MW-2

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** ok **Locked?** yes

Well Depth 49.10" **DTW (from TOC)** 35.00 **Volume H2O in well** 2.2

Other Information

Sunny, 40°

5 mph wind

sub-pump

started purge @ 1647

Date	Time	Volume (gallons)	pH (SU)	Temp (◦C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/20/2016	1648	2.5	6.59	16.5	1.30	-22.2	603	7.85
	1650	5.0	6.62	16.5	0.72	-35.1	612	4.45
	1652	7.5	6.60	16.5	0.51	-40.9	622	2.22

Sampling Date & Time 12/20/2016 @ 1652

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/21/2016 **Well No.** MW-3

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** ok **Locked?** yes

Well Depth 39.60" **DTW (from TOC)** 26.84 **Volume H2O in well** 1.2

Other Information

Sunny, 30°

5 mph wind

sub pump

start purging @ 0759

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/21/2016	0800	1.5	6.49	17.3	2.03	-7.4	841	14.7
	0802	3.0	6.50	17.3	1.23	-5.2	864	3.24
	0804	4.5	6.32	17.3	1.17	-4.9	868	3.07

Sampling Date & Time 12/21/2016 @ 0804

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/21/2016 **Well No.** MW-4

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** ok **Locked?** yes

Well Depth 100.60 **DTW (from TOC)** 87.02 **Volume H2O in well** 2.2

Other Information

Sunny 30°

5 mph wind

sub-pump

start purging @ 0826

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/21/2016	0834	2.5	6.46	15.5	5.79	1.3	658	152
	0837	5.0	6.35	15.8	5.69	-1.5	655	66.6
	0842	7.5	6.32	15.7	5.77	-3.2	650	37.8
	0847	10.0	6.34	15.7	5.59	-2.5	655	10.7

Sampling Date & Time 12/21/2016 @ 0847

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** **12/21/2016** **Well No.** **MW-5**

Sampling Personnel **Darren Motley**

Casing Diameter **2.0"** **Condition of well** **ok** **Locked?** **yes**

Well Depth **89.75"** **DTW (from TOC)** ***77.77*** **Volume H2O in well** **1.9**

Other Information

Sunny 30°

5 mph wind

sub-pump

start purging @ 0911

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/21/2016	**0902**	**2.0**	**6.44**	**14.8**	**6.11**	**6.8**	**692**	**43.9**
	0907	**4.0**	**6.36**	**15.2**	**5.28**	**1.7**	**702**	**21.1**
	0911	**6.0**	**6.27**	**15.2**	**4.98**	**-1.3**	**702**	**8.47**

Sampling Date & Time **12/21/2016 @ 0911**

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/21/2016 **Well No.** MW-6

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** ok **Locked?** yes

Well Depth 68.90 **DTW (from TOC)** 56.75 **Volume H2O in well** 1.9

Other Information

Sunny, 35°

5 mph wind

sub-pump

start purging @ 0935

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/21/2016	0942	2.0	6.68	14.2	6.30	18.1	725	12.1
	0944	4.0	6.55	14.3	4.95	16.0	731	8.69
	0946	6.0	6.50	14.3	4.15	12.5	733	5.68

Sampling Date & Time 12/21/2016 @ 0946

Notes

Dup @ 0950

F.B @ 0955

E.B @ 0957

T.B @ 1000

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/20/2016 **Well No.** MW-7

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** ok **Locked?** yes

Well Depth 23.0 **DTW (from TOC)** 8.69 **Volume H2O in well** 2.3

Other Information

Sunny, 42°

5 mph wind

sub-pump

start purging @ 1426

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/20/2016	1431	2.5	6.25	15.1	6.33	29.8	575	18.8
	1432	5.0	6.33	15.2	5.39	29.3	576	44.9
	1435	7.5	6.29	15.3	5.86	28.0	577	15.3
	1438	10.0	6.25	15.4	5.86	28.1	578	8.4

Sampling Date & Time 12/20/2016 @ 1438

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/21/2016 **Well No.** CAO-1

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** ok **Locked?** yes

Well Depth 37.0 **DTW (from TOC)** 26.45 **Volume H2O in well** 1.7

Other Information

Sunny, 40°

5 mph wind

sub-pump

start purging @ 1006

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/21/2016	1011	2.0	6.32	16.6	1.92	45.8	1413	110
	1016	4.0	6.22	16.7	1.64	21.3	1455	15.4
	1022	6.0	5.91	16.7	1.46	8.4	1466	5.65

Sampling Date & Time 12/21/2016 @ 1022

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/21/2016 **Well No.** CAO-2

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** ok **Locked?** yes

Well Depth 39.65 **DTW (from TOC)** 28.23 **Volume H2O in well** 1.8

Other Information

Sunny 30°

5 mph wind

sub-pump

start purging @ 0730

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/21/2016	0736	2.0	6.52	17.2	2.27	45.3	713	35.0
	0739	4.0	6.48	17.1	2.87	40.3	710	40.7
	0744	6.0	6.39	17.0	1.99	38.2	456	20.5
	0748	8.0	6.35	16.5	1.16	36.4	489	12.7

Sampling Date & Time 12/21/2016 @ 0748

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/21/2016 **Well No.** CAO-3

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** ok **Locked?** yes

Well Depth 23.0 **DTW (from TOC)** 10.62 **Volume H2O in well** 2.0

Other Information

Sunny 30°

5 mph wind

sub-pump

start purging @ 0650

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/21/2016	0700	2.0	6.95	15.5	2.13	32.7	476	34.3
	0715	4.0	6.89	15.5	1.07	27.7	469	6.85
	0720	6.0	6.60	15.6	0.77	25.9	469	4.13

Sampling Date & Time 12/21/2016 @ 0720

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/20/2016 **Well No.** NAB-2

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** ok **Locked?** no

Well Depth 93.0 **DTW (from TOC)** 79.26 **Volume H2O in well** 2.2

Other Information

Cloudy 40°

5 mph wind

sub-pump

start purging @ 1221

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/20/2016	1226	2.5	6.40	15.5	1.60	55.0	683	82.6
	1229	5.0	6.36	15.7	1.06	52.4	714	56.1
	1233	7.5	6.31	15.8	0.60	39.8	718	45.3
	1237	10.0	6.39	15.8	0.52	38.2	722	37.9
	1240	12.5	6.30	15.8	0.46	35.5	730	28.0
	1242	15.0	6.28	15.8	0.41	33.3	737	33.5
	1247	17.5	6.26	16.0	0.43	32.6	749	25.9
	1253	20.0	6.24	16.0	0.44	34.1	744	19.8

Sampling Date & Time 12/20/2016 @ 1253

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/20/2016 **Well No.** NAB-3

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** ok **Locked?** yes

Well Depth 47.0 **DTW (from TOC)** 25.18 **Volume H2O in well** 3.5

Other Information

Sunny 12°

5 mph wind

sub-pump

start purging @ 0800

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/20/2016	0805	3.5	7.43	15.3	2.09	39.3	746	98.1
	0809	7.0	7.27	15.4	2.10	41.3	731	59.4
	0813	10.5	7.52	15.3	3.48	42.7	727	42.7
	0819	14.0	7.75	15.3	6.36	43.5	721	13.5

Sampling Date & Time 12/20/2016 @ 0819

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** **9/20/2016** **Well No.** **NAB-4**

Sampling Personnel **Darren Motley**

Casing Diameter **2.0"** **Condition of well** **ok** **Locked?** **yes**

Well Depth **132.0** **DTW (from TOC)** **108.07** **Volume H2O in well** **3.9**

Other Information

Sunny 23°

5 mph wind

sub-pump

start purging @ 0946

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/20/2016	1009	4.0	6.40	16.1	1.49	44.9	741	233
	1021	8.0	6.33	16.0	2.02	35.0	744	83.5
	1034	12.0	6.28	16.3	0.86	39.0	738	40.2
	1042	16.0	6.26	15.8	0.78	38.2	743	13.4

Sampling Date & Time **12/20/2016 @ 1042**

Notes

pad is lifted off of ground

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/20/2016 **Well No.** NAB-7

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** **Locked?**

Well Depth 44.0 **DTW (from TOC)** 22.51 **Volume H2O in well** 3.5

Other Information

Sunny 95°

5 mph wind

sub-pump

start purging @ 1447

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/20/2016	1455	3.5	6.46	16.0	1.01	18.8	626	27.2
	1505	7.0	6.25	16.0	0.77	18.9	632	3.23
	1515	10.5	6.30	16.0	0.90	19.2	636	2.71

Sampling Date & Time 12/20/2016 @ 1515

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** **12/19/2016** **Well No.** **NAB-8**

Sampling Personnel **Darren Motley**

Casing Diameter **2.0"** **Condition of well** **ok** **Locked?** **yes**

Well Depth **95.0** **DTW (from TOC)** **78.54** **Volume H2O in well** **2.6**

Other Information

Sunny 20°

5 mph wind

Bailer

start purging @ 1230

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/19/2016	**1240**	**3.0**	**8.61**	**11.9**	**3.90**	**72.8**	**605**	**21.3**
Dry @ 3 gallons								
12/20/2016	**0700**		**4.00**	**5.2**	**4.36**	**93.6**	**624**	**17.8**

Sampling Date & Time **12/20/2016 @ 0700**

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/20/2016 **Well No.** MW-509D

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** **Locked?**

Well Depth 39.65 **DTW (from TOC)** 24.25 **Volume H2O in well** 2.5

Other Information

Sunny, 42°

5 mph wind

sub-pump

start purging @ 1300

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/20/2016	1310	2.5	6.33	17.2	5.35	28.1	656	326
	1316	5.0	6.36	17.3	5.18	28.8	653	115
	1324	7.5	6.31	17.2	5.27	30.7	653	83.5
	1330	10.0	6.28	17.2	5.24	30.4	655	106
	1345	12.5	6.28	17.1	4.99	31.7	660	109
	1353	15.0	6.26	17.0	4.54	31.8	662	103
	1358	17.5	6.23	17.0	4.51	32.0	664	153
	1415	24.5	6.20	16.8	3.60	32.1	667	59.1

Sampling Date & Time 12/20/2016 @ 1415

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** **12/20/2016** **Well No.** **MW-577**

Sampling Personnel **Darren Motley**

Casing Diameter **2.0"** **Condition of well** **ok** **Locked?** **yes**

Well Depth **55.6** **DTW (from TOC)** **42.25** **Volume H2O in well** **2.1**

Other Information

Sunny, 32°

5 mph wind

sub-pump

start purging @ 1149

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/20/2016	1156	2.0	6.44	16.0	2.20	36.9	693	25.5
	1204	4.0	6.47	16.1	1.48	27.8	676	5.72
	1210	6.0	6.28	16.1	1.40	27.7	674	4.79

Sampling Date & Time **12/20/2016 @ 1210**

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/20/2016 **Well No.** MW-633D

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** OK **Locked?** YES

Well Depth 87.92 **DTW (from TOC)** 64.40 **Volume H2O in well** 3.8

Other Information

Sunny 30°

5 mph wind

sub-pump

start purging @ 1102

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/20/2016	1112	4.0	6.60	16.5	1.49	42.8	719	64.2
	1221	8.0	6.54	16.8	1.22	39.6	718	6.54
	1130	12.0	6.41	17.0	1.13	38.4	719	4.99

Sampling Date & Time 12/20/2016 @ 1130

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/20/2016 **Well No.** MW-689D

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** ok **Locked?** yes

Well Depth 44.45 **DTW (from TOC)** 26.08 **Volume H2O in well** 2.9

Other Information

Sunny 18°

5 mph wind

sub-pump

start purging @ 0836

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/20/2016	0845	3.0	10.81	16.8	2.24	38.0	663	551
	0851	6.0	6.43	16.8	1.33	34.6	669	141
	0859	9.0	6.24	16.9	0.92	34.1	669	69.3
	0906	12.0	6.30	16.8	2.36	32.3	668	43.4
	0913	15.0	6.35	16.9	0.62	32.0	669	31.0
	0920	18.0	6.30	16.9	0.52	32.2	671	25.4
	0925	21.0	6.33	16.9	0.448	32.7	672	20.1
	0931	24.0	6.38	16.9	0.42	32.1	672	16.8

Sampling Date & Time 12/20/2016 @ 0931

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** **12/19/2016** **Well No.** **NE-2**

Sampling Personnel **Darren Motley**

Casing Diameter **2.0"** **Condition of well** **ok** **Locked?** **yes**

Well Depth **64.0** **DTW (from TOC)** **53.10** **Volume H2O in well** **1.5**

Other Information

Sunny, 18°

5 mph wind

bailer

start purging @ 1150

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/19/2016	**1150**	**1.0**	**8.0**	**13.5**	**2.50**	**141.5**	**2192**	**19.2**

Sampling Date & Time **12/19/2016 @ 1150**

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** **12/19/2016** **Well No.** **NE-3**

Sampling Personnel **Darren Motley**

Casing Diameter **2.0"** **Condition of well** **ok** **Locked?** **yes**

Well Depth **27.7** **DTW (from TOC)** **8.30** **Volume H2O in well** **3.1**

Other Information

Sunny 20°

5 mph wind

Sub-pump

started purging @ 1210

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/19/2016	**1215**	**3.0**	**9.55**	**11.1**	**4.94**	**81.1**	**668**	**237**
Dry @ 4.5 gallons								
12/20/2016	**0730**		**4.64**	**7.1**	**5.36**	**88.2**	**592**	**13.8**

Sampling Date & Time **12/20/2016 @ 0730**

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** 12/21/2016 **Well No.** NE-6

Sampling Personnel Darren Motley

Casing Diameter 2.0" **Condition of well** ok **Locked?** yes

Well Depth 18.15 **DTW (from TOC)** 9.35 **Volume H2O in well** 1.4

Other Information

Sunny 30°

5 mph wind

sub-pump

start purging @ 0812

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/21/2016	0814	1.5	6.43	14.9	7.29	-0.1	606	11.8
	0817	3.0	6.35	15.5	2.53	-6.6	635	5.17
	0820	4.5	6.32	15.5	2.36	-6.9	623	3.36

Sampling Date & Time 12/21/2016 @ 0820

Notes

SCS FIELD SERVICES

Field Groundwater Sampling Record

Facility NABORS **Date** **12/20/2016** **Sample.** **Leachate**

Sampling Personnel **Darren Motley**

Other Information

Sunny 40°

5 mph wind

Date	Time	Volume (gallons)	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
12/20/2016	**1705**		***7.59***	**3.1**	**8.73**	**90.8**	**3605**	**156**

Sampling Date & Time **1705**

Notes

SCS FIELD SERVICES

Field Spring Sampling Record

Facility NABORS **Date** **12/21/2016**

Sampling Personnel **Darren Motley**

Other Information

Sunny 45

5 mph wind

Sample ID	Sample Date	Time	pH (SU)	Temp (°C)	D.O. (mg/L)	ORP (mV)	S.C (µSm)	Turbidity (NTU)
SP-NE-3	**1055**	**Dry**						
TSP-3	**1050**	**Dry**						
TSP-4	**1045**	**Dry**						
Spring-A	**1140**		**6.81**	**14.0**	**5.47**	**-25.6**	**573**	**6.28**
Class I Draw	**1105**		**6.65**	**7.1**	**5.67**	**-8.4**	**557**	**8.83**
Class IV Draw	**1110**	**Dry**						
SP-7	**1130**		**6.75**	**11.7**	**6.02**	**-16.9**	**472**	**23.4**
Spring B	**1100**	**Dry**						
Entrance Sweep	**1120**		**6.95**	**8.1**	**5.55**	**-14.2**	**576**	**10.8**

Notes

APPENDIX C

LABORATORY ANALYTICAL RESULTS

ANALYTICAL REPORT

January 03, 2017

SCS Engineers - Little Rock, AR

Sample Delivery Group:	L880396
Samples Received:	12/22/2016
Project Number:	
Description:	Nabors Landfill

Report To: Stacie Whitmer
11219 Richardson Drive
North Little Rock, AR 72113

Entire Report Reviewed By:

Mark W. Beasley
Technical Service Representative

MW-1 L880396-01 GW

Collected by: Darren Motley
Collected date/time: 12/20/16 15:58
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938474	1	12/27/16 13:24	12/27/16 15:37	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 14:50	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/29/16 23:04	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 04:17	12/26/16 04:17	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938032	1	12/22/16 22:24	12/22/16 22:24	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:29	DR
Wet Chemistry by Method 9056A	WG938352	1	12/24/16 16:17	12/24/16 16:17	NJM
Wet Chemistry by Method 9056A	WG938352	5	12/24/16 20:16	12/24/16 20:16	NJM
Wet Chemistry by Method 9060A	WG938825	1	12/29/16 12:12	12/29/16 12:12	SJM

MW-2 L880396-02 GW

Collected by: Darren Motley
Collected date/time: 12/20/16 16:52
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938474	1	12/27/16 13:24	12/27/16 15:37	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 14:57	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/29/16 23:07	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 04:41	12/26/16 04:41	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938032	1	12/22/16 22:24	12/22/16 22:24	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:30	DR
Wet Chemistry by Method 9056A	WG938352	1	12/24/16 16:32	12/24/16 16:32	NJM
Wet Chemistry by Method 9060A	WG938824	1	12/28/16 13:39	12/28/16 13:39	SJM

MW-3 L880396-03 GW

Collected by: Darren Motley
Collected date/time: 12/21/16 08:04
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938476	1	12/27/16 13:32	12/27/16 14:20	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 14:59	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/29/16 23:11	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 05:04	12/26/16 05:04	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938032	1	12/22/16 22:24	12/22/16 22:24	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:31	DR
Wet Chemistry by Method 9056A	WG938352	1	12/24/16 15:47	12/24/16 15:47	NJM
Wet Chemistry by Method 9060A	WG938824	1	12/28/16 13:50	12/28/16 13:50	SJM

MW-4 L880396-04 GW

Collected by: Darren Motley
Collected date/time: 12/21/16 08:47
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938476	1	12/27/16 13:32	12/27/16 14:20	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 15:02	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/29/16 23:14	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 05:27	12/26/16 05:27	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938032	1	12/22/16 22:24	12/22/16 22:24	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:32	DR
Wet Chemistry by Method 9056A	WG938352	1	12/24/16 17:32	12/24/16 17:32	NJM
Wet Chemistry by Method 9060A	WG938824	1	12/28/16 14:02	12/28/16 14:02	SJM

MW-5 L880396-05 GW

Collected by: Darren Motley
Collected date/time: 12/21/16 09:11
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938476	1	12/27/16 13:32	12/27/16 14:20	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 15:04	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/29/16 23:18	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 05:51	12/26/16 05:51	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938032	1	12/22/16 22:24	12/22/16 22:24	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:33	DR
Wet Chemistry by Method 9056A	WG938352	1	12/24/16 17:47	12/24/16 17:47	NJM
Wet Chemistry by Method 9060A	WG938824	1	12/28/16 14:13	12/28/16 14:13	SJM

MW-6 L880396-06 GW

Collected by: Darren Motley
Collected date/time: 12/21/16 09:46
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938476	1	12/27/16 13:32	12/27/16 14:20	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 15:06	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/29/16 23:22	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 06:14	12/26/16 06:14	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938032	1	12/22/16 22:25	12/22/16 22:25	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:36	DR
Wet Chemistry by Method 9056A	WG938352	1	12/24/16 18:31	12/24/16 18:31	NJM
Wet Chemistry by Method 9060A	WG938824	1	12/28/16 14:24	12/28/16 14:24	SJM

MW-7 L880396-07 GW

Collected by: Darren Motley
Collected date/time: 12/20/16 14:38
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938474	1	12/27/16 13:24	12/27/16 15:37	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 15:15	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/29/16 23:25	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 06:37	12/26/16 06:37	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938032	1	12/22/16 22:26	12/22/16 22:26	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:37	DR
Wet Chemistry by Method 9056A	WG938352	1	12/24/16 18:46	12/24/16 18:46	NJM
Wet Chemistry by Method 9060A	WG938824	1	12/28/16 14:35	12/28/16 14:35	SJM

CAO-1 L880396-08 GW

Collected by: Darren Motley
Collected date/time: 12/21/16 10:22
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938476	1	12/27/16 13:32	12/27/16 14:20	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 15:18	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/29/16 23:29	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 07:00	12/26/16 07:00	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938032	1	12/22/16 22:26	12/22/16 22:26	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:38	DR
Wet Chemistry by Method 9056A	WG938352	1	12/24/16 19:01	12/24/16 19:01	NJM
Wet Chemistry by Method 9060A	WG938825	2	12/29/16 12:37	12/29/16 12:37	SJM

CAO-2 L880396-09 GW

Collected by: Darren Motley
Collected date/time: 12/21/16 07:48
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938690	1	12/28/16 13:53	12/28/16 14:22	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 15:20	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/29/16 23:32	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 07:24	12/26/16 07:24	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938032	1	12/22/16 22:26	12/22/16 22:26	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:39	DR
Wet Chemistry by Method 9056A	WG938352	1	12/24/16 19:16	12/24/16 19:16	NJM
Wet Chemistry by Method 9060A	WG938824	1	12/28/16 14:58	12/28/16 14:58	SJM

CAO-3 L880396-10 GW

Collected by: Darren Motley
Collected date/time: 12/21/16 07:20
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938690	1	12/28/16 13:53	12/28/16 14:22	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 15:22	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/29/16 23:36	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 07:47	12/26/16 07:47	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938032	1	12/22/16 22:26	12/22/16 22:26	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:41	DR
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 16:01	12/27/16 16:01	KCF
Wet Chemistry by Method 9060A	WG938824	1	12/28/16 15:10	12/28/16 15:10	SJM

NAB-2 L880396-11 GW

Collected by: Darren Motley
Collected date/time: 12/20/16 12:53
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938474	1	12/27/16 13:24	12/27/16 15:37	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 15:24	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/29/16 23:50	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 08:10	12/26/16 08:10	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938032	1	12/22/16 22:26	12/22/16 22:26	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:42	DR
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 16:21	12/27/16 16:21	KCF
Wet Chemistry by Method 9060A	WG938824	1	12/28/16 16:16	12/28/16 16:16	SJM

NAB-3 L880396-12 GW

Collected by: Darren Motley
Collected date/time: 12/20/16 08:19
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938474	1	12/27/16 13:24	12/27/16 15:37	JER
Mercury by Method 7470A	WG938411	1	12/24/16 05:59	12/27/16 14:46	NJB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/29/16 23:54	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 08:33	12/26/16 08:33	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938032	1	12/22/16 22:26	12/22/16 22:26	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:43	DR
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 16:31	12/27/16 16:31	KCF
Wet Chemistry by Method 9060A	WG938824	1	12/28/16 16:41	12/28/16 16:41	SJM

NAB-4 L880396-13 GW

Collected by: Darren Motley
Collected date/time: 12/20/16 10:42
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938474	1	12/27/16 13:24	12/27/16 15:37	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 15:27	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/29/16 23:57	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 08:57	12/26/16 08:57	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938032	1	12/22/16 22:27	12/22/16 22:27	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:44	DR
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 16:41	12/27/16 16:41	KCF
Wet Chemistry by Method 9060A	WG938824	1	12/28/16 16:55	12/28/16 16:55	SJM

NAB-7 L880396-14 GW

Collected by: Darren Motley
Collected date/time: 12/20/16 15:15
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938474	1	12/27/16 13:24	12/27/16 15:37	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 15:29	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/30/16 00:01	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 09:20	12/26/16 09:20	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938032	1	12/22/16 22:27	12/22/16 22:27	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:47	DR
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 16:52	12/27/16 16:52	KCF
Wet Chemistry by Method 9060A	WG938824	1	12/28/16 17:06	12/28/16 17:06	SJM

NAB-8 L880396-15 GW

Collected by: Darren Motley
Collected date/time: 12/20/16 07:00
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938474	1	12/27/16 13:24	12/27/16 15:37	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 15:31	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/30/16 00:04	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 09:43	12/26/16 09:43	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938033	1	12/22/16 22:09	12/22/16 22:09	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:48	DR
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 17:02	12/27/16 17:02	KCF
Wet Chemistry by Method 9060A	WG938824	1	12/28/16 17:32	12/28/16 17:32	SJM

MW-509D L880396-16 GW

Collected by: Darren Motley
Collected date/time: 12/20/16 14:15
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938474	1	12/27/16 13:24	12/27/16 15:37	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 15:33	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/30/16 00:08	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 10:07	12/26/16 10:07	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938033	1	12/22/16 22:10	12/22/16 22:10	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:51	DR
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 17:42	12/27/16 17:42	KCF
Wet Chemistry by Method 9060A	WG938824	1	12/28/16 17:43	12/28/16 17:43	SJM

MW-577 L880396-17 GW

Collected by: Darren Motley
Collected date/time: 12/20/16 12:10
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938474	1	12/27/16 13:24	12/27/16 15:37	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 15:36	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/30/16 00:12	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 10:30	12/26/16 10:30	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938033	1	12/22/16 22:10	12/22/16 22:10	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:52	DR
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 17:53	12/27/16 17:53	KCF
Wet Chemistry by Method 9060A	WG938824	1	12/28/16 17:55	12/28/16 17:55	SJM

MW-689D L880396-18 GW

Collected by: Darren Motley
Collected date/time: 12/20/16 09:31
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938475	1	12/27/16 13:28	12/27/16 14:58	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 15:42	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/30/16 00:15	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 10:53	12/26/16 10:53	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938033	1	12/22/16 22:10	12/22/16 22:10	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:53	DR
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 18:03	12/27/16 18:03	KCF
Wet Chemistry by Method 9060A	WG938825	1	12/29/16 12:57	12/29/16 12:57	SJM

MW-633D L880396-19 GW

Collected by: Darren Motley
Collected date/time: 12/20/16 11:30
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938475	1	12/27/16 13:28	12/27/16 14:58	JER
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 15:45	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/30/16 00:19	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 11:17	12/26/16 11:17	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938033	1	12/22/16 22:11	12/22/16 22:11	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:54	DR
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 18:13	12/27/16 18:13	KCF
Wet Chemistry by Method 9060A	WG938825	1	12/29/16 13:46	12/29/16 13:46	SJM

NE-2 L880396-20 GW

Collected by: Darren Motley
Collected date/time: 12/19/16 11:50
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938167	1	12/24/16 05:22	12/25/16 00:15	JM
Mercury by Method 7470A	WG938137	1	12/23/16 09:21	12/23/16 15:47	TRB
Metals (ICPMS) by Method 6020	WG938047	1	12/27/16 11:18	12/29/16 22:24	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938408	1	12/26/16 11:40	12/26/16 11:40	ACG
Wet Chemistry by Method 4500S2 D-2011	WG938033	1	12/22/16 22:11	12/22/16 22:11	MZ
Wet Chemistry by Method 9012B	WG938805	1	12/28/16 09:52	12/29/16 09:58	DR
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 18:23	12/27/16 18:23	KCF
Wet Chemistry by Method 9056A	WG939061	20	12/28/16 12:47	12/28/16 12:47	KCF
Wet Chemistry by Method 9060A	WG938825	1	12/29/16 14:15	12/29/16 14:15	SJM

NE-3 L880396-21 GW

Collected by: Darren Motley
Collected date/time: 12/20/16 07:30
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938475	1	12/27/16 13:28	12/27/16 14:58	JER
Mercury by Method 7470A	WG938138	1	12/23/16 09:22	12/23/16 17:37	TRB
Metals (ICPMS) by Method 6020	WG938049	1	12/27/16 13:37	12/30/16 01:36	VSS
Metals (ICPMS) by Method 6020	WG939602	1	12/30/16 07:06	12/30/16 10:18	JPD
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938588	1	12/26/16 00:57	12/26/16 00:57	JHH
Wet Chemistry by Method 4500S2 D-2011	WG938033	1	12/22/16 22:11	12/22/16 22:11	MZ
Wet Chemistry by Method 9012B	WG939301	1	12/29/16 14:18	12/29/16 17:43	ASK
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 18:33	12/27/16 18:33	KCF
Wet Chemistry by Method 9060A	WG939610	1	12/31/16 11:59	12/31/16 11:59	SJM

NE-6 L880396-22 GW

Collected by: Darren Motley
Collected date/time: 12/21/16 08:20
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938690	1	12/28/16 13:53	12/28/16 14:22	JER
Mercury by Method 7470A	WG938138	1	12/23/16 09:22	12/23/16 17:39	TRB
Metals (ICPMS) by Method 6020	WG938049	1	12/27/16 13:37	12/30/16 01:40	VSS
Metals (ICPMS) by Method 6020	WG939602	1	12/30/16 07:06	12/30/16 10:22	JPD
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938588	1	12/26/16 01:21	12/26/16 01:21	JHH
Wet Chemistry by Method 4500S2 D-2011	WG938033	1	12/22/16 22:12	12/22/16 22:12	MZ
Wet Chemistry by Method 9012B	WG939301	1	12/29/16 14:18	12/29/16 17:44	ASK
Wet Chemistry by Method 9056A	WG939061	1	12/28/16 12:57	12/28/16 12:57	KCF
Wet Chemistry by Method 9060A	WG939610	1	12/31/16 12:34	12/31/16 12:34	SJM

MW-1R L880396-23 GW

Collected by: Darren Motley
Collected date/time: 12/20/16 16:29
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938475	1	12/27/16 13:28	12/27/16 14:58	JER
Mercury by Method 7470A	WG938138	1	12/23/16 09:22	12/23/16 17:41	TRB
Metals (ICPMS) by Method 6020	WG938049	1	12/27/16 13:37	12/30/16 01:43	VSS
Metals (ICPMS) by Method 6020	WG939602	2	12/30/16 07:06	12/30/16 11:12	JPD
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938588	1	12/26/16 01:45	12/26/16 01:45	JHH
Wet Chemistry by Method 4500S2 D-2011	WG938033	1	12/22/16 22:12	12/22/16 22:12	MZ
Wet Chemistry by Method 9012B	WG939301	1	12/29/16 14:18	12/29/16 17:47	ASK
Wet Chemistry by Method 9056A	WG938354	2	12/27/16 19:34	12/27/16 19:34	KCF
Wet Chemistry by Method 9060A	WG939610	1	12/31/16 12:52	12/31/16 12:52	SJM

SP-7 L880396-24 GW

Collected by: Darren Motley
Collected date/time: 12/21/16 11:30
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938690	1	12/28/16 13:53	12/28/16 14:22	JER
Mercury by Method 7470A	WG938138	1	12/23/16 09:22	12/23/16 17:43	TRB
Metals (ICPMS) by Method 6020	WG938049	1	12/27/16 13:37	12/30/16 01:47	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938588	1	12/26/16 02:09	12/26/16 02:09	JHH
Wet Chemistry by Method 4500S2 D-2011	WG938033	1	12/22/16 22:12	12/22/16 22:12	MZ
Wet Chemistry by Method 9012B	WG939301	1	12/29/16 14:18	12/29/16 17:48	ASK
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 19:44	12/27/16 19:44	KCF
Wet Chemistry by Method 9060A	WG939610	1	12/31/16 13:10	12/31/16 13:10	SJM

SPRING-A L880396-25 GW

Collected by: Darren Motley
Collected date/time: 12/21/16 11:40
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938690	1	12/28/16 13:53	12/28/16 14:22	JER
Mercury by Method 7470A	WG938138	1	12/23/16 09:22	12/23/16 17:46	TRB
Metals (ICPMS) by Method 6020	WG938049	1	12/27/16 13:37	12/30/16 01:50	VSS
Metals (ICPMS) by Method 6020	WG939602	1	12/30/16 07:06	12/30/16 10:36	JPD
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938676	1	12/26/16 15:30	12/26/16 15:30	BMB
Wet Chemistry by Method 4500S2 D-2011	WG938033	1	12/22/16 22:12	12/22/16 22:12	MZ
Wet Chemistry by Method 9012B	WG939301	1	12/29/16 14:18	12/29/16 17:49	ASK
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 19:54	12/27/16 19:54	KCF
Wet Chemistry by Method 9060A	WG939610	1	12/31/16 13:29	12/31/16 13:29	SJM

LANDFILL ENTRANCE SEEP L880396-26 GW

Collected by: Darren Motley
Collected date/time: 12/21/16 11:20
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938690	1	12/28/16 13:53	12/28/16 14:22	JER
Mercury by Method 7470A	WG938138	1	12/23/16 09:22	12/23/16 17:48	TRB
Metals (ICPMS) by Method 6020	WG938049	1	12/27/16 13:37	12/30/16 01:54	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938676	1	12/26/16 15:43	12/26/16 15:43	BMB
Wet Chemistry by Method 4500S2 D-2011	WG938033	1	12/22/16 22:13	12/22/16 22:13	MZ
Wet Chemistry by Method 9012B	WG939301	1	12/29/16 14:18	12/29/16 17:52	ASK
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 20:05	12/27/16 20:05	KCF
Wet Chemistry by Method 9060A	WG939610	1	12/31/16 13:50	12/31/16 13:50	SJM

CLASS I DRAW L880396-27 GW

Collected by: Darren Motley
Collected date/time: 12/21/16 11:05
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938690	1	12/28/16 13:53	12/28/16 14:22	JER
Mercury by Method 7470A	WG938138	1	12/23/16 09:22	12/23/16 17:50	TRB
Metals (ICPMS) by Method 6020	WG938049	1	12/27/16 13:37	12/30/16 01:57	VSS
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938676	1	12/26/16 15:57	12/26/16 15:57	BMB
Wet Chemistry by Method 4500S2 D-2011	WG938033	1	12/22/16 22:14	12/22/16 22:14	MZ
Wet Chemistry by Method 9012B	WG939301	1	12/29/16 14:18	12/29/16 17:54	ASK
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 20:15	12/27/16 20:15	KCF
Wet Chemistry by Method 9060A	WG939610	1	12/31/16 15:26	12/31/16 15:26	SJM

LEACHATE L880396-28 GW

Collected by: Darren Motley
Collected date/time: 12/20/16 17:05
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938475	1	12/27/16 13:28	12/27/16 14:58	JER
Mercury by Method 7470A	WG938138	1	12/23/16 09:22	12/23/16 17:52	TRB
Metals (ICPMS) by Method 6020	WG938049	1	12/27/16 13:37	12/30/16 02:01	VSS
Metals (ICPMS) by Method 6020	WG938049	5	12/27/16 13:37	12/30/16 02:24	VSS
Metals (ICPMS) by Method 6020	WG939602	1	12/30/16 07:06	12/30/16 10:39	JPD
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938676	1	12/26/16 16:10	12/26/16 16:10	BMB
Wet Chemistry by Method 4500S2 D-2011	WG938033	1	12/22/16 22:14	12/22/16 22:14	MZ
Wet Chemistry by Method 9012B	WG939301	1	12/29/16 14:18	12/29/16 17:55	ASK
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 20:25	12/27/16 20:25	KCF
Wet Chemistry by Method 9056A	WG938354	10	12/27/16 20:35	12/27/16 20:35	KCF
Wet Chemistry by Method 9060A	WG939610	1	12/31/16 15:46	12/31/16 15:46	SJM

DUPLICATE L880396-29 GW

Collected by: Darren Motley
Collected date/time: 12/21/16 09:50
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938690	1	12/28/16 13:53	12/28/16 14:22	JER
Mercury by Method 7470A	WG938138	1	12/23/16 09:22	12/23/16 17:55	TRB
Metals (ICPMS) by Method 6020	WG938049	1	12/27/16 13:37	12/30/16 02:05	VSS
Metals (ICPMS) by Method 6020	WG939602	1	12/30/16 07:06	12/30/16 10:43	JPD
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938676	1	12/26/16 16:24	12/26/16 16:24	BMB
Wet Chemistry by Method 4500S2 D-2011	WG938033	1	12/22/16 22:14	12/22/16 22:14	MZ
Wet Chemistry by Method 9012B	WG939301	1	12/29/16 14:18	12/29/16 17:56	ASK
Wet Chemistry by Method 9056A	WG938354	1	12/27/16 20:45	12/27/16 20:45	KCF
Wet Chemistry by Method 9060A	WG939610	1	12/31/16 16:01	12/31/16 16:01	SJM

FIELD BLANK L880396-30 GW

Collected by: Darren Motley
Collected date/time: 12/21/16 09:55
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Gravimetric Analysis by Method 2540 C-2011	WG938690	1	12/28/16 13:53	12/28/16 14:22	JER
Mercury by Method 7470A	WG938138	1	12/23/16 09:22	12/23/16 18:04	TRB
Metals (ICPMS) by Method 6020	WG938049	1	12/27/16 13:37	12/30/16 01:00	VSS
Metals (ICPMS) by Method 6020	WG939602	1	12/30/16 07:06	12/30/16 10:46	JPD
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938676	1	12/26/16 16:37	12/26/16 16:37	BMB
Wet Chemistry by Method 4500S2 D-2011	WG938033	1	12/22/16 22:15	12/22/16 22:15	MZ
Wet Chemistry by Method 9012B	WG939301	1	12/29/16 14:18	12/29/16 17:58	ASK
Wet Chemistry by Method 9056A	WG938517	1	12/27/16 18:29	12/27/16 18:29	KCF
Wet Chemistry by Method 9060A	WG939610	1	12/31/16 16:20	12/31/16 16:20	SJM

EQUIPMENT BLANK L880396-31 GW

Collected by: Darren Motley
Collected date/time: 12/21/16 09:57
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938676	1	12/26/16 16:50	12/26/16 16:50	BMB

TRIP BLANK L880396-32 GW

Collected by: Darren Motley
Collected date/time: 12/21/16 10:00
Received date/time: 12/22/16 09:30

Method	Batch	Dilution	Preparation date/time	Analysis date/time	Analyst
Volatile Organic Compounds (GC/MS) by Method 8260B	WG938588	1	12/25/16 19:07	12/25/16 19:07	JHH

All sample aliquots were received at the correct temperature, in the proper containers, with the appropriate preservatives, and within method specified holding times. All MDL (LOD) and RDL (LOQ) values reported for environmental samples have been corrected for the dilution factor used in the analysis. All Method and Batch Quality Control are within established criteria except where addressed in this case narrative, a non-conformance form or properly qualified within the sample results. By my digital signature below, I affirm to the best of my knowledge, all problems/anomalies observed by the laboratory as having the potential to affect the quality of the data have been identified by the laboratory, and no information or data have been knowingly withheld that would affect the quality of the data.

Mark W. Beasley
Technical Service Representative

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	972000		2820	10000	1	12/27/2016 15:37	WG938474

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:24	WG938032

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:29	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	122000		260	5000	5	12/24/2016 20:16	WG938352
Sulfate	23700		77.4	5000	1	12/24/2016 16:17	WG938352

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	9230		102	1000	1	12/29/2016 12:12	WG938825

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 14:50	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	1.15	J	0.754	2.00	1	12/29/2016 23:04	WG938047
Arsenic	71.0		0.250	2.00	1	12/29/2016 23:04	WG938047
Barium	201		0.360	5.00	1	12/29/2016 23:04	WG938047
Beryllium	U		0.120	2.00	1	12/29/2016 23:04	WG938047
Cadmium	0.654	J	0.160	1.00	1	12/29/2016 23:04	WG938047
Chromium	5.88		0.540	2.00	1	12/29/2016 23:04	WG938047
Copper	10.4	B	0.520	5.00	1	12/29/2016 23:04	WG938047
Cobalt	109		0.260	2.00	1	12/29/2016 23:04	WG938047
Iron	6300		15.0	100	1	12/29/2016 23:04	WG938047
Lead	2.90	B	0.240	2.00	1	12/29/2016 23:04	WG938047
Manganese	780		0.250	5.00	1	12/29/2016 23:04	WG938047
Nickel	158		0.350	2.00	1	12/29/2016 23:04	WG938047
Selenium	U		0.380	2.00	1	12/29/2016 23:04	WG938047
Silver	U		0.310	2.00	1	12/29/2016 23:04	WG938047
Thallium	0.372	J	0.190	2.00	1	12/29/2016 23:04	WG938047
Tin	0.693	B J	0.300	2.00	1	12/29/2016 23:04	WG938047
Vanadium	1.78	J	0.180	5.00	1	12/29/2016 23:04	WG938047
Zinc	619		2.56	25.0	1	12/29/2016 23:04	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 04:17	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 04:17	WG938408
Benzene	0.572	J	0.331	1.00	1	12/26/2016 04:17	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 04:17	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 04:17	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 04:17	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 04:17	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 04:17	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 04:17	WG938408
Chlorobenzene	1.20		0.348	1.00	1	12/26/2016 04:17	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 04:17	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 04:17	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 04:17	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 04:17	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 04:17	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 04:17	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 04:17	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 04:17	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 04:17	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 04:17	WG938408
1,1-Dichloroethane	11.1		0.259	1.00	1	12/26/2016 04:17	WG938408
1,2-Dichloroethane	0.399	J	0.361	1.00	1	12/26/2016 04:17	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 04:17	WG938408
cis-1,2-Dichloroethene	7.76		0.260	1.00	1	12/26/2016 04:17	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 04:17	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 04:17	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 04:17	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 04:17	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 04:17	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 04:17	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 04:17	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 04:17	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 04:17	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 04:17	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 04:17	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 04:17	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 04:17	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 04:17	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 04:17	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 04:17	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 04:17	WG938408
Trichloroethene	0.607	J	0.398	1.00	1	12/26/2016 04:17	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 04:17	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 04:17	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 04:17	WG938408
Vinyl chloride	1.37		0.259	1.00	1	12/26/2016 04:17	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 04:17	WG938408
(S) Toluene-d8	*100*			*90.0-115*		*12/26/2016 04:17*	WG938408
(S) Dibromofluoromethane	*102*			*79.0-121*		*12/26/2016 04:17*	WG938408
(S) 4-Bromofluorobenzene	*94.4*			*80.1-120*		*12/26/2016 04:17*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	381000		2820	10000	1	12/27/2016 15:37	WG938474

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:24	WG938032

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:30	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	3610		51.9	1000	1	12/24/2016 16:32	WG938352
Sulfate	14700		77.4	5000	1	12/24/2016 16:32	WG938352

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	519	J	102	1000	1	12/28/2016 13:39	WG938824

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 14:57	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/29/2016 23:07	WG938047
Arsenic	4.66		0.250	2.00	1	12/29/2016 23:07	WG938047
Barium	31.5		0.360	5.00	1	12/29/2016 23:07	WG938047
Beryllium	U		0.120	2.00	1	12/29/2016 23:07	WG938047
Cadmium	U		0.160	1.00	1	12/29/2016 23:07	WG938047
Chromium	U		0.540	2.00	1	12/29/2016 23:07	WG938047
Copper	0.563	B J	0.520	5.00	1	12/29/2016 23:07	WG938047
Cobalt	0.295	J	0.260	2.00	1	12/29/2016 23:07	WG938047
Iron	118		15.0	100	1	12/29/2016 23:07	WG938047
Lead	0.636	B J	0.240	2.00	1	12/29/2016 23:07	WG938047
Manganese	10.4		0.250	5.00	1	12/29/2016 23:07	WG938047
Nickel	9.79		0.350	2.00	1	12/29/2016 23:07	WG938047
Selenium	U		0.380	2.00	1	12/29/2016 23:07	WG938047
Silver	U		0.310	2.00	1	12/29/2016 23:07	WG938047
Thallium	U		0.190	2.00	1	12/29/2016 23:07	WG938047
Tin	U		0.300	2.00	1	12/29/2016 23:07	WG938047
Vanadium	0.361	J	0.180	5.00	1	12/29/2016 23:07	WG938047
Zinc	3.66	B J	2.56	25.0	1	12/29/2016 23:07	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 04:41	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 04:41	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 04:41	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 04:41	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 04:41	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 04:41	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 04:41	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 04:41	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 04:41	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 04:41	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 04:41	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 04:41	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 04:41	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 04:41	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 04:41	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 04:41	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 04:41	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 04:41	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 04:41	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 04:41	WG938408
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 04:41	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 04:41	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 04:41	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 04:41	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 04:41	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 04:41	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 04:41	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 04:41	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 04:41	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 04:41	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 04:41	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 04:41	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 04:41	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 04:41	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 04:41	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 04:41	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 04:41	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 04:41	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 04:41	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 04:41	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 04:41	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 04:41	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 04:41	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 04:41	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 04:41	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 04:41	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 04:41	WG938408
(S) Toluene-d8	*101*			*90.0-115*		*12/26/2016 04:41*	WG938408
(S) Dibromofluoromethane	*103*			*79.0-121*		*12/26/2016 04:41*	WG938408
(S) 4-Bromofluorobenzene	*94.5*			*80.1-120*		*12/26/2016 04:41*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	525000		2820	10000	1	12/27/2016 14:20	WG938476

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:24	WG938032

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:31	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	9540		51.9	1000	1	12/24/2016 15:47	WG938352
Sulfate	12300		77.4	5000	1	12/24/2016 15:47	WG938352

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	2320		102	1000	1	12/28/2016 13:50	WG938824

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 14:59	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/29/2016 23:11	WG938047
Arsenic	2.31		0.250	2.00	1	12/29/2016 23:11	WG938047
Barium	81.2		0.360	5.00	1	12/29/2016 23:11	WG938047
Beryllium	U		0.120	2.00	1	12/29/2016 23:11	WG938047
Cadmium	0.430	J	0.160	1.00	1	12/29/2016 23:11	WG938047
Chromium	0.594	J	0.540	2.00	1	12/29/2016 23:11	WG938047
Copper	1.08	B J	0.520	5.00	1	12/29/2016 23:11	WG938047
Cobalt	4.10		0.260	2.00	1	12/29/2016 23:11	WG938047
Iron	473		15.0	100	1	12/29/2016 23:11	WG938047
Lead	0.937	B J	0.240	2.00	1	12/29/2016 23:11	WG938047
Manganese	2230		0.250	5.00	1	12/29/2016 23:11	WG938047
Nickel	62.2		0.350	2.00	1	12/29/2016 23:11	WG938047
Selenium	U		0.380	2.00	1	12/29/2016 23:11	WG938047
Silver	U		0.310	2.00	1	12/29/2016 23:11	WG938047
Thallium	0.661	J	0.190	2.00	1	12/29/2016 23:11	WG938047
Tin	U		0.300	2.00	1	12/29/2016 23:11	WG938047
Vanadium	0.426	J	0.180	5.00	1	12/29/2016 23:11	WG938047
Zinc	146		2.56	25.0	1	12/29/2016 23:11	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 05:04	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 05:04	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 05:04	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 05:04	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 05:04	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 05:04	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 05:04	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 05:04	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 05:04	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 05:04	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 05:04	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 05:04	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 05:04	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 05:04	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 05:04	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 05:04	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 05:04	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 05:04	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 05:04	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 05:04	WG938408
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 05:04	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 05:04	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 05:04	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 05:04	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 05:04	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 05:04	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 05:04	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 05:04	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 05:04	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 05:04	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 05:04	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 05:04	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 05:04	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 05:04	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 05:04	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 05:04	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 05:04	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 05:04	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 05:04	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 05:04	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 05:04	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 05:04	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 05:04	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 05:04	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 05:04	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 05:04	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 05:04	WG938408
(S) Toluene-d8	*101*			*90.0-115*		*12/26/2016 05:04*	WG938408
(S) Dibromofluoromethane	*101*			*79.0-121*		*12/26/2016 05:04*	WG938408
(S) 4-Bromofluorobenzene	*93.1*			*80.1-120*		*12/26/2016 05:04*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	376000		2820	10000	1	12/27/2016 14:20	WG938476

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:24	WG938032

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:32	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	7310		51.9	1000	1	12/24/2016 17:32	WG938352
Sulfate	14900		77.4	5000	1	12/24/2016 17:32	WG938352

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	1090		102	1000	1	12/28/2016 14:02	WG938824

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 15:02	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/29/2016 23:14	WG938047
Arsenic	0.505	J	0.250	2.00	1	12/29/2016 23:14	WG938047
Barium	35.8		0.360	5.00	1	12/29/2016 23:14	WG938047
Beryllium	U		0.120	2.00	1	12/29/2016 23:14	WG938047
Cadmium	0.883	J	0.160	1.00	1	12/29/2016 23:14	WG938047
Chromium	0.895	J	0.540	2.00	1	12/29/2016 23:14	WG938047
Copper	1.22	B J	0.520	5.00	1	12/29/2016 23:14	WG938047
Cobalt	U		0.260	2.00	1	12/29/2016 23:14	WG938047
Iron	41.1	J	15.0	100	1	12/29/2016 23:14	WG938047
Lead	0.888	B J	0.240	2.00	1	12/29/2016 23:14	WG938047
Manganese	2.09	B J	0.250	5.00	1	12/29/2016 23:14	WG938047
Nickel	1.14	J	0.350	2.00	1	12/29/2016 23:14	WG938047
Selenium	U		0.380	2.00	1	12/29/2016 23:14	WG938047
Silver	U		0.310	2.00	1	12/29/2016 23:14	WG938047
Thallium	U		0.190	2.00	1	12/29/2016 23:14	WG938047
Tin	U		0.300	2.00	1	12/29/2016 23:14	WG938047
Vanadium	0.400	J	0.180	5.00	1	12/29/2016 23:14	WG938047
Zinc	105		2.56	25.0	1	12/29/2016 23:14	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 05:27	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 05:27	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 05:27	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 05:27	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 05:27	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 05:27	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 05:27	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 05:27	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 05:27	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 05:27	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 05:27	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 05:27	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 05:27	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 05:27	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 05:27	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 05:27	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 05:27	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 05:27	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 05:27	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 05:27	WG938408
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 05:27	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 05:27	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 05:27	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 05:27	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 05:27	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 05:27	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 05:27	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 05:27	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 05:27	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 05:27	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 05:27	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 05:27	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 05:27	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 05:27	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 05:27	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 05:27	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 05:27	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 05:27	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 05:27	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 05:27	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 05:27	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 05:27	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 05:27	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 05:27	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 05:27	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 05:27	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 05:27	WG938408
(S) Toluene-d8	*101*			*90.0-115*		*12/26/2016 05:27*	WG938408
(S) Dibromofluoromethane	*104*			*79.0-121*		*12/26/2016 05:27*	WG938408
(S) 4-Bromofluorobenzene	*95.3*			*80.1-120*		*12/26/2016 05:27*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	416000		2820	10000	1	12/27/2016 14:20	WG938476

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U	J5	6.50	50.0	1	12/22/2016 22:24	WG938032

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:33	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	5760		51.9	1000	1	12/24/2016 17:47	WG938352
Sulfate	7070		77.4	5000	1	12/24/2016 17:47	WG938352

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	1200		102	1000	1	12/28/2016 14:13	WG938824

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 15:04	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/29/2016 23:18	WG938047
Arsenic	0.536	J	0.250	2.00	1	12/29/2016 23:18	WG938047
Barium	34.1		0.360	5.00	1	12/29/2016 23:18	WG938047
Beryllium	U		0.120	2.00	1	12/29/2016 23:18	WG938047
Cadmium	U		0.160	1.00	1	12/29/2016 23:18	WG938047
Chromium	3.28		0.540	2.00	1	12/29/2016 23:18	WG938047
Copper	3.77	B J	0.520	5.00	1	12/29/2016 23:18	WG938047
Cobalt	U		0.260	2.00	1	12/29/2016 23:18	WG938047
Iron	60.6	J	15.0	100	1	12/29/2016 23:18	WG938047
Lead	0.765	B J	0.240	2.00	1	12/29/2016 23:18	WG938047
Manganese	1.85	B J	0.250	5.00	1	12/29/2016 23:18	WG938047
Nickel	2.81		0.350	2.00	1	12/29/2016 23:18	WG938047
Selenium	U		0.380	2.00	1	12/29/2016 23:18	WG938047
Silver	U		0.310	2.00	1	12/29/2016 23:18	WG938047
Thallium	U		0.190	2.00	1	12/29/2016 23:18	WG938047
Tin	0.313	B J	0.300	2.00	1	12/29/2016 23:18	WG938047
Vanadium	0.385	J	0.180	5.00	1	12/29/2016 23:18	WG938047
Zinc	26.8	B	2.56	25.0	1	12/29/2016 23:18	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 05:51	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 05:51	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 05:51	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 05:51	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 05:51	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 05:51	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 05:51	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 05:51	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 05:51	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 05:51	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 05:51	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 05:51	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 05:51	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 05:51	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 05:51	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 05:51	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 05:51	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 05:51	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 05:51	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 05:51	WG938408
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 05:51	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 05:51	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 05:51	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 05:51	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 05:51	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 05:51	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 05:51	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 05:51	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 05:51	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 05:51	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 05:51	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 05:51	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 05:51	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 05:51	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 05:51	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 05:51	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 05:51	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 05:51	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 05:51	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 05:51	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 05:51	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 05:51	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 05:51	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 05:51	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 05:51	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 05:51	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 05:51	WG938408
(S) Toluene-d8	*101*			*90.0-115*		*12/26/2016 05:51*	WG938408
(S) Dibromofluoromethane	*104*			*79.0-121*		*12/26/2016 05:51*	WG938408
(S) 4-Bromofluorobenzene	*93.8*			*80.1-120*		*12/26/2016 05:51*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	432000		2820	10000	1	12/27/2016 14:20	WG938476

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:25	WG938032

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:36	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	21000		51.9	1000	1	12/24/2016 18:31	WG938352
Sulfate	7760		77.4	5000	1	12/24/2016 18:31	WG938352

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	1750		102	1000	1	12/28/2016 14:24	WG938824

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 15:06	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/29/2016 23:22	WG938047
Arsenic	0.423	J	0.250	2.00	1	12/29/2016 23:22	WG938047
Barium	46.5		0.360	5.00	1	12/29/2016 23:22	WG938047
Beryllium	U		0.120	2.00	1	12/29/2016 23:22	WG938047
Cadmium	0.439	J	0.160	1.00	1	12/29/2016 23:22	WG938047
Chromium	0.805	J	0.540	2.00	1	12/29/2016 23:22	WG938047
Copper	1.21	B J	0.520	5.00	1	12/29/2016 23:22	WG938047
Cobalt	U		0.260	2.00	1	12/29/2016 23:22	WG938047
Iron	33.4	J	15.0	100	1	12/29/2016 23:22	WG938047
Lead	0.601	B J	0.240	2.00	1	12/29/2016 23:22	WG938047
Manganese	0.867	B J	0.250	5.00	1	12/29/2016 23:22	WG938047
Nickel	0.957	J	0.350	2.00	1	12/29/2016 23:22	WG938047
Selenium	U		0.380	2.00	1	12/29/2016 23:22	WG938047
Silver	U		0.310	2.00	1	12/29/2016 23:22	WG938047
Thallium	U		0.190	2.00	1	12/29/2016 23:22	WG938047
Tin	0.311	B J	0.300	2.00	1	12/29/2016 23:22	WG938047
Vanadium	0.427	J	0.180	5.00	1	12/29/2016 23:22	WG938047
Zinc	41.5	B	2.56	25.0	1	12/29/2016 23:22	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 06:14	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 06:14	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 06:14	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 06:14	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 06:14	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 06:14	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 06:14	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 06:14	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 06:14	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 06:14	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 06:14	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 06:14	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 06:14	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 06:14	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 06:14	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 06:14	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 06:14	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 06:14	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 06:14	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 06:14	WG938408
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 06:14	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 06:14	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 06:14	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 06:14	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 06:14	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 06:14	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 06:14	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 06:14	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 06:14	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 06:14	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 06:14	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 06:14	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 06:14	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 06:14	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 06:14	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 06:14	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 06:14	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 06:14	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 06:14	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 06:14	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 06:14	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 06:14	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 06:14	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 06:14	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 06:14	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 06:14	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 06:14	WG938408
(S) Toluene-d8	*100*			*90.0-115*		*12/26/2016 06:14*	WG938408
(S) Dibromofluoromethane	*103*			*79.0-121*		*12/26/2016 06:14*	WG938408
(S) 4-Bromofluorobenzene	*93.6*			*80.1-120*		*12/26/2016 06:14*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	338000		2820	10000	1	12/27/2016 15:37	WG938474

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:26	WG938032

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:37	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	1500		51.9	1000	1	12/24/2016 18:46	WG938352
Sulfate	6390		77.4	5000	1	12/24/2016 18:46	WG938352

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	1070		102	1000	1	12/28/2016 14:35	WG938824

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 15:15	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/29/2016 23:25	WG938047
Arsenic	0.687	J	0.250	2.00	1	12/29/2016 23:25	WG938047
Barium	46.3		0.360	5.00	1	12/29/2016 23:25	WG938047
Beryllium	U		0.120	2.00	1	12/29/2016 23:25	WG938047
Cadmium	0.307	J	0.160	1.00	1	12/29/2016 23:25	WG938047
Chromium	U		0.540	2.00	1	12/29/2016 23:25	WG938047
Copper	1.70	B J	0.520	5.00	1	12/29/2016 23:25	WG938047
Cobalt	3.56		0.260	2.00	1	12/29/2016 23:25	WG938047
Iron	153		15.0	100	1	12/29/2016 23:25	WG938047
Lead	1.53	B J	0.240	2.00	1	12/29/2016 23:25	WG938047
Manganese	363		0.250	5.00	1	12/29/2016 23:25	WG938047
Nickel	10.6		0.350	2.00	1	12/29/2016 23:25	WG938047
Selenium	U		0.380	2.00	1	12/29/2016 23:25	WG938047
Silver	U		0.310	2.00	1	12/29/2016 23:25	WG938047
Thallium	U		0.190	2.00	1	12/29/2016 23:25	WG938047
Tin	0.318	B J	0.300	2.00	1	12/29/2016 23:25	WG938047
Vanadium	0.497	J	0.180	5.00	1	12/29/2016 23:25	WG938047
Zinc	15.1	B J	2.56	25.0	1	12/29/2016 23:25	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 06:37	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 06:37	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 06:37	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 06:37	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 06:37	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 06:37	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 06:37	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 06:37	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 06:37	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 06:37	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 06:37	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 06:37	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 06:37	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 06:37	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 06:37	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 06:37	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 06:37	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 06:37	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 06:37	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 06:37	WG938408
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 06:37	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 06:37	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 06:37	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 06:37	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 06:37	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 06:37	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 06:37	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 06:37	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 06:37	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 06:37	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 06:37	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 06:37	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 06:37	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 06:37	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 06:37	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 06:37	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 06:37	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 06:37	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 06:37	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 06:37	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 06:37	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 06:37	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 06:37	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 06:37	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 06:37	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 06:37	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 06:37	WG938408
(S) Toluene-d8	*101*			*90.0-115*		*12/26/2016 06:37*	WG938408
(S) Dibromofluoromethane	*102*			*79.0-121*		*12/26/2016 06:37*	WG938408
(S) 4-Bromofluorobenzene	*92.2*			*80.1-120*		*12/26/2016 06:37*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	794000		2820	10000	1	12/27/2016 14:20	WG938476

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:26	WG938032

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:38	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	95900		51.9	1000	1	12/24/2016 19:01	WG938352
Sulfate	104	J	77.4	5000	1	12/24/2016 19:01	WG938352

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	10600		204	2000	2	12/29/2016 12:37	WG938825

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 15:18	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/29/2016 23:29	WG938047
Arsenic	123		0.250	2.00	1	12/29/2016 23:29	WG938047
Barium	250		0.360	5.00	1	12/29/2016 23:29	WG938047
Beryllium	U		0.120	2.00	1	12/29/2016 23:29	WG938047
Cadmium	U		0.160	1.00	1	12/29/2016 23:29	WG938047
Chromium	1.32	J	0.540	2.00	1	12/29/2016 23:29	WG938047
Copper	0.868	B J	0.520	5.00	1	12/29/2016 23:29	WG938047
Cobalt	28.6		0.260	2.00	1	12/29/2016 23:29	WG938047
Iron	36300		15.0	100	1	12/29/2016 23:29	WG938047
Lead	0.886	B J	0.240	2.00	1	12/29/2016 23:29	WG938047
Manganese	1200		0.250	5.00	1	12/29/2016 23:29	WG938047
Nickel	57.1		0.350	2.00	1	12/29/2016 23:29	WG938047
Selenium	U		0.380	2.00	1	12/29/2016 23:29	WG938047
Silver	U		0.310	2.00	1	12/29/2016 23:29	WG938047
Thallium	U		0.190	2.00	1	12/29/2016 23:29	WG938047
Tin	0.812	B J	0.300	2.00	1	12/29/2016 23:29	WG938047
Vanadium	1.15	J	0.180	5.00	1	12/29/2016 23:29	WG938047
Zinc	27.2	B	2.56	25.0	1	12/29/2016 23:29	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 07:00	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 07:00	WG938408
Benzene	4.34		0.331	1.00	1	12/26/2016 07:00	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 07:00	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 07:00	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 07:00	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 07:00	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 07:00	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 07:00	WG938408
Chlorobenzene	0.849	J	0.348	1.00	1	12/26/2016 07:00	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 07:00	WG938408
Chloroethane	2.00	J J4	0.453	5.00	1	12/26/2016 07:00	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 07:00	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 07:00	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 07:00	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 07:00	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 07:00	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 07:00	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 07:00	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 07:00	WG938408
1,1-Dichloroethane	5.43		0.259	1.00	1	12/26/2016 07:00	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 07:00	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 07:00	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 07:00	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 07:00	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 07:00	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 07:00	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 07:00	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 07:00	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 07:00	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 07:00	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 07:00	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 07:00	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 07:00	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 07:00	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 07:00	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 07:00	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 07:00	WG938408
Toluene	0.443	J	0.412	1.00	1	12/26/2016 07:00	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 07:00	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 07:00	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 07:00	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 07:00	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 07:00	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 07:00	WG938408
Vinyl chloride	2.60		0.259	1.00	1	12/26/2016 07:00	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 07:00	WG938408
(S) Toluene-d8	*101*			*90.0-115*		*12/26/2016 07:00*	WG938408
(S) Dibromofluoromethane	*103*			*79.0-121*		*12/26/2016 07:00*	WG938408
(S) 4-Bromofluorobenzene	*93.6*			*80.1-120*		*12/26/2016 07:00*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	433000		2820	10000	1	12/28/2016 14:22	WG938690

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:26	WG938032

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:39	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	66200		51.9	1000	1	12/24/2016 19:16	WG938352
Sulfate	261	J	77.4	5000	1	12/24/2016 19:16	WG938352

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	5770		102	1000	1	12/28/2016 14:58	WG938824

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 15:20	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/29/2016 23:32	WG938047
Arsenic	1.35	J	0.250	2.00	1	12/29/2016 23:32	WG938047
Barium	99.7		0.360	5.00	1	12/29/2016 23:32	WG938047
Beryllium	U		0.120	2.00	1	12/29/2016 23:32	WG938047
Cadmium	1.48		0.160	1.00	1	12/29/2016 23:32	WG938047
Chromium	0.821	J	0.540	2.00	1	12/29/2016 23:32	WG938047
Copper	1.50	B J	0.520	5.00	1	12/29/2016 23:32	WG938047
Cobalt	11.1		0.260	2.00	1	12/29/2016 23:32	WG938047
Iron	305		15.0	100	1	12/29/2016 23:32	WG938047
Lead	1.30	B J	0.240	2.00	1	12/29/2016 23:32	WG938047
Manganese	4350		0.250	5.00	1	12/29/2016 23:32	WG938047
Nickel	43.8		0.350	2.00	1	12/29/2016 23:32	WG938047
Selenium	U		0.380	2.00	1	12/29/2016 23:32	WG938047
Silver	U		0.310	2.00	1	12/29/2016 23:32	WG938047
Thallium	0.612	J	0.190	2.00	1	12/29/2016 23:32	WG938047
Tin	0.432	B J	0.300	2.00	1	12/29/2016 23:32	WG938047
Vanadium	0.711	J	0.180	5.00	1	12/29/2016 23:32	WG938047
Zinc	42.9	B	2.56	25.0	1	12/29/2016 23:32	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 07:24	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 07:24	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 07:24	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 07:24	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 07:24	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 07:24	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 07:24	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 07:24	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 07:24	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 07:24	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 07:24	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 07:24	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 07:24	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 07:24	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 07:24	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 07:24	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 07:24	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 07:24	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 07:24	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 07:24	WG938408
1,1-Dichloroethane	0.488	J	0.259	1.00	1	12/26/2016 07:24	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 07:24	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 07:24	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 07:24	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 07:24	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 07:24	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 07:24	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 07:24	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 07:24	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 07:24	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 07:24	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 07:24	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 07:24	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 07:24	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 07:24	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 07:24	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 07:24	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 07:24	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 07:24	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 07:24	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 07:24	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 07:24	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 07:24	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 07:24	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 07:24	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 07:24	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 07:24	WG938408
(S) Toluene-d8	*99.7*			*90.0-115*		*12/26/2016 07:24*	WG938408
(S) Dibromofluoromethane	*102*			*79.0-121*		*12/26/2016 07:24*	WG938408
(S) 4-Bromofluorobenzene	*94.0*			*80.1-120*		*12/26/2016 07:24*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	305000		2820	10000	1	12/28/2016 14:22	WG938690

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:26	WG938032

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:41	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	17900		51.9	1000	1	12/27/2016 16:01	WG938354
Sulfate	6010		77.4	5000	1	12/27/2016 16:01	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	1640		102	1000	1	12/28/2016 15:10	WG938824

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 15:22	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/29/2016 23:36	WG938047
Arsenic	U		0.250	2.00	1	12/29/2016 23:36	WG938047
Barium	41.2		0.360	5.00	1	12/29/2016 23:36	WG938047
Beryllium	U		0.120	2.00	1	12/29/2016 23:36	WG938047
Cadmium	0.765	J	0.160	1.00	1	12/29/2016 23:36	WG938047
Chromium	U		0.540	2.00	1	12/29/2016 23:36	WG938047
Copper	1.05	B J	0.520	5.00	1	12/29/2016 23:36	WG938047
Cobalt	U		0.260	2.00	1	12/29/2016 23:36	WG938047
Iron	39.0	J	15.0	100	1	12/29/2016 23:36	WG938047
Lead	1.06	B J	0.240	2.00	1	12/29/2016 23:36	WG938047
Manganese	8.40		0.250	5.00	1	12/29/2016 23:36	WG938047
Nickel	1.96	J	0.350	2.00	1	12/29/2016 23:36	WG938047
Selenium	U		0.380	2.00	1	12/29/2016 23:36	WG938047
Silver	U		0.310	2.00	1	12/29/2016 23:36	WG938047
Thallium	U		0.190	2.00	1	12/29/2016 23:36	WG938047
Tin	0.506	B J	0.300	2.00	1	12/29/2016 23:36	WG938047
Vanadium	0.514	J	0.180	5.00	1	12/29/2016 23:36	WG938047
Zinc	21.3	B J	2.56	25.0	1	12/29/2016 23:36	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 07:47	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 07:47	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 07:47	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 07:47	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 07:47	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 07:47	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 07:47	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 07:47	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 07:47	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 07:47	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 07:47	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 07:47	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 07:47	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 07:47	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 07:47	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 07:47	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 07:47	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 07:47	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 07:47	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 07:47	WG938408
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 07:47	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 07:47	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 07:47	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 07:47	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 07:47	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 07:47	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 07:47	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 07:47	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 07:47	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 07:47	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 07:47	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 07:47	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 07:47	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 07:47	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 07:47	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 07:47	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 07:47	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 07:47	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 07:47	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 07:47	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 07:47	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 07:47	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 07:47	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 07:47	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 07:47	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 07:47	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 07:47	WG938408
(S) Toluene-d8	*100*			*90.0-115*		*12/26/2016 07:47*	WG938408
(S) Dibromofluoromethane	*104*			*79.0-121*		*12/26/2016 07:47*	WG938408
(S) 4-Bromofluorobenzene	*92.7*			*80.1-120*		*12/26/2016 07:47*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	443000		2820	10000	1	12/27/2016 15:37	WG938474

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:26	WG938032

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:42	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	10400		51.9	1000	1	12/27/2016 16:21	WG938354
Sulfate	13500		77.4	5000	1	12/27/2016 16:21	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	1390		102	1000	1	12/28/2016 16:16	WG938824

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 15:24	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	0.806	J	0.754	2.00	1	12/29/2016 23:50	WG938047
Arsenic	6.28		0.250	2.00	1	12/29/2016 23:50	WG938047
Barium	39.4		0.360	5.00	1	12/29/2016 23:50	WG938047
Beryllium	U		0.120	2.00	1	12/29/2016 23:50	WG938047
Cadmium	U		0.160	1.00	1	12/29/2016 23:50	WG938047
Chromium	5.77		0.540	2.00	1	12/29/2016 23:50	WG938047
Copper	1.85	B J	0.520	5.00	1	12/29/2016 23:50	WG938047
Cobalt	U		0.260	2.00	1	12/29/2016 23:50	WG938047
Iron	536		15.0	100	1	12/29/2016 23:50	WG938047
Lead	3.16	B	0.240	2.00	1	12/29/2016 23:50	WG938047
Manganese	17.5		0.250	5.00	1	12/29/2016 23:50	WG938047
Nickel	5.86		0.350	2.00	1	12/29/2016 23:50	WG938047
Selenium	U		0.380	2.00	1	12/29/2016 23:50	WG938047
Silver	U		0.310	2.00	1	12/29/2016 23:50	WG938047
Thallium	U		0.190	2.00	1	12/29/2016 23:50	WG938047
Tin	0.638	B J	0.300	2.00	1	12/29/2016 23:50	WG938047
Vanadium	1.28	J	0.180	5.00	1	12/29/2016 23:50	WG938047
Zinc	60.2		2.56	25.0	1	12/29/2016 23:50	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 08:10	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 08:10	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 08:10	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 08:10	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 08:10	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 08:10	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 08:10	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 08:10	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 08:10	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 08:10	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 08:10	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 08:10	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 08:10	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 08:10	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 08:10	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 08:10	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 08:10	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 08:10	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 08:10	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 08:10	WG938408
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 08:10	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 08:10	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 08:10	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 08:10	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 08:10	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 08:10	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 08:10	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 08:10	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 08:10	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 08:10	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 08:10	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 08:10	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 08:10	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 08:10	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 08:10	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 08:10	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 08:10	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 08:10	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 08:10	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 08:10	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 08:10	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 08:10	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 08:10	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 08:10	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 08:10	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 08:10	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 08:10	WG938408
(S) Toluene-d8	*101*			*90.0-115*		*12/26/2016 08:10*	WG938408
(S) Dibromofluoromethane	*102*			*79.0-121*		*12/26/2016 08:10*	WG938408
(S) 4-Bromofluorobenzene	*92.3*			*80.1-120*		*12/26/2016 08:10*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	425000		2820	10000	1	12/27/2016 15:37	WG938474

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:26	WG938032

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:43	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	9170		51.9	1000	1	12/27/2016 16:31	WG938354
Sulfate	8970		77.4	5000	1	12/27/2016 16:31	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	94700		102	1000	1	12/28/2016 16:41	WG938824

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	0.134	J	0.0490	0.200	1	12/27/2016 14:46	WG938411

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/29/2016 23:54	WG938047
Arsenic	0.561	J	0.250	2.00	1	12/29/2016 23:54	WG938047
Barium	50.2		0.360	5.00	1	12/29/2016 23:54	WG938047
Beryllium	U		0.120	2.00	1	12/29/2016 23:54	WG938047
Cadmium	0.995	J	0.160	1.00	1	12/29/2016 23:54	WG938047
Chromium	1.21	J	0.540	2.00	1	12/29/2016 23:54	WG938047
Copper	2.63	B J	0.520	5.00	1	12/29/2016 23:54	WG938047
Cobalt	1.25	J	0.260	2.00	1	12/29/2016 23:54	WG938047
Iron	108		15.0	100	1	12/29/2016 23:54	WG938047
Lead	0.842	B J	0.240	2.00	1	12/29/2016 23:54	WG938047
Manganese	48.3		0.250	5.00	1	12/29/2016 23:54	WG938047
Nickel	2.68		0.350	2.00	1	12/29/2016 23:54	WG938047
Selenium	U		0.380	2.00	1	12/29/2016 23:54	WG938047
Silver	U		0.310	2.00	1	12/29/2016 23:54	WG938047
Thallium	U		0.190	2.00	1	12/29/2016 23:54	WG938047
Tin	0.407	B J	0.300	2.00	1	12/29/2016 23:54	WG938047
Vanadium	0.542	J	0.180	5.00	1	12/29/2016 23:54	WG938047
Zinc	82.3		2.56	25.0	1	12/29/2016 23:54	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 08:33	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 08:33	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 08:33	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 08:33	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 08:33	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 08:33	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 08:33	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 08:33	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 08:33	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 08:33	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 08:33	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 08:33	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 08:33	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 08:33	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 08:33	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 08:33	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 08:33	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 08:33	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 08:33	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 08:33	WG938408
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 08:33	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 08:33	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 08:33	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 08:33	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 08:33	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 08:33	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 08:33	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 08:33	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 08:33	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 08:33	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 08:33	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 08:33	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 08:33	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 08:33	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 08:33	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 08:33	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 08:33	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 08:33	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 08:33	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 08:33	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 08:33	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 08:33	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 08:33	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 08:33	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 08:33	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 08:33	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 08:33	WG938408
(S) Toluene-d8	*100*			*90.0-115*		*12/26/2016 08:33*	WG938408
(S) Dibromofluoromethane	*102*			*79.0-121*		*12/26/2016 08:33*	WG938408
(S) 4-Bromofluorobenzene	*92.4*			*80.1-120*		*12/26/2016 08:33*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	458000		2820	10000	1	12/27/2016 15:37	WG938474

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:27	WG938032

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:44	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	4110		51.9	1000	1	12/27/2016 16:41	WG938354
Sulfate	41900		77.4	5000	1	12/27/2016 16:41	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	1790		102	1000	1	12/28/2016 16:55	WG938824

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 15:27	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/29/2016 23:57	WG938047
Arsenic	4.58		0.250	2.00	1	12/29/2016 23:57	WG938047
Barium	41.9		0.360	5.00	1	12/29/2016 23:57	WG938047
Beryllium	U		0.120	2.00	1	12/29/2016 23:57	WG938047
Cadmium	1.28		0.160	1.00	1	12/29/2016 23:57	WG938047
Chromium	5.71		0.540	2.00	1	12/29/2016 23:57	WG938047
Copper	4.00	B J	0.520	5.00	1	12/29/2016 23:57	WG938047
Cobalt	1.06	J	0.260	2.00	1	12/29/2016 23:57	WG938047
Iron	247		15.0	100	1	12/29/2016 23:57	WG938047
Lead	1.48	B J	0.240	2.00	1	12/29/2016 23:57	WG938047
Manganese	16.6		0.250	5.00	1	12/29/2016 23:57	WG938047
Nickel	18.7		0.350	2.00	1	12/29/2016 23:57	WG938047
Selenium	U		0.380	2.00	1	12/29/2016 23:57	WG938047
Silver	U		0.310	2.00	1	12/29/2016 23:57	WG938047
Thallium	0.547	J	0.190	2.00	1	12/29/2016 23:57	WG938047
Tin	0.531	B J	0.300	2.00	1	12/29/2016 23:57	WG938047
Vanadium	0.569	J	0.180	5.00	1	12/29/2016 23:57	WG938047
Zinc	656		2.56	25.0	1	12/29/2016 23:57	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 08:57	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 08:57	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 08:57	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 08:57	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 08:57	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 08:57	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 08:57	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 08:57	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 08:57	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 08:57	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 08:57	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 08:57	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 08:57	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 08:57	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 08:57	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 08:57	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 08:57	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 08:57	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 08:57	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 08:57	WG938408
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 08:57	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 08:57	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 08:57	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 08:57	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 08:57	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 08:57	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 08:57	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 08:57	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 08:57	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 08:57	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 08:57	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 08:57	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 08:57	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 08:57	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 08:57	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 08:57	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 08:57	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 08:57	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 08:57	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 08:57	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 08:57	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 08:57	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 08:57	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 08:57	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 08:57	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 08:57	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 08:57	WG938408
(S) Toluene-d8	*101*			*90.0-115*		*12/26/2016 08:57*	WG938408
(S) Dibromofluoromethane	*102*			*79.0-121*		*12/26/2016 08:57*	WG938408
(S) 4-Bromofluorobenzene	*93.1*			*80.1-120*		*12/26/2016 08:57*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	399000		2820	10000	1	12/27/2016 15:37	WG938474

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:27	WG938032

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:47	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	3050		51.9	1000	1	12/27/2016 16:52	WG938354
Sulfate	29700		77.4	5000	1	12/27/2016 16:52	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	1610	P1	102	1000	1	12/28/2016 17:06	WG938824

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 15:29	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/30/2016 00:01	WG938047
Arsenic	5.74		0.250	2.00	1	12/30/2016 00:01	WG938047
Barium	23.0		0.360	5.00	1	12/30/2016 00:01	WG938047
Beryllium	U		0.120	2.00	1	12/30/2016 00:01	WG938047
Cadmium	U		0.160	1.00	1	12/30/2016 00:01	WG938047
Chromium	7.44		0.540	2.00	1	12/30/2016 00:01	WG938047
Copper	5.29	B	0.520	5.00	1	12/30/2016 00:01	WG938047
Cobalt	0.774	J	0.260	2.00	1	12/30/2016 00:01	WG938047
Iron	400		15.0	100	1	12/30/2016 00:01	WG938047
Lead	5.55	B	0.240	2.00	1	12/30/2016 00:01	WG938047
Manganese	7.01		0.250	5.00	1	12/30/2016 00:01	WG938047
Nickel	19.5		0.350	2.00	1	12/30/2016 00:01	WG938047
Selenium	U		0.380	2.00	1	12/30/2016 00:01	WG938047
Silver	U		0.310	2.00	1	12/30/2016 00:01	WG938047
Thallium	U		0.190	2.00	1	12/30/2016 00:01	WG938047
Tin	U		0.300	2.00	1	12/30/2016 00:01	WG938047
Vanadium	0.359	J	0.180	5.00	1	12/30/2016 00:01	WG938047
Zinc	43.0	B	2.56	25.0	1	12/30/2016 00:01	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 09:20	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 09:20	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 09:20	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 09:20	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 09:20	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 09:20	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 09:20	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 09:20	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 09:20	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 09:20	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 09:20	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 09:20	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 09:20	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 09:20	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 09:20	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 09:20	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 09:20	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 09:20	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 09:20	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 09:20	WG938408
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 09:20	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 09:20	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 09:20	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 09:20	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 09:20	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 09:20	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 09:20	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 09:20	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 09:20	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 09:20	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 09:20	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 09:20	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 09:20	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 09:20	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 09:20	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 09:20	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 09:20	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 09:20	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 09:20	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 09:20	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 09:20	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 09:20	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 09:20	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 09:20	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 09:20	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 09:20	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 09:20	WG938408
(S) Toluene-d8	*103*			*90.0-115*		*12/26/2016 09:20*	WG938408
(S) Dibromofluoromethane	*102*			*79.0-121*		*12/26/2016 09:20*	WG938408
(S) 4-Bromofluorobenzene	*93.5*			*80.1-120*		*12/26/2016 09:20*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	371000		2820	10000	1	12/27/2016 15:37	WG938474

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:09	WG938033

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:48	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	2230		51.9	1000	1	12/27/2016 17:02	WG938354
Sulfate	11200		77.4	5000	1	12/27/2016 17:02	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	23300		102	1000	1	12/28/2016 17:32	WG938824

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 15:31	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/30/2016 00:04	WG938047
Arsenic	3.54		0.250	2.00	1	12/30/2016 00:04	WG938047
Barium	31.9		0.360	5.00	1	12/30/2016 00:04	WG938047
Beryllium	U		0.120	2.00	1	12/30/2016 00:04	WG938047
Cadmium	0.291	J	0.160	1.00	1	12/30/2016 00:04	WG938047
Chromium	U		0.540	2.00	1	12/30/2016 00:04	WG938047
Copper	0.959	B J	0.520	5.00	1	12/30/2016 00:04	WG938047
Cobalt	U		0.260	2.00	1	12/30/2016 00:04	WG938047
Iron	37.8	J	15.0	100	1	12/30/2016 00:04	WG938047
Lead	0.583	B J	0.240	2.00	1	12/30/2016 00:04	WG938047
Manganese	4.05	J	0.250	5.00	1	12/30/2016 00:04	WG938047
Nickel	U		0.350	2.00	1	12/30/2016 00:04	WG938047
Selenium	U		0.380	2.00	1	12/30/2016 00:04	WG938047
Silver	U		0.310	2.00	1	12/30/2016 00:04	WG938047
Thallium	U		0.190	2.00	1	12/30/2016 00:04	WG938047
Tin	U		0.300	2.00	1	12/30/2016 00:04	WG938047
Vanadium	0.351	J	0.180	5.00	1	12/30/2016 00:04	WG938047
Zinc	7.99	B J	2.56	25.0	1	12/30/2016 00:04	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 09:43	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 09:43	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 09:43	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 09:43	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 09:43	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 09:43	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 09:43	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 09:43	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 09:43	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 09:43	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 09:43	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 09:43	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 09:43	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 09:43	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 09:43	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 09:43	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 09:43	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 09:43	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 09:43	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 09:43	WG938408
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 09:43	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 09:43	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 09:43	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 09:43	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 09:43	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 09:43	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 09:43	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 09:43	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 09:43	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 09:43	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 09:43	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 09:43	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 09:43	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 09:43	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 09:43	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 09:43	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 09:43	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 09:43	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 09:43	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 09:43	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 09:43	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 09:43	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 09:43	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 09:43	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 09:43	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 09:43	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 09:43	WG938408
(S) Toluene-d8	*101*			*90.0-115*		*12/26/2016 09:43*	WG938408
(S) Dibromofluoromethane	*105*			*79.0-121*		*12/26/2016 09:43*	WG938408
(S) 4-Bromofluorobenzene	*95.8*			*80.1-120*		*12/26/2016 09:43*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	395000		2820	10000	1	12/27/2016 15:37	WG938474

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:10	WG938033

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:51	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	6420		51.9	1000	1	12/27/2016 17:42	WG938354
Sulfate	16000		77.4	5000	1	12/27/2016 17:42	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	1230		102	1000	1	12/28/2016 17:43	WG938824

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 15:33	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/30/2016 00:08	WG938047
Arsenic	1.36	J	0.250	2.00	1	12/30/2016 00:08	WG938047
Barium	34.0		0.360	5.00	1	12/30/2016 00:08	WG938047
Beryllium	U		0.120	2.00	1	12/30/2016 00:08	WG938047
Cadmium	6.21		0.160	1.00	1	12/30/2016 00:08	WG938047
Chromium	2.33		0.540	2.00	1	12/30/2016 00:08	WG938047
Copper	2.62	B J	0.520	5.00	1	12/30/2016 00:08	WG938047
Cobalt	0.342	J	0.260	2.00	1	12/30/2016 00:08	WG938047
Iron	882		15.0	100	1	12/30/2016 00:08	WG938047
Lead	2.90	B	0.240	2.00	1	12/30/2016 00:08	WG938047
Manganese	14.9		0.250	5.00	1	12/30/2016 00:08	WG938047
Nickel	2.83		0.350	2.00	1	12/30/2016 00:08	WG938047
Selenium	U		0.380	2.00	1	12/30/2016 00:08	WG938047
Silver	U		0.310	2.00	1	12/30/2016 00:08	WG938047
Thallium	U		0.190	2.00	1	12/30/2016 00:08	WG938047
Tin	0.307	B J	0.300	2.00	1	12/30/2016 00:08	WG938047
Vanadium	2.56	J	0.180	5.00	1	12/30/2016 00:08	WG938047
Zinc	457		2.56	25.0	1	12/30/2016 00:08	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 10:07	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 10:07	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 10:07	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 10:07	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 10:07	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 10:07	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 10:07	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 10:07	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 10:07	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 10:07	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 10:07	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 10:07	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 10:07	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 10:07	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 10:07	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 10:07	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 10:07	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 10:07	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 10:07	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 10:07	WG938408
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 10:07	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 10:07	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 10:07	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 10:07	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 10:07	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 10:07	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 10:07	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 10:07	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 10:07	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 10:07	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 10:07	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 10:07	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 10:07	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 10:07	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 10:07	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 10:07	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 10:07	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 10:07	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 10:07	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 10:07	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 10:07	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 10:07	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 10:07	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 10:07	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 10:07	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 10:07	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 10:07	WG938408
(S) Toluene-d8	*103*			*90.0-115*		*12/26/2016 10:07*	WG938408
(S) Dibromofluoromethane	*104*			*79.0-121*		*12/26/2016 10:07*	WG938408
(S) 4-Bromofluorobenzene	*97.3*			*80.1-120*		*12/26/2016 10:07*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	398000		2820	10000	1	12/27/2016 15:37	WG938474

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:10	WG938033

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:52	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	2480		51.9	1000	1	12/27/2016 17:53	WG938354
Sulfate	20200		77.4	5000	1	12/27/2016 17:53	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	1140		102	1000	1	12/28/2016 17:55	WG938824

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 15:36	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/30/2016 00:12	WG938047
Arsenic	1.39	J	0.250	2.00	1	12/30/2016 00:12	WG938047
Barium	23.3		0.360	5.00	1	12/30/2016 00:12	WG938047
Beryllium	U		0.120	2.00	1	12/30/2016 00:12	WG938047
Cadmium	U		0.160	1.00	1	12/30/2016 00:12	WG938047
Chromium	U		0.540	2.00	1	12/30/2016 00:12	WG938047
Copper	0.986	B J	0.520	5.00	1	12/30/2016 00:12	WG938047
Cobalt	U		0.260	2.00	1	12/30/2016 00:12	WG938047
Iron	189		15.0	100	1	12/30/2016 00:12	WG938047
Lead	0.479	B J	0.240	2.00	1	12/30/2016 00:12	WG938047
Manganese	4.84	J	0.250	5.00	1	12/30/2016 00:12	WG938047
Nickel	4.80		0.350	2.00	1	12/30/2016 00:12	WG938047
Selenium	U		0.380	2.00	1	12/30/2016 00:12	WG938047
Silver	U		0.310	2.00	1	12/30/2016 00:12	WG938047
Thallium	U		0.190	2.00	1	12/30/2016 00:12	WG938047
Tin	U		0.300	2.00	1	12/30/2016 00:12	WG938047
Vanadium	0.297	J	0.180	5.00	1	12/30/2016 00:12	WG938047
Zinc	U		2.56	25.0	1	12/30/2016 00:12	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 10:30	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 10:30	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 10:30	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 10:30	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 10:30	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 10:30	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 10:30	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 10:30	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 10:30	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 10:30	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 10:30	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 10:30	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 10:30	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 10:30	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 10:30	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 10:30	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 10:30	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 10:30	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 10:30	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 10:30	WG938408
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 10:30	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 10:30	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 10:30	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 10:30	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 10:30	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 10:30	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 10:30	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 10:30	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 10:30	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 10:30	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 10:30	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 10:30	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 10:30	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 10:30	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 10:30	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 10:30	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 10:30	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 10:30	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 10:30	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 10:30	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 10:30	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 10:30	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 10:30	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 10:30	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 10:30	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 10:30	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 10:30	WG938408
(S) Toluene-d8	*101*			*90.0-115*		*12/26/2016 10:30*	WG938408
(S) Dibromofluoromethane	*103*			*79.0-121*		*12/26/2016 10:30*	WG938408
(S) 4-Bromofluorobenzene	*97.7*			*80.1-120*		*12/26/2016 10:30*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	408000		2820	10000	1	12/27/2016 14:58	WG938475

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:10	WG938033

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:53	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	2490		51.9	1000	1	12/27/2016 18:03	WG938354
Sulfate	14200		77.4	5000	1	12/27/2016 18:03	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	579	B J	102	1000	1	12/29/2016 12:57	WG938825

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 15:42	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/30/2016 00:15	WG938047
Arsenic	0.733	J	0.250	2.00	1	12/30/2016 00:15	WG938047
Barium	27.7		0.360	5.00	1	12/30/2016 00:15	WG938047
Beryllium	U		0.120	2.00	1	12/30/2016 00:15	WG938047
Cadmium	U		0.160	1.00	1	12/30/2016 00:15	WG938047
Chromium	1.18	J	0.540	2.00	1	12/30/2016 00:15	WG938047
Copper	2.56	B J	0.520	5.00	1	12/30/2016 00:15	WG938047
Cobalt	U		0.260	2.00	1	12/30/2016 00:15	WG938047
Iron	143		15.0	100	1	12/30/2016 00:15	WG938047
Lead	1.24	B J	0.240	2.00	1	12/30/2016 00:15	WG938047
Manganese	8.03		0.250	5.00	1	12/30/2016 00:15	WG938047
Nickel	1.53	J	0.350	2.00	1	12/30/2016 00:15	WG938047
Selenium	U		0.380	2.00	1	12/30/2016 00:15	WG938047
Silver	U		0.310	2.00	1	12/30/2016 00:15	WG938047
Thallium	U		0.190	2.00	1	12/30/2016 00:15	WG938047
Tin	0.472	B J	0.300	2.00	1	12/30/2016 00:15	WG938047
Vanadium	0.508	J	0.180	5.00	1	12/30/2016 00:15	WG938047
Zinc	5.23	B J	2.56	25.0	1	12/30/2016 00:15	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 10:53	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 10:53	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 10:53	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 10:53	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 10:53	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 10:53	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 10:53	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 10:53	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 10:53	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 10:53	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 10:53	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 10:53	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 10:53	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 10:53	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 10:53	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 10:53	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 10:53	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 10:53	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 10:53	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 10:53	WG938408
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 10:53	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 10:53	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 10:53	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 10:53	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 10:53	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 10:53	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 10:53	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 10:53	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 10:53	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 10:53	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 10:53	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 10:53	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 10:53	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 10:53	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 10:53	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 10:53	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 10:53	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 10:53	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 10:53	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 10:53	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 10:53	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 10:53	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 10:53	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 10:53	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 10:53	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 10:53	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 10:53	WG938408
(S) Toluene-d8	*102*			*90.0-115*		*12/26/2016 10:53*	WG938408
(S) Dibromofluoromethane	*104*			*79.0-121*		*12/26/2016 10:53*	WG938408
(S) 4-Bromofluorobenzene	*94.2*			*80.1-120*		*12/26/2016 10:53*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	366000		2820	10000	1	12/27/2016 14:58	WG938475

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:11	WG938033

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:54	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	15400		51.9	1000	1	12/27/2016 18:13	WG938354
Sulfate	10600		77.4	5000	1	12/27/2016 18:13	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	1530	B	102	1000	1	12/29/2016 13:46	WG938825

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 15:45	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/30/2016 00:19	WG938047
Arsenic	0.414	J	0.250	2.00	1	12/30/2016 00:19	WG938047
Barium	41.0		0.360	5.00	1	12/30/2016 00:19	WG938047
Beryllium	U		0.120	2.00	1	12/30/2016 00:19	WG938047
Cadmium	1.67		0.160	1.00	1	12/30/2016 00:19	WG938047
Chromium	1.74	J	0.540	2.00	1	12/30/2016 00:19	WG938047
Copper	3.51	B J	0.520	5.00	1	12/30/2016 00:19	WG938047
Cobalt	U		0.260	2.00	1	12/30/2016 00:19	WG938047
Iron	54.8	J	15.0	100	1	12/30/2016 00:19	WG938047
Lead	0.434	B J	0.240	2.00	1	12/30/2016 00:19	WG938047
Manganese	3.50	J	0.250	5.00	1	12/30/2016 00:19	WG938047
Nickel	4.65		0.350	2.00	1	12/30/2016 00:19	WG938047
Selenium	U		0.380	2.00	1	12/30/2016 00:19	WG938047
Silver	U		0.310	2.00	1	12/30/2016 00:19	WG938047
Thallium	U		0.190	2.00	1	12/30/2016 00:19	WG938047
Tin	0.552	B J	0.300	2.00	1	12/30/2016 00:19	WG938047
Vanadium	0.561	J	0.180	5.00	1	12/30/2016 00:19	WG938047
Zinc	283		2.56	25.0	1	12/30/2016 00:19	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 11:17	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 11:17	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 11:17	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 11:17	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 11:17	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 11:17	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 11:17	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 11:17	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 11:17	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 11:17	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 11:17	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 11:17	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 11:17	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 11:17	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 11:17	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 11:17	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 11:17	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 11:17	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 11:17	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 11:17	WG938408
1,1-Dichloroethane	1.97		0.259	1.00	1	12/26/2016 11:17	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 11:17	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 11:17	WG938408
cis-1,2-Dichloroethene	0.607	J	0.260	1.00	1	12/26/2016 11:17	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 11:17	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 11:17	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 11:17	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 11:17	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 11:17	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 11:17	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 11:17	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 11:17	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 11:17	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 11:17	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 11:17	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 11:17	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 11:17	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 11:17	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 11:17	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 11:17	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 11:17	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 11:17	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 11:17	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 11:17	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 11:17	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 11:17	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 11:17	WG938408
(S) Toluene-d8	*103*			*90.0-115*		*12/26/2016 11:17*	WG938408
(S) Dibromofluoromethane	*103*			*79.0-121*		*12/26/2016 11:17*	WG938408
(S) 4-Bromofluorobenzene	*96.8*			*80.1-120*		*12/26/2016 11:17*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	1660000		2820	10000	1	12/25/2016 00:15	WG938167

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:11	WG938033

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 09:58	WG938805

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	24900		51.9	1000	1	12/27/2016 18:23	WG938354
Sulfate	737000		1550	100000	20	12/28/2016 12:47	WG939061

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	12400		102	1000	1	12/29/2016 14:15	WG938825

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 15:47	WG938137

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/29/2016 22:24	WG938047
Arsenic	1.99	J	0.250	2.00	1	12/29/2016 22:24	WG938047
Barium	18.5	O1	0.360	5.00	1	12/29/2016 22:24	WG938047
Beryllium	U		0.120	2.00	1	12/29/2016 22:24	WG938047
Cadmium	U		0.160	1.00	1	12/29/2016 22:24	WG938047
Chromium	U		0.540	2.00	1	12/29/2016 22:24	WG938047
Copper	1.46	B J	0.520	5.00	1	12/29/2016 22:24	WG938047
Cobalt	0.603	J	0.260	2.00	1	12/29/2016 22:24	WG938047
Iron	600		15.0	100	1	12/29/2016 22:24	WG938047
Lead	1.23	B J	0.240	2.00	1	12/29/2016 22:24	WG938047
Manganese	22.0		0.250	5.00	1	12/29/2016 22:24	WG938047
Nickel	7.19		0.350	2.00	1	12/29/2016 22:24	WG938047
Selenium	0.467	J	0.380	2.00	1	12/29/2016 22:24	WG938047
Silver	U		0.310	2.00	1	12/29/2016 22:24	WG938047
Thallium	U		0.190	2.00	1	12/29/2016 22:24	WG938047
Tin	0.305	B J	0.300	2.00	1	12/29/2016 22:24	WG938047
Vanadium	0.924	J	0.180	5.00	1	12/29/2016 22:24	WG938047
Zinc	79.7		2.56	25.0	1	12/29/2016 22:24	WG938047

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 11:40	WG938408
Acrylonitrile	U		1.87	10.0	1	12/26/2016 11:40	WG938408
Benzene	U		0.331	1.00	1	12/26/2016 11:40	WG938408
Bromochloromethane	U	J4	0.520	1.00	1	12/26/2016 11:40	WG938408
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 11:40	WG938408
Bromoform	U		0.469	1.00	1	12/26/2016 11:40	WG938408
Bromomethane	U		0.866	5.00	1	12/26/2016 11:40	WG938408
Carbon disulfide	U		0.275	1.00	1	12/26/2016 11:40	WG938408
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 11:40	WG938408
Chlorobenzene	U		0.348	1.00	1	12/26/2016 11:40	WG938408
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 11:40	WG938408
Chloroethane	U	J4	0.453	5.00	1	12/26/2016 11:40	WG938408
Chloroform	U		0.324	5.00	1	12/26/2016 11:40	WG938408
Chloromethane	U		0.276	2.50	1	12/26/2016 11:40	WG938408
Dibromomethane	U		0.346	1.00	1	12/26/2016 11:40	WG938408
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 11:40	WG938408
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 11:40	WG938408
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 11:40	WG938408
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 11:40	WG938408
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 11:40	WG938408
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 11:40	WG938408
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 11:40	WG938408
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 11:40	WG938408
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 11:40	WG938408
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 11:40	WG938408
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 11:40	WG938408
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 11:40	WG938408
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 11:40	WG938408
Ethylbenzene	U		0.384	1.00	1	12/26/2016 11:40	WG938408
2-Hexanone	U		3.82	10.0	1	12/26/2016 11:40	WG938408
Iodomethane	U		1.71	10.0	1	12/26/2016 11:40	WG938408
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 11:40	WG938408
Methylene Chloride	U		1.00	5.00	1	12/26/2016 11:40	WG938408
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 11:40	WG938408
Styrene	U		0.307	1.00	1	12/26/2016 11:40	WG938408
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 11:40	WG938408
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 11:40	WG938408
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 11:40	WG938408
Toluene	U		0.412	1.00	1	12/26/2016 11:40	WG938408
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 11:40	WG938408
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 11:40	WG938408
Trichloroethene	U		0.398	1.00	1	12/26/2016 11:40	WG938408
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 11:40	WG938408
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 11:40	WG938408
Vinyl acetate	U		1.63	10.0	1	12/26/2016 11:40	WG938408
Vinyl chloride	U		0.259	1.00	1	12/26/2016 11:40	WG938408
Xylenes, Total	U		1.06	3.00	1	12/26/2016 11:40	WG938408
(S) Toluene-d8	*102*			*90.0-115*		*12/26/2016 11:40*	WG938408
(S) Dibromofluoromethane	*103*			*79.0-121*		*12/26/2016 11:40*	WG938408
(S) 4-Bromofluorobenzene	*95.6*			*80.1-120*		*12/26/2016 11:40*	WG938408

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	347000		2820	10000	1	12/27/2016 14:58	WG938475

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:11	WG938033

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 17:43	WG939301

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	5540		51.9	1000	1	12/27/2016 18:33	WG938354
Sulfate	15300		77.4	5000	1	12/27/2016 18:33	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	1600		102	1000	1	12/31/2016 11:59	WG939610

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 17:37	WG938138

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/30/2016 01:36	WG938049
Arsenic	0.766	J	0.250	2.00	1	12/30/2016 01:36	WG938049
Barium	128		0.360	5.00	1	12/30/2016 01:36	WG938049
Beryllium	U		0.120	2.00	1	12/30/2016 01:36	WG938049
Cadmium	0.181	J	0.160	1.00	1	12/30/2016 01:36	WG938049
Chromium	U		0.540	2.00	1	12/30/2016 01:36	WG938049
Copper	0.727	B J	0.520	5.00	1	12/30/2016 01:36	WG938049
Cobalt	0.394	J	0.260	2.00	1	12/30/2016 01:36	WG938049
Iron	62.0	B J	15.0	100	1	12/30/2016 01:36	WG938049
Lead	0.515	B J	0.240	2.00	1	12/30/2016 01:36	WG938049
Manganese	32.9		0.250	5.00	1	12/30/2016 01:36	WG938049
Nickel	3.64		0.350	2.00	1	12/30/2016 01:36	WG938049
Selenium	U		0.380	2.00	1	12/30/2016 01:36	WG938049
Silver	U		0.310	2.00	1	12/30/2016 01:36	WG938049
Thallium	U		0.190	2.00	1	12/30/2016 01:36	WG938049
Tin	0.476	B J	0.300	2.00	1	12/30/2016 01:36	WG938049
Vanadium	0.460	J	0.180	5.00	1	12/30/2016 01:36	WG938049
Zinc	41.1		2.56	25.0	1	12/30/2016 10:18	WG939602

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 00:57	WG938588
Acrylonitrile	U		1.87	10.0	1	12/26/2016 00:57	WG938588
Benzene	U		0.331	1.00	1	12/26/2016 00:57	WG938588
Bromochloromethane	U		0.520	1.00	1	12/26/2016 00:57	WG938588
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 00:57	WG938588
Bromoform	U		0.469	1.00	1	12/26/2016 00:57	WG938588
Bromomethane	U		0.866	5.00	1	12/26/2016 00:57	WG938588
Carbon disulfide	U		0.275	1.00	1	12/26/2016 00:57	WG938588
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 00:57	WG938588
Chlorobenzene	U		0.348	1.00	1	12/26/2016 00:57	WG938588
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 00:57	WG938588
Chloroethane	U		0.453	5.00	1	12/26/2016 00:57	WG938588
Chloroform	U		0.324	5.00	1	12/26/2016 00:57	WG938588
Chloromethane	U		0.276	2.50	1	12/26/2016 00:57	WG938588
Dibromomethane	U		0.346	1.00	1	12/26/2016 00:57	WG938588
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 00:57	WG938588
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 00:57	WG938588
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 00:57	WG938588
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 00:57	WG938588
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 00:57	WG938588
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 00:57	WG938588
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 00:57	WG938588
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 00:57	WG938588
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 00:57	WG938588
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 00:57	WG938588
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 00:57	WG938588
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 00:57	WG938588
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 00:57	WG938588
Ethylbenzene	U		0.384	1.00	1	12/26/2016 00:57	WG938588
2-Hexanone	U		3.82	10.0	1	12/26/2016 00:57	WG938588
Iodomethane	U		1.71	10.0	1	12/26/2016 00:57	WG938588
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 00:57	WG938588
Methylene Chloride	U		1.00	5.00	1	12/26/2016 00:57	WG938588
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 00:57	WG938588
Styrene	U		0.307	1.00	1	12/26/2016 00:57	WG938588
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 00:57	WG938588
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 00:57	WG938588
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 00:57	WG938588
Toluene	U		0.412	1.00	1	12/26/2016 00:57	WG938588
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 00:57	WG938588
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 00:57	WG938588
Trichloroethene	U		0.398	1.00	1	12/26/2016 00:57	WG938588
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 00:57	WG938588
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 00:57	WG938588
Vinyl acetate	U		1.63	10.0	1	12/26/2016 00:57	WG938588
Vinyl chloride	U		0.259	1.00	1	12/26/2016 00:57	WG938588
Xylenes, Total	U		1.06	3.00	1	12/26/2016 00:57	WG938588
(S) Toluene-d8	*98.5*			*90.0-115*		*12/26/2016 00:57*	WG938588
(S) Dibromofluoromethane	*101*			*79.0-121*		*12/26/2016 00:57*	WG938588
(S) 4-Bromofluorobenzene	*100*			*80.1-120*		*12/26/2016 00:57*	WG938588

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	398000		2820	10000	1	12/28/2016 14:22	WG938690

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:12	WG938033

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 17:44	WG939301

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfate	22400		77.4	5000	1	12/28/2016 12:57	WG939061

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	663	J	102	1000	1	12/31/2016 12:34	WG939610

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 17:39	WG938138

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/30/2016 01:40	WG938049
Arsenic	1.57	J	0.250	2.00	1	12/30/2016 01:40	WG938049
Barium	32.8		0.360	5.00	1	12/30/2016 01:40	WG938049
Beryllium	U		0.120	2.00	1	12/30/2016 01:40	WG938049
Cadmium	U		0.160	1.00	1	12/30/2016 01:40	WG938049
Chromium	3.12		0.540	2.00	1	12/30/2016 01:40	WG938049
Copper	1.89	B J	0.520	5.00	1	12/30/2016 01:40	WG938049
Cobalt	0.682	J	0.260	2.00	1	12/30/2016 01:40	WG938049
Iron	345		15.0	100	1	12/30/2016 01:40	WG938049
Lead	1.02	B J	0.240	2.00	1	12/30/2016 01:40	WG938049
Manganese	15.3		0.250	5.00	1	12/30/2016 01:40	WG938049
Nickel	7.82		0.350	2.00	1	12/30/2016 01:40	WG938049
Selenium	U		0.380	2.00	1	12/30/2016 01:40	WG938049
Silver	U		0.310	2.00	1	12/30/2016 01:40	WG938049
Thallium	U		0.190	2.00	1	12/30/2016 01:40	WG938049
Tin	0.323	B J	0.300	2.00	1	12/30/2016 01:40	WG938049
Vanadium	0.521	J	0.180	5.00	1	12/30/2016 01:40	WG938049
Zinc	202		2.56	25.0	1	12/30/2016 10:22	WG939602

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 01:21	WG938588
Acrylonitrile	U		1.87	10.0	1	12/26/2016 01:21	WG938588
Benzene	U		0.331	1.00	1	12/26/2016 01:21	WG938588
Bromochloromethane	U		0.520	1.00	1	12/26/2016 01:21	WG938588
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 01:21	WG938588
Bromoform	U		0.469	1.00	1	12/26/2016 01:21	WG938588
Bromomethane	U		0.866	5.00	1	12/26/2016 01:21	WG938588
Carbon disulfide	U		0.275	1.00	1	12/26/2016 01:21	WG938588
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 01:21	WG938588
Chlorobenzene	U		0.348	1.00	1	12/26/2016 01:21	WG938588
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 01:21	WG938588
Chloroethane	U		0.453	5.00	1	12/26/2016 01:21	WG938588
Chloroform	U		0.324	5.00	1	12/26/2016 01:21	WG938588
Chloromethane	U		0.276	2.50	1	12/26/2016 01:21	WG938588
Dibromomethane	U		0.346	1.00	1	12/26/2016 01:21	WG938588
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 01:21	WG938588
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 01:21	WG938588
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 01:21	WG938588
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 01:21	WG938588
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 01:21	WG938588
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 01:21	WG938588
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 01:21	WG938588
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 01:21	WG938588
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 01:21	WG938588
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 01:21	WG938588
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 01:21	WG938588
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 01:21	WG938588
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 01:21	WG938588
Ethylbenzene	U		0.384	1.00	1	12/26/2016 01:21	WG938588
2-Hexanone	U		3.82	10.0	1	12/26/2016 01:21	WG938588
Iodomethane	U		1.71	10.0	1	12/26/2016 01:21	WG938588
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 01:21	WG938588
Methylene Chloride	U		1.00	5.00	1	12/26/2016 01:21	WG938588
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 01:21	WG938588
Styrene	U		0.307	1.00	1	12/26/2016 01:21	WG938588
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 01:21	WG938588
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 01:21	WG938588
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 01:21	WG938588
Toluene	U		0.412	1.00	1	12/26/2016 01:21	WG938588
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 01:21	WG938588
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 01:21	WG938588
Trichloroethene	U		0.398	1.00	1	12/26/2016 01:21	WG938588
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 01:21	WG938588
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 01:21	WG938588
Vinyl acetate	U		1.63	10.0	1	12/26/2016 01:21	WG938588
Vinyl chloride	U		0.259	1.00	1	12/26/2016 01:21	WG938588
Xylenes, Total	U		1.06	3.00	1	12/26/2016 01:21	WG938588
(S) Toluene-d8	*99.6*			*90.0-115*		*12/26/2016 01:21*	WG938588
(S) Dibromofluoromethane	*98.8*			*79.0-121*		*12/26/2016 01:21*	WG938588
(S) 4-Bromofluorobenzene	*102*			*80.1-120*		*12/26/2016 01:21*	WG938588

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	928000		2820	10000	1	12/27/2016 14:58	WG938475

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:12	WG938033

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 17:47	WG939301

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	102000		104	2000	2	12/27/2016 19:34	WG938354
Sulfate	23700		155	10000	2	12/27/2016 19:34	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	7520		102	1000	1	12/31/2016 12:52	WG939610

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	0.715		0.0490	0.200	1	12/23/2016 17:41	WG938138

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/30/2016 01:43	WG938049
Arsenic	62.7		0.250	2.00	1	12/30/2016 01:43	WG938049
Barium	173		0.360	5.00	1	12/30/2016 01:43	WG938049
Beryllium	U		0.120	2.00	1	12/30/2016 01:43	WG938049
Cadmium	0.239	J	0.160	1.00	1	12/30/2016 01:43	WG938049
Chromium	1.68	J	0.540	2.00	1	12/30/2016 01:43	WG938049
Copper	4.93	B J	0.520	5.00	1	12/30/2016 01:43	WG938049
Cobalt	69.2		0.260	2.00	1	12/30/2016 01:43	WG938049
Iron	23700		15.0	100	1	12/30/2016 01:43	WG938049
Lead	2.41	B	0.240	2.00	1	12/30/2016 01:43	WG938049
Manganese	346		0.250	5.00	1	12/30/2016 01:43	WG938049
Nickel	80.2		0.350	2.00	1	12/30/2016 01:43	WG938049
Selenium	U		0.380	2.00	1	12/30/2016 01:43	WG938049
Silver	U		0.310	2.00	1	12/30/2016 01:43	WG938049
Thallium	1.64	J	0.190	2.00	1	12/30/2016 01:43	WG938049
Tin	1.43	B J	0.300	2.00	1	12/30/2016 01:43	WG938049
Vanadium	0.450	J	0.180	5.00	1	12/30/2016 01:43	WG938049
Zinc	1600		5.12	50.0	2	12/30/2016 11:12	WG939602

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 01:45	WG938588
Acrylonitrile	U		1.87	10.0	1	12/26/2016 01:45	WG938588
Benzene	0.857	J	0.331	1.00	1	12/26/2016 01:45	WG938588
Bromochloromethane	U		0.520	1.00	1	12/26/2016 01:45	WG938588
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 01:45	WG938588
Bromoform	U		0.469	1.00	1	12/26/2016 01:45	WG938588
Bromomethane	U		0.866	5.00	1	12/26/2016 01:45	WG938588
Carbon disulfide	U		0.275	1.00	1	12/26/2016 01:45	WG938588
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 01:45	WG938588
Chlorobenzene	1.46		0.348	1.00	1	12/26/2016 01:45	WG938588
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 01:45	WG938588
Chloroethane	2.07	J	0.453	5.00	1	12/26/2016 01:45	WG938588
Chloroform	U		0.324	5.00	1	12/26/2016 01:45	WG938588
Chloromethane	U		0.276	2.50	1	12/26/2016 01:45	WG938588
Dibromomethane	U		0.346	1.00	1	12/26/2016 01:45	WG938588
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 01:45	WG938588
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 01:45	WG938588
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 01:45	WG938588
1,4-Dichlorobenzene	2.63		0.274	1.00	1	12/26/2016 01:45	WG938588
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 01:45	WG938588
1,1-Dichloroethane	16.2		0.259	1.00	1	12/26/2016 01:45	WG938588
1,2-Dichloroethane	0.369	J	0.361	1.00	1	12/26/2016 01:45	WG938588
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 01:45	WG938588
cis-1,2-Dichloroethene	10.2		0.260	1.00	1	12/26/2016 01:45	WG938588
trans-1,2-Dichloroethene	0.405	J	0.396	1.00	1	12/26/2016 01:45	WG938588
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 01:45	WG938588
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 01:45	WG938588
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 01:45	WG938588
Ethylbenzene	U		0.384	1.00	1	12/26/2016 01:45	WG938588
2-Hexanone	U		3.82	10.0	1	12/26/2016 01:45	WG938588
Iodomethane	U		1.71	10.0	1	12/26/2016 01:45	WG938588
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 01:45	WG938588
Methylene Chloride	U		1.00	5.00	1	12/26/2016 01:45	WG938588
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 01:45	WG938588
Styrene	U		0.307	1.00	1	12/26/2016 01:45	WG938588
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 01:45	WG938588
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 01:45	WG938588
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 01:45	WG938588
Toluene	U		0.412	1.00	1	12/26/2016 01:45	WG938588
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 01:45	WG938588
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 01:45	WG938588
Trichloroethene	1.11		0.398	1.00	1	12/26/2016 01:45	WG938588
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 01:45	WG938588
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 01:45	WG938588
Vinyl acetate	U		1.63	10.0	1	12/26/2016 01:45	WG938588
Vinyl chloride	3.69		0.259	1.00	1	12/26/2016 01:45	WG938588
Xylenes, Total	U		1.06	3.00	1	12/26/2016 01:45	WG938588
(S) Toluene-d8	*99.9*			*90.0-115*		*12/26/2016 01:45*	WG938588
(S) Dibromofluoromethane	*97.2*			*79.0-121*		*12/26/2016 01:45*	WG938588
(S) 4-Bromofluorobenzene	*101*			*80.1-120*		*12/26/2016 01:45*	WG938588

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	275000		2820	10000	1	12/28/2016 14:22	WG938690

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:12	WG938033

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 17:48	WG939301

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	16100		51.9	1000	1	12/27/2016 19:44	WG938354
Sulfate	5170		77.4	5000	1	12/27/2016 19:44	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	2370		102	1000	1	12/31/2016 13:10	WG939610

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 17:43	WG938138

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/30/2016 01:47	WG938049
Arsenic	4.73		0.250	2.00	1	12/30/2016 01:47	WG938049
Barium	72.7		0.360	5.00	1	12/30/2016 01:47	WG938049
Beryllium	U		0.120	2.00	1	12/30/2016 01:47	WG938049
Cadmium	U		0.160	1.00	1	12/30/2016 01:47	WG938049
Chromium	U		0.540	2.00	1	12/30/2016 01:47	WG938049
Copper	0.792	B J	0.520	5.00	1	12/30/2016 01:47	WG938049
Cobalt	2.71		0.260	2.00	1	12/30/2016 01:47	WG938049
Iron	2050		15.0	100	1	12/30/2016 01:47	WG938049
Lead	0.428	B J	0.240	2.00	1	12/30/2016 01:47	WG938049
Manganese	1140		0.250	5.00	1	12/30/2016 01:47	WG938049
Nickel	3.35		0.350	2.00	1	12/30/2016 01:47	WG938049
Selenium	U		0.380	2.00	1	12/30/2016 01:47	WG938049
Silver	U		0.310	2.00	1	12/30/2016 01:47	WG938049
Thallium	U		0.190	2.00	1	12/30/2016 01:47	WG938049
Tin	U		0.300	2.00	1	12/30/2016 01:47	WG938049
Vanadium	0.281	J	0.180	5.00	1	12/30/2016 01:47	WG938049
Zinc	6.14	B J J3 J4	2.56	25.0	1	12/30/2016 01:47	WG938049

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 02:09	WG938588
Acrylonitrile	U		1.87	10.0	1	12/26/2016 02:09	WG938588
Benzene	U		0.331	1.00	1	12/26/2016 02:09	WG938588
Bromochloromethane	U		0.520	1.00	1	12/26/2016 02:09	WG938588
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 02:09	WG938588
Bromoform	U		0.469	1.00	1	12/26/2016 02:09	WG938588
Bromomethane	U		0.866	5.00	1	12/26/2016 02:09	WG938588
Carbon disulfide	U		0.275	1.00	1	12/26/2016 02:09	WG938588
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 02:09	WG938588
Chlorobenzene	U		0.348	1.00	1	12/26/2016 02:09	WG938588
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 02:09	WG938588
Chloroethane	U		0.453	5.00	1	12/26/2016 02:09	WG938588
Chloroform	U		0.324	5.00	1	12/26/2016 02:09	WG938588
Chloromethane	U		0.276	2.50	1	12/26/2016 02:09	WG938588
Dibromomethane	U		0.346	1.00	1	12/26/2016 02:09	WG938588
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 02:09	WG938588
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 02:09	WG938588
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 02:09	WG938588
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 02:09	WG938588
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 02:09	WG938588
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 02:09	WG938588
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 02:09	WG938588
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 02:09	WG938588
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 02:09	WG938588
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 02:09	WG938588
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 02:09	WG938588
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 02:09	WG938588
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 02:09	WG938588
Ethylbenzene	U		0.384	1.00	1	12/26/2016 02:09	WG938588
2-Hexanone	U		3.82	10.0	1	12/26/2016 02:09	WG938588
Iodomethane	U		1.71	10.0	1	12/26/2016 02:09	WG938588
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 02:09	WG938588
Methylene Chloride	U		1.00	5.00	1	12/26/2016 02:09	WG938588
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 02:09	WG938588
Styrene	U		0.307	1.00	1	12/26/2016 02:09	WG938588
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 02:09	WG938588
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 02:09	WG938588
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 02:09	WG938588
Toluene	U		0.412	1.00	1	12/26/2016 02:09	WG938588
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 02:09	WG938588
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 02:09	WG938588
Trichloroethene	U		0.398	1.00	1	12/26/2016 02:09	WG938588
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 02:09	WG938588
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 02:09	WG938588
Vinyl acetate	U		1.63	10.0	1	12/26/2016 02:09	WG938588
Vinyl chloride	U		0.259	1.00	1	12/26/2016 02:09	WG938588
Xylenes, Total	U		1.06	3.00	1	12/26/2016 02:09	WG938588
(S) Toluene-d8	*100*			*90.0-115*		*12/26/2016 02:09*	WG938588
(S) Dibromofluoromethane	*99.2*			*79.0-121*		*12/26/2016 02:09*	WG938588
(S) 4-Bromofluorobenzene	*102*			*80.1-120*		*12/26/2016 02:09*	WG938588

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	359000		2820	10000	1	12/28/2016 14:22	WG938690

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:12	WG938033

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 17:49	WG939301

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	2370		51.9	1000	1	12/27/2016 19:54	WG938354
Sulfate	8170		77.4	5000	1	12/27/2016 19:54	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	967	J	102	1000	1	12/31/2016 13:29	WG939610

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 17:46	WG938138

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/30/2016 01:50	WG938049
Arsenic	0.375	J	0.250	2.00	1	12/30/2016 01:50	WG938049
Barium	41.8		0.360	5.00	1	12/30/2016 01:50	WG938049
Beryllium	U		0.120	2.00	1	12/30/2016 01:50	WG938049
Cadmium	1.61		0.160	1.00	1	12/30/2016 01:50	WG938049
Chromium	U		0.540	2.00	1	12/30/2016 01:50	WG938049
Copper	1.19	B J	0.520	5.00	1	12/30/2016 01:50	WG938049
Cobalt	1.69	J	0.260	2.00	1	12/30/2016 01:50	WG938049
Iron	118	B	15.0	100	1	12/30/2016 01:50	WG938049
Lead	0.597	B J	0.240	2.00	1	12/30/2016 01:50	WG938049
Manganese	327		0.250	5.00	1	12/30/2016 01:50	WG938049
Nickel	4.46		0.350	2.00	1	12/30/2016 01:50	WG938049
Selenium	U		0.380	2.00	1	12/30/2016 01:50	WG938049
Silver	U		0.310	2.00	1	12/30/2016 01:50	WG938049
Thallium	U		0.190	2.00	1	12/30/2016 01:50	WG938049
Tin	U		0.300	2.00	1	12/30/2016 01:50	WG938049
Vanadium	0.706	J	0.180	5.00	1	12/30/2016 01:50	WG938049
Zinc	122		2.56	25.0	1	12/30/2016 10:36	WG939602

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 15:30	WG938676
Acrylonitrile	U		1.87	10.0	1	12/26/2016 15:30	WG938676
Benzene	U		0.331	1.00	1	12/26/2016 15:30	WG938676
Bromochloromethane	U		0.520	1.00	1	12/26/2016 15:30	WG938676
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 15:30	WG938676
Bromoform	U		0.469	1.00	1	12/26/2016 15:30	WG938676
Bromomethane	U		0.866	5.00	1	12/26/2016 15:30	WG938676
Carbon disulfide	U		0.275	1.00	1	12/26/2016 15:30	WG938676
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 15:30	WG938676
Chlorobenzene	U		0.348	1.00	1	12/26/2016 15:30	WG938676
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 15:30	WG938676
Chloroethane	U		0.453	5.00	1	12/26/2016 15:30	WG938676
Chloroform	U		0.324	5.00	1	12/26/2016 15:30	WG938676
Chloromethane	U		0.276	2.50	1	12/26/2016 15:30	WG938676
Dibromomethane	U		0.346	1.00	1	12/26/2016 15:30	WG938676
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 15:30	WG938676
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 15:30	WG938676
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 15:30	WG938676
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 15:30	WG938676
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 15:30	WG938676
1,1-Dichloroethane	2.92		0.259	1.00	1	12/26/2016 15:30	WG938676
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 15:30	WG938676
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 15:30	WG938676
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 15:30	WG938676
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 15:30	WG938676
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 15:30	WG938676
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 15:30	WG938676
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 15:30	WG938676
Ethylbenzene	U		0.384	1.00	1	12/26/2016 15:30	WG938676
2-Hexanone	U		3.82	10.0	1	12/26/2016 15:30	WG938676
Iodomethane	U		1.71	10.0	1	12/26/2016 15:30	WG938676
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 15:30	WG938676
Methylene Chloride	U		1.00	5.00	1	12/26/2016 15:30	WG938676
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 15:30	WG938676
Styrene	U		0.307	1.00	1	12/26/2016 15:30	WG938676
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 15:30	WG938676
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 15:30	WG938676
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 15:30	WG938676
Toluene	U		0.412	1.00	1	12/26/2016 15:30	WG938676
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 15:30	WG938676
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 15:30	WG938676
Trichloroethene	U		0.398	1.00	1	12/26/2016 15:30	WG938676
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 15:30	WG938676
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 15:30	WG938676
Vinyl acetate	U		1.63	10.0	1	12/26/2016 15:30	WG938676
Vinyl chloride	U		0.259	1.00	1	12/26/2016 15:30	WG938676
Xylenes, Total	U		1.06	3.00	1	12/26/2016 15:30	WG938676
(S) Toluene-d8	*103*			*90.0-115*		*12/26/2016 15:30*	WG938676
(S) Dibromofluoromethane	*105*			*79.0-121*		*12/26/2016 15:30*	WG938676
(S) 4-Bromofluorobenzene	*113*			*80.1-120*		*12/26/2016 15:30*	WG938676

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	335000		2820	10000	1	12/28/2016 14:22	WG938690

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:13	WG938033

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 17:52	WG939301

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	12700		51.9	1000	1	12/27/2016 20:05	WG938354
Sulfate	4410	J	77.4	5000	1	12/27/2016 20:05	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	2380		102	1000	1	12/31/2016 13:50	WG939610

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 17:48	WG938138

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/30/2016 01:54	WG938049
Arsenic	3.00		0.250	2.00	1	12/30/2016 01:54	WG938049
Barium	79.8		0.360	5.00	1	12/30/2016 01:54	WG938049
Beryllium	U		0.120	2.00	1	12/30/2016 01:54	WG938049
Cadmium	U		0.160	1.00	1	12/30/2016 01:54	WG938049
Chromium	U		0.540	2.00	1	12/30/2016 01:54	WG938049
Copper	U		0.520	5.00	1	12/30/2016 01:54	WG938049
Cobalt	2.41		0.260	2.00	1	12/30/2016 01:54	WG938049
Iron	2010		15.0	100	1	12/30/2016 01:54	WG938049
Lead	0.245	B J	0.240	2.00	1	12/30/2016 01:54	WG938049
Manganese	914		0.250	5.00	1	12/30/2016 01:54	WG938049
Nickel	2.36		0.350	2.00	1	12/30/2016 01:54	WG938049
Selenium	U		0.380	2.00	1	12/30/2016 01:54	WG938049
Silver	U		0.310	2.00	1	12/30/2016 01:54	WG938049
Thallium	U		0.190	2.00	1	12/30/2016 01:54	WG938049
Tin	0.379	B J	0.300	2.00	1	12/30/2016 01:54	WG938049
Vanadium	0.288	J	0.180	5.00	1	12/30/2016 01:54	WG938049
Zinc	U	J3 J4	2.56	25.0	1	12/30/2016 01:54	WG938049

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 15:43	WG938676
Acrylonitrile	U		1.87	10.0	1	12/26/2016 15:43	WG938676
Benzene	U		0.331	1.00	1	12/26/2016 15:43	WG938676
Bromochloromethane	U		0.520	1.00	1	12/26/2016 15:43	WG938676
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 15:43	WG938676
Bromoform	U		0.469	1.00	1	12/26/2016 15:43	WG938676
Bromomethane	U		0.866	5.00	1	12/26/2016 15:43	WG938676
Carbon disulfide	U		0.275	1.00	1	12/26/2016 15:43	WG938676
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 15:43	WG938676
Chlorobenzene	U		0.348	1.00	1	12/26/2016 15:43	WG938676
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 15:43	WG938676
Chloroethane	2.81	J	0.453	5.00	1	12/26/2016 15:43	WG938676
Chloroform	U		0.324	5.00	1	12/26/2016 15:43	WG938676
Chloromethane	U		0.276	2.50	1	12/26/2016 15:43	WG938676
Dibromomethane	U		0.346	1.00	1	12/26/2016 15:43	WG938676
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 15:43	WG938676
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 15:43	WG938676
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 15:43	WG938676
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 15:43	WG938676
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 15:43	WG938676
1,1-Dichloroethane	0.862	J	0.259	1.00	1	12/26/2016 15:43	WG938676
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 15:43	WG938676
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 15:43	WG938676
cis-1,2-Dichloroethene	0.477	J	0.260	1.00	1	12/26/2016 15:43	WG938676
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 15:43	WG938676
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 15:43	WG938676
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 15:43	WG938676
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 15:43	WG938676
Ethylbenzene	U		0.384	1.00	1	12/26/2016 15:43	WG938676
2-Hexanone	U		3.82	10.0	1	12/26/2016 15:43	WG938676
Iodomethane	U		1.71	10.0	1	12/26/2016 15:43	WG938676
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 15:43	WG938676
Methylene Chloride	U		1.00	5.00	1	12/26/2016 15:43	WG938676
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 15:43	WG938676
Styrene	U		0.307	1.00	1	12/26/2016 15:43	WG938676
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 15:43	WG938676
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 15:43	WG938676
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 15:43	WG938676
Toluene	U		0.412	1.00	1	12/26/2016 15:43	WG938676
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 15:43	WG938676
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 15:43	WG938676
Trichloroethene	U		0.398	1.00	1	12/26/2016 15:43	WG938676
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 15:43	WG938676
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 15:43	WG938676
Vinyl acetate	U		1.63	10.0	1	12/26/2016 15:43	WG938676
Vinyl chloride	U		0.259	1.00	1	12/26/2016 15:43	WG938676
Xylenes, Total	U		1.06	3.00	1	12/26/2016 15:43	WG938676
(S) Toluene-d8	*102*			*90.0-115*		*12/26/2016 15:43*	WG938676
(S) Dibromofluoromethane	*105*			*79.0-121*		*12/26/2016 15:43*	WG938676
(S) 4-Bromofluorobenzene	*115*			*80.1-120*		*12/26/2016 15:43*	WG938676

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	319000		2820	10000	1	12/28/2016 14:22	WG938690

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:14	WG938033

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 17:54	WG939301

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	45000		51.9	1000	1	12/27/2016 20:15	WG938354
Sulfate	2800	J	77.4	5000	1	12/27/2016 20:15	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	4230		102	1000	1	12/31/2016 15:26	WG939610

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 17:50	WG938138

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/30/2016 01:57	WG938049
Arsenic	0.799	J	0.250	2.00	1	12/30/2016 01:57	WG938049
Barium	66.1		0.360	5.00	1	12/30/2016 01:57	WG938049
Beryllium	U		0.120	2.00	1	12/30/2016 01:57	WG938049
Cadmium	U		0.160	1.00	1	12/30/2016 01:57	WG938049
Chromium	U		0.540	2.00	1	12/30/2016 01:57	WG938049
Copper	0.709	B J	0.520	5.00	1	12/30/2016 01:57	WG938049
Cobalt	1.08	J	0.260	2.00	1	12/30/2016 01:57	WG938049
Iron	567		15.0	100	1	12/30/2016 01:57	WG938049
Lead	0.762	B J	0.240	2.00	1	12/30/2016 01:57	WG938049
Manganese	464		0.250	5.00	1	12/30/2016 01:57	WG938049
Nickel	3.33		0.350	2.00	1	12/30/2016 01:57	WG938049
Selenium	U		0.380	2.00	1	12/30/2016 01:57	WG938049
Silver	U		0.310	2.00	1	12/30/2016 01:57	WG938049
Thallium	U		0.190	2.00	1	12/30/2016 01:57	WG938049
Tin	0.322	B J	0.300	2.00	1	12/30/2016 01:57	WG938049
Vanadium	0.474	J	0.180	5.00	1	12/30/2016 01:57	WG938049
Zinc	2.99	B J J3 J4	2.56	25.0	1	12/30/2016 01:57	WG938049

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 15:57	WG938676
Acrylonitrile	U		1.87	10.0	1	12/26/2016 15:57	WG938676
Benzene	U		0.331	1.00	1	12/26/2016 15:57	WG938676
Bromochloromethane	U		0.520	1.00	1	12/26/2016 15:57	WG938676
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 15:57	WG938676
Bromoform	U		0.469	1.00	1	12/26/2016 15:57	WG938676
Bromomethane	U		0.866	5.00	1	12/26/2016 15:57	WG938676
Carbon disulfide	U		0.275	1.00	1	12/26/2016 15:57	WG938676
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 15:57	WG938676
Chlorobenzene	U		0.348	1.00	1	12/26/2016 15:57	WG938676
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 15:57	WG938676
Chloroethane	U		0.453	5.00	1	12/26/2016 15:57	WG938676
Chloroform	U		0.324	5.00	1	12/26/2016 15:57	WG938676
Chloromethane	U		0.276	2.50	1	12/26/2016 15:57	WG938676
Dibromomethane	U		0.346	1.00	1	12/26/2016 15:57	WG938676
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 15:57	WG938676
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 15:57	WG938676
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 15:57	WG938676
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 15:57	WG938676
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 15:57	WG938676
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 15:57	WG938676
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 15:57	WG938676
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 15:57	WG938676
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 15:57	WG938676
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 15:57	WG938676
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 15:57	WG938676
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 15:57	WG938676
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 15:57	WG938676
Ethylbenzene	U		0.384	1.00	1	12/26/2016 15:57	WG938676
2-Hexanone	U		3.82	10.0	1	12/26/2016 15:57	WG938676
Iodomethane	U		1.71	10.0	1	12/26/2016 15:57	WG938676
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 15:57	WG938676
Methylene Chloride	U		1.00	5.00	1	12/26/2016 15:57	WG938676
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 15:57	WG938676
Styrene	U		0.307	1.00	1	12/26/2016 15:57	WG938676
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 15:57	WG938676
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 15:57	WG938676
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 15:57	WG938676
Toluene	U		0.412	1.00	1	12/26/2016 15:57	WG938676
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 15:57	WG938676
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 15:57	WG938676
Trichloroethene	U		0.398	1.00	1	12/26/2016 15:57	WG938676
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 15:57	WG938676
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 15:57	WG938676
Vinyl acetate	U		1.63	10.0	1	12/26/2016 15:57	WG938676
Vinyl chloride	U		0.259	1.00	1	12/26/2016 15:57	WG938676
Xylenes, Total	U		1.06	3.00	1	12/26/2016 15:57	WG938676
(S) Toluene-d8	*103*			*90.0-115*		*12/26/2016 15:57*	WG938676
(S) Dibromofluoromethane	*104*			*79.0-121*		*12/26/2016 15:57*	WG938676
(S) 4-Bromofluorobenzene	*113*			*80.1-120*		*12/26/2016 15:57*	WG938676

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	1830000		2820	10000	1	12/27/2016 14:58	WG938475

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:14	WG938033

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 17:55	WG939301

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	659000		519	10000	10	12/27/2016 20:35	WG938354
Sulfate	11200		77.4	5000	1	12/27/2016 20:25	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	60000		102	1000	1	12/31/2016 15:46	WG939610

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 17:52	WG938138

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/30/2016 02:01	WG938049
Arsenic	9.12		0.250	2.00	1	12/30/2016 02:01	WG938049
Barium	2560		1.80	25.0	5	12/30/2016 02:24	WG938049
Beryllium	U		0.120	2.00	1	12/30/2016 02:01	WG938049
Cadmium	U		0.160	1.00	1	12/30/2016 02:01	WG938049
Chromium	30.7		0.540	2.00	1	12/30/2016 02:01	WG938049
Copper	12.2	B	0.520	5.00	1	12/30/2016 02:01	WG938049
Cobalt	6.24		0.260	2.00	1	12/30/2016 02:01	WG938049
Iron	49400		15.0	100	1	12/30/2016 02:01	WG938049
Lead	3.15	B	0.240	2.00	1	12/30/2016 02:01	WG938049
Manganese	1220		0.250	5.00	1	12/30/2016 02:01	WG938049
Nickel	49.9		0.350	2.00	1	12/30/2016 02:01	WG938049
Selenium	0.627	J	0.380	2.00	1	12/30/2016 02:01	WG938049
Silver	U		0.310	2.00	1	12/30/2016 02:01	WG938049
Thallium	U		0.190	2.00	1	12/30/2016 02:01	WG938049
Tin	6.32		0.300	2.00	1	12/30/2016 02:01	WG938049
Vanadium	2.08	J	0.180	5.00	1	12/30/2016 02:01	WG938049
Zinc	66.6		2.56	25.0	1	12/30/2016 10:39	WG939602

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	23.6	J	10.0	50.0	1	12/26/2016 16:10	WG938676
Acrylonitrile	U		1.87	10.0	1	12/26/2016 16:10	WG938676
Benzene	U		0.331	1.00	1	12/26/2016 16:10	WG938676
Bromochloromethane	U		0.520	1.00	1	12/26/2016 16:10	WG938676
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 16:10	WG938676
Bromoform	U		0.469	1.00	1	12/26/2016 16:10	WG938676
Bromomethane	U		0.866	5.00	1	12/26/2016 16:10	WG938676
Carbon disulfide	U		0.275	1.00	1	12/26/2016 16:10	WG938676
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 16:10	WG938676
Chlorobenzene	U		0.348	1.00	1	12/26/2016 16:10	WG938676
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 16:10	WG938676
Chloroethane	0.565	J	0.453	5.00	1	12/26/2016 16:10	WG938676
Chloroform	U		0.324	5.00	1	12/26/2016 16:10	WG938676
Chloromethane	U		0.276	2.50	1	12/26/2016 16:10	WG938676
Dibromomethane	U		0.346	1.00	1	12/26/2016 16:10	WG938676
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 16:10	WG938676
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 16:10	WG938676
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 16:10	WG938676
1,4-Dichlorobenzene	1.40		0.274	1.00	1	12/26/2016 16:10	WG938676
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 16:10	WG938676
1,1-Dichloroethane	0.565	J	0.259	1.00	1	12/26/2016 16:10	WG938676
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 16:10	WG938676
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 16:10	WG938676
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 16:10	WG938676
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 16:10	WG938676
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 16:10	WG938676
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 16:10	WG938676
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 16:10	WG938676
Ethylbenzene	U		0.384	1.00	1	12/26/2016 16:10	WG938676
2-Hexanone	U		3.82	10.0	1	12/26/2016 16:10	WG938676
Iodomethane	U		1.71	10.0	1	12/26/2016 16:10	WG938676
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 16:10	WG938676
Methylene Chloride	U		1.00	5.00	1	12/26/2016 16:10	WG938676
4-Methyl-2-pentanone (MIBK)	2.87	J	2.14	10.0	1	12/26/2016 16:10	WG938676
Styrene	U		0.307	1.00	1	12/26/2016 16:10	WG938676
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 16:10	WG938676
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 16:10	WG938676
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 16:10	WG938676
Toluene	U		0.412	1.00	1	12/26/2016 16:10	WG938676
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 16:10	WG938676
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 16:10	WG938676
Trichloroethene	U		0.398	1.00	1	12/26/2016 16:10	WG938676
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 16:10	WG938676
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 16:10	WG938676
Vinyl acetate	U		1.63	10.0	1	12/26/2016 16:10	WG938676
Vinyl chloride	0.463	J	0.259	1.00	1	12/26/2016 16:10	WG938676
Xylenes, Total	2.04	J	1.06	3.00	1	12/26/2016 16:10	WG938676
(S) Toluene-d8	*104*			*90.0-115*		*12/26/2016 16:10*	WG938676
(S) Dibromofluoromethane	*105*			*79.0-121*		*12/26/2016 16:10*	WG938676
(S) 4-Bromofluorobenzene	*114*			*80.1-120*		*12/26/2016 16:10*	WG938676

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	459000		2820	10000	1	12/28/2016 14:22	WG938690

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:14	WG938033

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 17:56	WG939301

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	20800		51.9	1000	1	12/27/2016 20:45	WG938354
Sulfate	8320		77.4	5000	1	12/27/2016 20:45	WG938354

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	1390		102	1000	1	12/31/2016 16:01	WG939610

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 17:55	WG938138

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/30/2016 02:05	WG938049
Arsenic	0.422	J	0.250	2.00	1	12/30/2016 02:05	WG938049
Barium	48.2		0.360	5.00	1	12/30/2016 02:05	WG938049
Beryllium	U		0.120	2.00	1	12/30/2016 02:05	WG938049
Cadmium	0.470	J	0.160	1.00	1	12/30/2016 02:05	WG938049
Chromium	U		0.540	2.00	1	12/30/2016 02:05	WG938049
Copper	1.60	B J	0.520	5.00	1	12/30/2016 02:05	WG938049
Cobalt	U		0.260	2.00	1	12/30/2016 02:05	WG938049
Iron	63.7	B J	15.0	100	1	12/30/2016 02:05	WG938049
Lead	0.635	B J	0.240	2.00	1	12/30/2016 02:05	WG938049
Manganese	1.34	B J	0.250	5.00	1	12/30/2016 02:05	WG938049
Nickel	0.804	J	0.350	2.00	1	12/30/2016 02:05	WG938049
Selenium	U		0.380	2.00	1	12/30/2016 02:05	WG938049
Silver	U		0.310	2.00	1	12/30/2016 02:05	WG938049
Thallium	U		0.190	2.00	1	12/30/2016 02:05	WG938049
Tin	0.326	B J	0.300	2.00	1	12/30/2016 02:05	WG938049
Vanadium	0.306	J	0.180	5.00	1	12/30/2016 02:05	WG938049
Zinc	41.6		2.56	25.0	1	12/30/2016 10:43	WG939602

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 16:24	WG938676
Acrylonitrile	U		1.87	10.0	1	12/26/2016 16:24	WG938676
Benzene	U		0.331	1.00	1	12/26/2016 16:24	WG938676
Bromochloromethane	U		0.520	1.00	1	12/26/2016 16:24	WG938676
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 16:24	WG938676
Bromoform	U		0.469	1.00	1	12/26/2016 16:24	WG938676
Bromomethane	U		0.866	5.00	1	12/26/2016 16:24	WG938676
Carbon disulfide	U		0.275	1.00	1	12/26/2016 16:24	WG938676
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 16:24	WG938676
Chlorobenzene	U		0.348	1.00	1	12/26/2016 16:24	WG938676
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 16:24	WG938676
Chloroethane	U		0.453	5.00	1	12/26/2016 16:24	WG938676
Chloroform	U		0.324	5.00	1	12/26/2016 16:24	WG938676
Chloromethane	U		0.276	2.50	1	12/26/2016 16:24	WG938676
Dibromomethane	U		0.346	1.00	1	12/26/2016 16:24	WG938676
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 16:24	WG938676
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 16:24	WG938676
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 16:24	WG938676
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 16:24	WG938676
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 16:24	WG938676
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 16:24	WG938676
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 16:24	WG938676
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 16:24	WG938676
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 16:24	WG938676
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 16:24	WG938676
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 16:24	WG938676
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 16:24	WG938676
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 16:24	WG938676
Ethylbenzene	U		0.384	1.00	1	12/26/2016 16:24	WG938676
2-Hexanone	U		3.82	10.0	1	12/26/2016 16:24	WG938676
Iodomethane	U		1.71	10.0	1	12/26/2016 16:24	WG938676
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 16:24	WG938676
Methylene Chloride	U		1.00	5.00	1	12/26/2016 16:24	WG938676
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 16:24	WG938676
Styrene	U		0.307	1.00	1	12/26/2016 16:24	WG938676
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 16:24	WG938676
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 16:24	WG938676
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 16:24	WG938676
Toluene	U		0.412	1.00	1	12/26/2016 16:24	WG938676
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 16:24	WG938676
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 16:24	WG938676
Trichloroethene	U		0.398	1.00	1	12/26/2016 16:24	WG938676
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 16:24	WG938676
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 16:24	WG938676
Vinyl acetate	U		1.63	10.0	1	12/26/2016 16:24	WG938676
Vinyl chloride	U		0.259	1.00	1	12/26/2016 16:24	WG938676
Xylenes, Total	U		1.06	3.00	1	12/26/2016 16:24	WG938676
(S) Toluene-d8	*104*			*90.0-115*		*12/26/2016 16:24*	WG938676
(S) Dibromofluoromethane	*101*			*79.0-121*		*12/26/2016 16:24*	WG938676
(S) 4-Bromofluorobenzene	*111*			*80.1-120*		*12/26/2016 16:24*	WG938676

Gravimetric Analysis by Method 2540 C-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Dissolved Solids	26000		2820	10000	1	12/28/2016 14:22	WG938690

Wet Chemistry by Method 4500S2 D-2011

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Sulfide	U		6.50	50.0	1	12/22/2016 22:15	WG938033

Wet Chemistry by Method 9012B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Cyanide	U		1.80	5.00	1	12/29/2016 17:58	WG939301

Wet Chemistry by Method 9056A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Chloride	122	J	51.9	1000	1	12/27/2016 18:29	WG938517
Sulfate	U		77.4	5000	1	12/27/2016 18:29	WG938517

Wet Chemistry by Method 9060A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
TOC (Total Organic Carbon)	397	J	102	1000	1	12/31/2016 16:20	WG939610

Mercury by Method 7470A

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Mercury	U		0.0490	0.200	1	12/23/2016 18:04	WG938138

Metals (ICPMS) by Method 6020

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Antimony	U		0.754	2.00	1	12/30/2016 01:00	WG938049
Arsenic	U		0.250	2.00	1	12/30/2016 01:00	WG938049
Barium	U		0.360	5.00	1	12/30/2016 01:00	WG938049
Beryllium	U		0.120	2.00	1	12/30/2016 01:00	WG938049
Cadmium	U		0.160	1.00	1	12/30/2016 01:00	WG938049
Chromium	U		0.540	2.00	1	12/30/2016 01:00	WG938049
Copper	6.19	B	0.520	5.00	1	12/30/2016 01:00	WG938049
Cobalt	U		0.260	2.00	1	12/30/2016 01:00	WG938049
Iron	46.6	B J	15.0	100	1	12/30/2016 01:00	WG938049
Lead	0.543	B J	0.240	2.00	1	12/30/2016 01:00	WG938049
Manganese	0.294	B J	0.250	5.00	1	12/30/2016 01:00	WG938049
Nickel	U		0.350	2.00	1	12/30/2016 01:00	WG938049
Selenium	U		0.380	2.00	1	12/30/2016 01:00	WG938049
Silver	U		0.310	2.00	1	12/30/2016 01:00	WG938049
Thallium	U		0.190	2.00	1	12/30/2016 01:00	WG938049
Tin	0.331	B J	0.300	2.00	1	12/30/2016 01:00	WG938049
Vanadium	U		0.180	5.00	1	12/30/2016 01:00	WG938049
Zinc	U		2.56	25.0	1	12/30/2016 01:00	WG938049
Zinc	U		2.56	25.0	1	12/30/2016 10:46	WG939602

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 16:37	WG938676
Acrylonitrile	U		1.87	10.0	1	12/26/2016 16:37	WG938676
Benzene	U		0.331	1.00	1	12/26/2016 16:37	WG938676
Bromochloromethane	U		0.520	1.00	1	12/26/2016 16:37	WG938676
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 16:37	WG938676
Bromoform	U		0.469	1.00	1	12/26/2016 16:37	WG938676
Bromomethane	U		0.866	5.00	1	12/26/2016 16:37	WG938676
Carbon disulfide	U		0.275	1.00	1	12/26/2016 16:37	WG938676
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 16:37	WG938676
Chlorobenzene	U		0.348	1.00	1	12/26/2016 16:37	WG938676
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 16:37	WG938676
Chloroethane	U		0.453	5.00	1	12/26/2016 16:37	WG938676
Chloroform	U		0.324	5.00	1	12/26/2016 16:37	WG938676
Chloromethane	U		0.276	2.50	1	12/26/2016 16:37	WG938676
Dibromomethane	U		0.346	1.00	1	12/26/2016 16:37	WG938676
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 16:37	WG938676
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 16:37	WG938676
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 16:37	WG938676
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 16:37	WG938676
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 16:37	WG938676
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 16:37	WG938676
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 16:37	WG938676
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 16:37	WG938676
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 16:37	WG938676
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 16:37	WG938676
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 16:37	WG938676
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 16:37	WG938676
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 16:37	WG938676
Ethylbenzene	U		0.384	1.00	1	12/26/2016 16:37	WG938676
2-Hexanone	U		3.82	10.0	1	12/26/2016 16:37	WG938676
Iodomethane	U		1.71	10.0	1	12/26/2016 16:37	WG938676
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 16:37	WG938676
Methylene Chloride	U		1.00	5.00	1	12/26/2016 16:37	WG938676
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 16:37	WG938676
Styrene	U		0.307	1.00	1	12/26/2016 16:37	WG938676
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 16:37	WG938676
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 16:37	WG938676
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 16:37	WG938676
Toluene	0.865	J	0.412	1.00	1	12/26/2016 16:37	WG938676
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 16:37	WG938676
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 16:37	WG938676
Trichloroethene	U		0.398	1.00	1	12/26/2016 16:37	WG938676
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 16:37	WG938676
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 16:37	WG938676
Vinyl acetate	U		1.63	10.0	1	12/26/2016 16:37	WG938676
Vinyl chloride	U		0.259	1.00	1	12/26/2016 16:37	WG938676
Xylenes, Total	U		1.06	3.00	1	12/26/2016 16:37	WG938676
(S) Toluene-d8	*103*			*90.0-115*		*12/26/2016 16:37*	WG938676
(S) Dibromofluoromethane	*104*			*79.0-121*		*12/26/2016 16:37*	WG938676
(S) 4-Bromofluorobenzene	*115*			*80.1-120*		*12/26/2016 16:37*	WG938676

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/26/2016 16:50	WG938676
Acrylonitrile	U		1.87	10.0	1	12/26/2016 16:50	WG938676
Benzene	U		0.331	1.00	1	12/26/2016 16:50	WG938676
Bromochloromethane	U		0.520	1.00	1	12/26/2016 16:50	WG938676
Bromodichloromethane	U		0.380	1.00	1	12/26/2016 16:50	WG938676
Bromoform	U		0.469	1.00	1	12/26/2016 16:50	WG938676
Bromomethane	U		0.866	5.00	1	12/26/2016 16:50	WG938676
Carbon disulfide	U		0.275	1.00	1	12/26/2016 16:50	WG938676
Carbon tetrachloride	U		0.379	1.00	1	12/26/2016 16:50	WG938676
Chlorobenzene	U		0.348	1.00	1	12/26/2016 16:50	WG938676
Chlorodibromomethane	U		0.327	1.00	1	12/26/2016 16:50	WG938676
Chloroethane	U		0.453	5.00	1	12/26/2016 16:50	WG938676
Chloroform	U		0.324	5.00	1	12/26/2016 16:50	WG938676
Chloromethane	U		0.276	2.50	1	12/26/2016 16:50	WG938676
Dibromomethane	U		0.346	1.00	1	12/26/2016 16:50	WG938676
1,2-Dibromoethane	U		0.381	1.00	1	12/26/2016 16:50	WG938676
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/26/2016 16:50	WG938676
1,2-Dichlorobenzene	U		0.349	1.00	1	12/26/2016 16:50	WG938676
1,4-Dichlorobenzene	U		0.274	1.00	1	12/26/2016 16:50	WG938676
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/26/2016 16:50	WG938676
1,1-Dichloroethane	U		0.259	1.00	1	12/26/2016 16:50	WG938676
1,2-Dichloroethane	U		0.361	1.00	1	12/26/2016 16:50	WG938676
1,1-Dichloroethene	U		0.398	1.00	1	12/26/2016 16:50	WG938676
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/26/2016 16:50	WG938676
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/26/2016 16:50	WG938676
1,2-Dichloropropane	U		0.306	1.00	1	12/26/2016 16:50	WG938676
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/26/2016 16:50	WG938676
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/26/2016 16:50	WG938676
Ethylbenzene	U		0.384	1.00	1	12/26/2016 16:50	WG938676
2-Hexanone	U		3.82	10.0	1	12/26/2016 16:50	WG938676
Iodomethane	U		1.71	10.0	1	12/26/2016 16:50	WG938676
2-Butanone (MEK)	U		3.93	15.0	1	12/26/2016 16:50	WG938676
Methylene Chloride	U		1.00	5.00	1	12/26/2016 16:50	WG938676
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/26/2016 16:50	WG938676
Styrene	U		0.307	1.00	1	12/26/2016 16:50	WG938676
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/26/2016 16:50	WG938676
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/26/2016 16:50	WG938676
Tetrachloroethene	U		0.372	1.00	1	12/26/2016 16:50	WG938676
Toluene	0.820	J	0.412	1.00	1	12/26/2016 16:50	WG938676
1,1,1-Trichloroethane	U		0.319	1.00	1	12/26/2016 16:50	WG938676
1,1,2-Trichloroethane	U		0.383	1.00	1	12/26/2016 16:50	WG938676
Trichloroethene	U		0.398	1.00	1	12/26/2016 16:50	WG938676
Trichlorofluoromethane	U		1.20	5.00	1	12/26/2016 16:50	WG938676
1,2,3-Trichloropropane	U		0.807	2.50	1	12/26/2016 16:50	WG938676
Vinyl acetate	U		1.63	10.0	1	12/26/2016 16:50	WG938676
Vinyl chloride	U		0.259	1.00	1	12/26/2016 16:50	WG938676
Xylenes, Total	U		1.06	3.00	1	12/26/2016 16:50	WG938676
(S) Toluene-d8	*104*			*90.0-115*		*12/26/2016 16:50*	WG938676
(S) Dibromofluoromethane	*105*			*79.0-121*		*12/26/2016 16:50*	WG938676
(S) 4-Bromofluorobenzene	*114*			*80.1-120*		*12/26/2016 16:50*	WG938676

Volatile Organic Compounds (GC/MS) by Method 8260B

Analyte	Result	Qualifier	MDL	RDL	Dilution	Analysis	Batch
	ug/l		ug/l	ug/l		date / time	
Acetone	U		10.0	50.0	1	12/25/2016 19:07	WG938588
Acrylonitrile	U		1.87	10.0	1	12/25/2016 19:07	WG938588
Benzene	U		0.331	1.00	1	12/25/2016 19:07	WG938588
Bromochloromethane	U		0.520	1.00	1	12/25/2016 19:07	WG938588
Bromodichloromethane	U		0.380	1.00	1	12/25/2016 19:07	WG938588
Bromoform	U		0.469	1.00	1	12/25/2016 19:07	WG938588
Bromomethane	U		0.866	5.00	1	12/25/2016 19:07	WG938588
Carbon disulfide	U		0.275	1.00	1	12/25/2016 19:07	WG938588
Carbon tetrachloride	U		0.379	1.00	1	12/25/2016 19:07	WG938588
Chlorobenzene	U		0.348	1.00	1	12/25/2016 19:07	WG938588
Chlorodibromomethane	U		0.327	1.00	1	12/25/2016 19:07	WG938588
Chloroethane	U		0.453	5.00	1	12/25/2016 19:07	WG938588
Chloroform	U		0.324	5.00	1	12/25/2016 19:07	WG938588
Chloromethane	U		0.276	2.50	1	12/25/2016 19:07	WG938588
Dibromomethane	U		0.346	1.00	1	12/25/2016 19:07	WG938588
1,2-Dibromoethane	U		0.381	1.00	1	12/25/2016 19:07	WG938588
1,2-Dibromo-3-Chloropropane	U		1.33	5.00	1	12/25/2016 19:07	WG938588
1,2-Dichlorobenzene	U		0.349	1.00	1	12/25/2016 19:07	WG938588
1,4-Dichlorobenzene	U		0.274	1.00	1	12/25/2016 19:07	WG938588
trans-1,4-Dichloro-2-butene	U		0.866	2.50	1	12/25/2016 19:07	WG938588
1,1-Dichloroethane	U		0.259	1.00	1	12/25/2016 19:07	WG938588
1,2-Dichloroethane	U		0.361	1.00	1	12/25/2016 19:07	WG938588
1,1-Dichloroethene	U		0.398	1.00	1	12/25/2016 19:07	WG938588
cis-1,2-Dichloroethene	U		0.260	1.00	1	12/25/2016 19:07	WG938588
trans-1,2-Dichloroethene	U		0.396	1.00	1	12/25/2016 19:07	WG938588
1,2-Dichloropropane	U		0.306	1.00	1	12/25/2016 19:07	WG938588
cis-1,3-Dichloropropene	U		0.418	1.00	1	12/25/2016 19:07	WG938588
trans-1,3-Dichloropropene	U		0.419	1.00	1	12/25/2016 19:07	WG938588
Ethylbenzene	U		0.384	1.00	1	12/25/2016 19:07	WG938588
2-Hexanone	U		3.82	10.0	1	12/25/2016 19:07	WG938588
Iodomethane	U		1.71	10.0	1	12/25/2016 19:07	WG938588
2-Butanone (MEK)	U		3.93	15.0	1	12/25/2016 19:07	WG938588
Methylene Chloride	U		1.00	5.00	1	12/25/2016 19:07	WG938588
4-Methyl-2-pentanone (MIBK)	U		2.14	10.0	1	12/25/2016 19:07	WG938588
Styrene	U		0.307	1.00	1	12/25/2016 19:07	WG938588
1,1,1,2-Tetrachloroethane	U		0.385	1.00	1	12/25/2016 19:07	WG938588
1,1,2,2-Tetrachloroethane	U		0.130	1.00	1	12/25/2016 19:07	WG938588
Tetrachloroethene	U		0.372	1.00	1	12/25/2016 19:07	WG938588
Toluene	U		0.412	1.00	1	12/25/2016 19:07	WG938588
1,1,1-Trichloroethane	U		0.319	1.00	1	12/25/2016 19:07	WG938588
1,1,2-Trichloroethane	U		0.383	1.00	1	12/25/2016 19:07	WG938588
Trichloroethene	U		0.398	1.00	1	12/25/2016 19:07	WG938588
Trichlorofluoromethane	U		1.20	5.00	1	12/25/2016 19:07	WG938588
1,2,3-Trichloropropane	U		0.807	2.50	1	12/25/2016 19:07	WG938588
Vinyl acetate	U		1.63	10.0	1	12/25/2016 19:07	WG938588
Vinyl chloride	U		0.259	1.00	1	12/25/2016 19:07	WG938588
Xylenes, Total	U		1.06	3.00	1	12/25/2016 19:07	WG938588
(S) Toluene-d8	*98.7*			*90.0-115*		*12/25/2016 19:07*	WG938588
(S) Dibromofluoromethane	*98.3*			*79.0-121*		*12/25/2016 19:07*	WG938588
(S) 4-Bromofluorobenzene	*99.5*			*80.1-120*		*12/25/2016 19:07*	WG938588

Abbreviations and Definitions

SDG	Sample Delivery Group.
MDL	Method Detection Limit.
RDL	Reported Detection Limit.
U	Not detected at the Reporting Limit (or MDL where applicable).
(S)	Surrogate (Surrogate Standard) - Analytes added to every blank, sample, Laboratory Control Sample/Duplicate and Matrix Spike/Duplicate; used to evaluate analytical efficiency by measuring recovery. Surrogates are not expected to be detected in all environmental media.

Qualifier	Description
B	The same analyte is found in the associated blank.
J	The identification of the analyte is acceptable; the reported value is an estimate.
J3	The associated batch QC was outside the established quality control range for precision.
J4	The associated batch QC was outside the established quality control range for accuracy.
J5	The sample matrix interfered with the ability to make any accurate determination; spike value is high.
O1	The analyte failed the method required serial dilution test and/or subsequent post-spike criteria. These failures indicate matrix interference.
P1	RPD value not applicable for sample concentrations less than 5 times the reporting limit.

ESC Lab Sciences is the only environmental laboratory accredited/certified to support your work nationwide from one location. One phone call, one point of contact, one laboratory. No other lab is as accessible or prepared to handle your needs throughout the country. Our capacity and capability from our single location laboratory is comparable to the collective totals of the network laboratories in our industry. The most significant benefit to our "one location" design is the design of our laboratory campus. The model is conducive to accelerated productivity, decreasing turn-around time, and preventing cross contamination, thus protecting sample integrity. Our focus on premium quality and prompt service allows us to be **YOUR LAB OF CHOICE.**

***** Not all certifications held by the laboratory are applicable to the results reported in the attached report.

State Accreditations

Alabama	40660	Nevada	TN-03-2002-34
Alaska	UST-080	New Hampshire	2975
Arizona	AZ0612	New Jersey–NELAP	TN002
Arkansas	88-0469	New Mexico	TN00003
California	01157CA	New York	11742
Colorado	TN00003	North Carolina	Env375
Conneticut	PH-0197	North Carolina [1]	DW21704
Florida	E87487	North Carolina [2]	41
Georgia	NELAP	North Dakota	R-140
Georgia [1]	923	Ohio–VAP	CL0069
Idaho	TN00003	Oklahoma	9915
Illinois	200008	Oregon	TN200002
Indiana	C-TN-01	Pennsylvania	68-02979
Iowa	364	Rhode Island	221
Kansas	E-10277	South Carolina	84004
Kentucky [1]	90010	South Dakota	n/a
Kentucky [2]	16	Tennessee [1 4]	2006
Louisiana	AI30792	Texas	T 104704245-07-TX
Maine	TN0002	Texas [5]	LAB0152
Maryland	324	Utah	6157585858
Massachusetts	M-TN003	Vermont	VT2006
Michigan	9958	Virginia	109
Minnesota	047-999-395	Washington	C1915
Mississippi	TN00003	West Virginia	233
Missouri	340	Wisconsin	9980939910
Montana	CERT0086	Wyoming	A2LA
Nebraska	NE-OS-15-05		

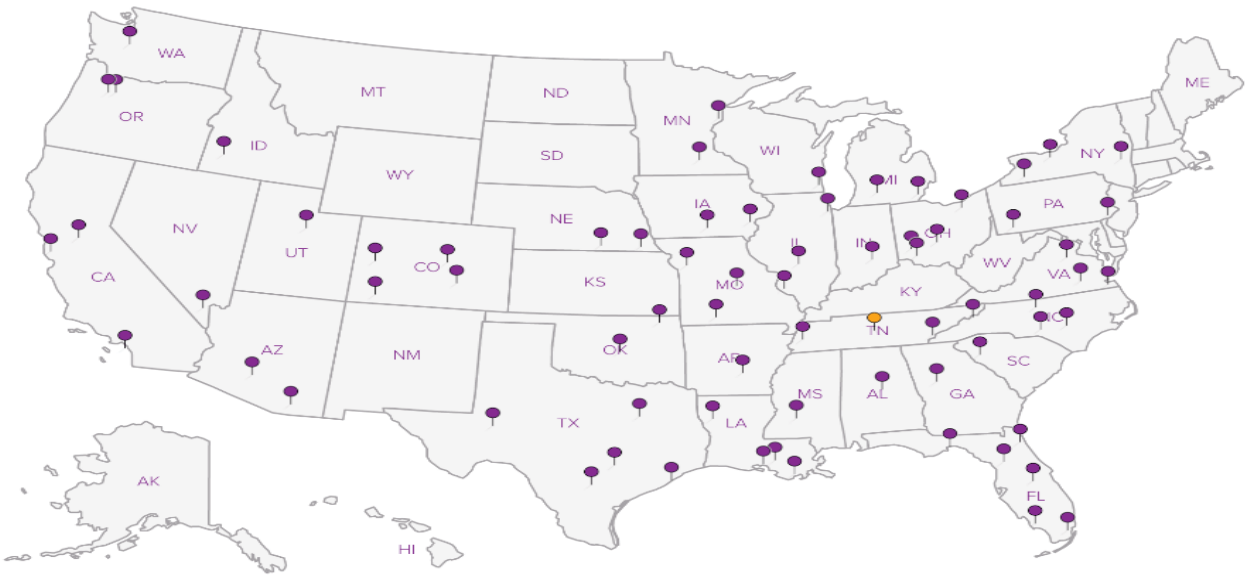
Third Party & Federal Accreditations

A2LA – ISO 17025	1461.01	AIHA	100789
A2LA – ISO 17025 [5]	1461.02	DOD	1461.01
Canada	1461.01	USDA	S-67674
EPA–Crypto	TN00003		

[1] Drinking Water [2] Underground Storage Tanks [3] Aquatic Toxicity [4] Chemical/Microbiological [5] Mold [n/a] Accreditation not applicable

Our Locations

ESC Lab Sciences has sixty-four client support centers that provide sample pickup and/or the delivery of sampling supplies. If you would like assistance from one of our support offices, please contact our main office. **ESC Lab Sciences performs all testing at our central laboratory.**



SCS Engineers - Little Rock, AR

11219 Richardson Drive
North Little Rock, AR 72113

Billing Information & Quote Number:
Accounts Payable
11219 Richardson Drive
N. Little Rock, AR 72113

Report to: **Stacie Whitmer**

Email To: **swhitmer@scsengineers.com**

Project Description: **Nabors Landfill**

City/State Collected: Mountain Home AR

Phone: **501-812-4551**
Fax:

Client Project #

Lab Project # **CHIROCKAR-NABORS**

Collected by (print): Darren Mothey

Site/Facility ID #

P.O. #

Collected by (signature):

Immediately Packed on Ice N ___ Y X

Rush? (Lab MUST Be Notified)
___ Same Day 200%
___ Next Day 100%
___ Two Day 50%
___ Three Day 25%

Date Results Needed

Email? ___No X_Yes
FAX? ___No ___Yes

Sample ID	Comp/Grab	Matrix *	Depth	Date	Time	No. of Cntrs	CN 250mlHDPEAmb-NaOH ≥12	Cl, SO4, TDS 250mlHDPE-NoPres	M6010AP1 + Fe,Mn,Sn 250mlHDPE-HNO3 ≤2	SULFIDE 125mlAmb-S-NaOH+ZnAc ≥12	TOC 250mlAmb-HCl ≤2	V8260AP1 40mlAmb-HCl	V8260AP1- Trip Blk 40mlAmb-HCl				Rem./Contaminant	Sample # (lab only)
MW-1	G	GW		12-20-16	1558	7	X	X	X	X	X	X						01
MW-2	G	GW		12-20-16	1652	7	X	X	X	X	X	X						02
MW-3	G	GW		12-21-16	804	7	X	X	X	X	X	X						03
MW-4	G	GW		12-21-16	847	7	X	X	X	X	X	X						04
MW-5	G	GW		12-21-16	911	7	X	X	X	X	X	X						05
MW-6	G	GW		12-21-16	946	7	X	X	X	X	X	X						06
MW-7	G	GW		12-20-16	1438	7	X	X	X	X	X	X						07
CAO-1	G	GW		12-21-16	1022	7	X	X	X	X	X	X						08
CAO-2	G	GW		12-21-16	748	7	X	X	X	X	X	X						09
CAO-3	G	GW		12-21-16	720	7	X	X	X	X	X	X						10

* Matrix: **SS** - Soil **GW** - Groundwater **WW** - WasteWater **DW** - Drinking Water **OT** - Other GW

Remarks:

pH ___ Temp ___
Flow ___ Other ___

Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)
Darren Mothey	12-21-16	1300	
Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)
Relinquished by : (Signature)	Date:	Time:	Received for lab by: (Signature)
			Mwt

Samples returned via: ☐ UPS ☑ FedEx ☐ Courier ☐ ___
Temp: 2.6 °C Bottles Received: 22

Date: 12-22-16 Time: 0930

Hold #

Condition: NB50 (lab use only) 02

COC Seal Intact: ___Y ___N ___NA

pH Checked: ≤2, ≥12 NCF:

Chain of Custody Page ___ of ___

ESC LABORATORY SCIENCES
YOUR LAB OF CHOICE
12065 Lebanon Rd
Mount Juliet, TN 37122
Phone: 615-758-5858
Phone: 800-767-5859
Fax: 615-758-5859

L# 870396

Table #

Acctnum: CHIROCKAR
Template: T98870
Prelogin: P577243
TSR: 134 - Mark W. Beasley
PB: JTA 11-23-16

Shipped Via: FedEX Ground

SCS Engineers - Little Rock, AR

11219 Richardson Drive
North Little Rock, AR 72113

Billing Information & Quote Number:
Accounts Payable
11219 Richardson Drive
N. Little Rock, AR 72113

Report to: **Stacie Whitmer**

Email To: swhitmer@scsengineers.com

Project Description: **Nabors Landfill**

City/State Collected: Mtn Home, AR

Phone: **501-812-4551**
Fax:

Client Project #

Lab Project # **CHIROCKAR-NABORS**

Collected by (print): Darren Mottley

Site/Facility ID #

P.O. #

Collected by (signature): Darren Mottley

Immediately Packed on Ice N ___ Y X

Rush? (Lab MUST Be Notified)
___ Same Day 200%
___ Next Day 100%
___ Two Day 50%
___ Three Day 25%

Date Results Needed

Email? ___No X__Yes
FAX? ___No ___Yes

Analysis / Container / Preservative

Sample ID	Comp/Grab	Matrix *	Depth	Date	Time	No. of Cntrs	CN 250mlHDPEAmb-NaOH +R	Cl, SO4, TDS 250mlHDPE-NoPres	M6010AP1 + Fe,Mn,Sn 250mlHDPE-HNO3 +R	SULFIDE 125mlAmb-S-NaOH+ZnAc +R	TOC 250mlAmb-HCl +R	V8260AP1 40mlAmb-HCl	V8260AP1- Trip Blk 40mlAmb-HCl				Rem./Contaminant	Sample # (lab only)
NAB-1		GW				7	X	X	X	X	X	X						
NAB-2	G	GW		12-20-16	1253	7	X	X	X	X	X	X						11
NAB-3	G	GW		12-20-16	819	7	X	X	X	X	X	X						12
NAB-4	G	GW		12-20-16	1042	7	X	X	X	X	X	X						13
NAB-7	G	GW		12-20-16	1515	7	X	X	X	X	X	X						14
NAB-8	G	GW		12-20-16	708	7	X	X	X	X	X	X						15
MW-509D	G	GW		12-20-16	1415	7	X	X	X	X	X	X						16
MW-577	G	GW		12-20-16	1210	7	X	X	X	X	X	X						17
MW-689D	G	GW		12-20-16	931	7	X	X	X	X	X	X						18
MW-633D	G	GW		12-20-16	1130	7	X	X	X	X	X	X						19

* Matrix: **SS** - Soil **GW** - Groundwater **WW** - WasteWater **DW** - Drinking Water **OT** - Other GW

Remarks:

pH _____ Temp _____

Flow _____ Other _____

70666 8129 16160

Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)
Darren Mottley	12-21-16	1300	
Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)
Relinquished by : (Signature)	Date:	Time:	Received for lab by: (Signature)
			MWL

Samples returned via: ☐ UPS ☑ FedEx ☐ Courier ☐ _____

Temp: 2-6 °C Bottles Received: 22

Date: 12-22-16 Time: 0930

Hold #

Condition: NB50 (lab use only) A

COC Seal Intact: ___Y ___N ___NA

pH Checked: 12, 712 NCF:

Chain of Custody Page ___ of ___

ESC LABORATORY SCIENCES
YOUR LAB OF CHOICE
12065 Lebanon Rd
Mount Juliet, TN 37122
Phone: 615-758-5858
Phone: 800-767-5859
Fax: 615-758-5859

L# 880396
1038

Acctnum: CHIROCKAR
Template: T98870
Prelogin: P577243
TSR: 134 - Mark W. Beasley
PB: JH 11-23-16
Shipped Via: FedEx Ground

SCS Engineers - Little Rock, AR

11219 Richardson Drive
North Little Rock, AR 72113

Billing Information & Quote Number:

Accounts Payable
11219 Richardson Drive
N. Little Rock, AR 72113

Report to: Stacie Whitmer

Email To: swhitmer@scsengineers.com

Project Description: Nabors Landfill

City/State Collected: Mt. Home AR

Phone: 501-812-4551
Fax:

Client Project #

Lab Project # CHIROCKAR-NABORS

Collected by (print): Darren Mottley

Site/Facility ID #

P.O. #

Collected by (signature):

Immediately Packed on Ice N ___ Y X

Rush? (Lab MUST Be Notified)
___ Same Day 200%
___ Next Day 100%
___ Two Day 50%
___ Three Day 25%

Date Results Needed

Email? ___No X_Yes
FAX? ___No ___Yes

Analysis / Container / Preservative

Sample ID	Comp/Grab	Matrix *	Depth	Date	Time	No. of Cntrs	CN 250mlHDPEAmb-NaOH (2,12)	Cl, SO4, TDS 250mlHDPE-NoPres	M6010AP1 + Fe,Mn,Sn 250mlHDPE-HNO3 (2)	SULFIDE 125mlAmb-S-NaOH+ZnAc (2,12)	TOC 250mlAmb-HCl (2)	V8260AP1 40mlAmb-HCl	V8260AP1- Trip Blk 40mlAmb-HCl				Rem./Contaminant	Sample # (lab only)
NE-2	G	GW		12-19-16	1150	7	X	X	X	X	X	X						20
NE-3	G	GW		12-20-16	730	7	X	X	X	X	X	X						21
NE-6	G	GW		12-21-16	820	7	X	X	X	X	X	X						22
MW-1R	G	GW		12-20-16	1629	7	X	X	X	X	X	X						23
TSP-1		GW				7	X	X	X	X	X	X						
TSP-2		GW				7	X	X	X	X	X	X						
TSP-3		GW				7	X	X	X	X	X	X						
TSP-4		GW				7	X	X	X	X	X	X						
SP-4		GW				7	X	X	X	X	X	X						
SP-5		GW				7	X	X	X	X	X	X						

* Matrix: SS - Soil GW - Groundwater WW - WasteWater DW - Drinking Water OT - Other GW

Remarks:

pH ___ Temp ___

Flow ___ Other ___

Chain of Custody Page ___ of ___

ESC LABORATORIES
YOUR LAB OF CHOICE
12065 Lebanon Rd
Mount Juliet, TN 37122
Phone: 615-758-5858
Phone: 800-767-5859
Fax: 615-758-5859

L# 880396

Table #

Acctnum: CHIROCKAR
Template: T98870
Prelogin: P577243
TSR: 134 - Mark W. Beasley
PB: JT 11-23-16
Shipped Via: FedEX Ground

Hold #

Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)
Darren Mottley	12-21-16	1300	
Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)
Relinquished by : (Signature)	Date:	Time:	Received for lab by: (Signature)
			Mwk

Samples returned via: ☐ UPS ☑ FedEx ☐ Courier ☐ ____

Temp: 2.6 °C Bottles Received: 22

Date: 12-22-16 Time: 0930

Condition: NBS0 (lab use only)

COC Seal Intact: ___Y ___N ___NA

pH Checked: C2,712 NCF:

SCS Engineers - Little Rock, AR

11219 Richardson Drive
North Little Rock, AR 72113

Billing Information & Quote Number:
Accounts Payable
11219 Richardson Drive
N. Little Rock, AR 72113

Report to: Stacie Whitmer

Email To: swhitmer@scsengineers.com

Project Description: Nabors Landfill

City/State Collected: MT. Home AR

Phone: 501-812-4551
Fax:

Client Project #

Lab Project # CHIROCKAR-NABORS

Collected by (print): Darren Mattey

Site/Facility ID #

P.O. #

Collected by (signature): Darren Mattey

Immediately Packed on Ice N ___ Y X

Rush? (Lab MUST Be Notified)
___ Same Day 200%
___ Next Day 100%
___ Two Day 50%
___ Three Day 25%

Date Results Needed

Email? ___No X_Yes
FAX? ___No ___Yes

Analysis / Container / Preservative

Sample ID	Comp/Grab	Matrix *	Depth	Date	Time	No. of Cntrs	CN 250mlHDPEAmb-NaOH (12)	Cl, SO4, TDS 250mlHDPE-NoPres	M6010AP1 + Fe,Mn,Sn 250mlHDPE-HNO3 (12)	SULFIDE 125mlAmb-S-NaOH+ZnAc (12)	TOC 250mlAmb-HCl (12)	V8260AP1 40mlAmb-HCl	V8260AP1- Trip Blk 40mlAmb-HCl	Rem./Contaminant	Sample # (lab only)
SP-7	G	GW		12-21-16	1130	7	X	X	X	X	X	X			24
SPRING-A	G	GW		12-21-16	1140	7	X	X	X	X	X	X			25
SPRING-B		GW				7	X	X	X	X	X	X			
SP-NE-3		GW				7	X	X	X	X	X	X			
LANDFILL ENTRANCE SEEP	G	GW		12-21-16	1120	7	X	X	X	X	X	X			26
CLASS IV DRAW		GW				7	X	X	X	X	X	X			
CLASS I DRAW	G	GW		12-21-16	1105	7	X	X	X	X	X	X			27
LEACHATE	G	GW		12-20-16	1705	7	X	X	X	X	X	X			28
DUPLICATE	G	GW		12-21-16	950	7	X	X	X	X	X	X			29
FIELD BLANK	G	GW		12-21-16	955	7	X	X	X	X	X	X			30

* Matrix: SS - Soil GW - Groundwater WW - WasteWater DW - Drinking Water OT - Other GW

Remarks:

pH ___ Temp ___
Flow ___ Other ___

Chain of Custody Page ___ of ___

ESC LAB SCIENCES
YOUR LAB OF CHOICE
12065 Lebanon Rd
Mount Juliet, TN 37122
Phone: 615-758-5858
Phone: 800-767-5859
Fax: 615-758-5859

L# 820396

Table #

Acctnum: CHIROCKAR
Template: T98870
Prelogin: P577243
TSR: 134 - Mark W. Beasley
PB: JG 11-23-16

Shipped Via: FedEX Ground

Hold #

Condition: NB50 (lab use only) Ga

Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)
Darren Mattey	12-21-16	1300	
Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)
Relinquished by : (Signature)	Date:	Time:	Received for lab by: (Signature)
			Mwat

Samples returned via: ☐ UPS ☑ FedEx ☐ Courier ☐ ___

Temp: 2.6 °C Bottles Received: 221

Date: 12-22-16 Time: 0930

COC Seal Intact: ___Y ___N ___NA

pH Checked: L2,712 NCF:

SCS Engineers - Little Rock, AR

11219 Richardson Drive
North Little Rock, AR 72113

Billing Information & Quote Number:
Accounts Payable
11219 Richardson Drive
N. Little Rock, AR 72113

Report to: Stacie Whitmer

Email To: swhitmer@scsengineers.com

Project Description: Nabors Landfill

City/State Collected: MT. Home AR

Phone: 501-812-4551
Fax:

Client Project #

Lab Project # CHIROCKAR-NABORS

Collected by (print): Denver Motley

Site/Facility ID #

P.O. #

Collected by (signature): Denver Motley

Immediately Packed on ice N ___ Y X

Rush? (Lab MUST Be Notified)
___ Same Day 200%
___ Next Day 100%
___ Two Day 50%
___ Three Day 25%

Date Results Needed

Email? ___No X_Yes
FAX? ___No ___Yes

Chain of Custody Page ___ of ___

ESC LABORATORY SCIENCES
YOUR LAB OF CHOICE
12065 Lebanon Rd
Mount Juliet, TN 37122
Phone: 615-758-5858
Phone: 800-767-5859
Fax: 615-758-5859

L# 880391

Table #

Acctnum: CHIROCKAR
Template: T98870
Prelogin: P577243
TSR: 134 - Mark W. Beasley
PB: JG 11-23-16
Shipped Via: FedEx Ground

Analysis / Container / Preservative

Sample ID	Comp/Grab	Matrix *	Depth	Date	Time	No. of Cntrs	CN 250mlHDPEAmb-NaOH	Cl, SO4, TDS 250mlHDPE-NoPres	M6010AP1 + Fe,Mn,Sn 250mlHDPE-HNO3	SULFIDE 125mlAmb-S-NaOH+ZnAc	TOC 250mlAmb-HCl	V8260AP1 40mlAmb-HCl	V8260AP1- Trip Blk 40mlAmb-HCl				Rem./Contaminant	Sample # (lab only)
		GW				7	X	X	X	X	X	X						
		GW				7	X	X	X	X	X	X						
		GW				7	X	X	X	X	X	X						
EQUIPMENT BLANK	6	GW		12-21-16	957	2						X						31
TRIP BLANK		GW		12-21-16	1000	1							X					32
		GW				7	X	X	X	X	X	X						
		GW				7	X	X	X	X	X	X						
		GW				7	X	X	X	X	X	X						
		GW				7	X	X	X	X	X	X						
		GW				7	X	X	X	X	X	X						

712

* Matrix: SS - Soil GW - Groundwater WW - WasteWater DW - Drinking Water OT - Other ___ MW 12-22-16

Remarks:

pH ___ Temp ___
Flow ___ Other ___

Hold #

Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)
Denver Motley	12-21-16	1300	
Relinquished by : (Signature)	Date:	Time:	Received by: (Signature)
Relinquished by : (Signature)	Date:	Time:	Received for lab by: (Signature)
			MWA

Samples returned via: ☐ UPS ☑ FedEx ☐ Courier ☐ ___
Temp: 2.6 °C Bottles Received: 22
Date: 12-22-16 Time: 0930

Condition: NB59 (lab use only) 02
COC Seal Intact: ✓ Y ___N ___NA
pH Checked: <2, 712 NCF:

ESC
L·A·B S·C·I·E·N·C·E·S
YOUR LAB OF CHOICE

Cooler Receipt Form

Client:	CH2M KAR	SDG#	880396
Cooler Received/Opened On: 12/22/16		Temperature Upon Receipt:	2.6 °c
Received By: Michael Witherspoon			
Signature: MWat			

Receipt Check List	Yes	No	N/A
Were custody seals on outside of cooler and intact?			
Were custody papers properly filled out?			
Did all bottles arrive in good condition?			
Were correct bottles used for the analyses requested?			
Was sufficient amount of sample sent in each bottle?			
Were all applicable sample containers correctly preserved and checked for preservation? (Any not in accepted range noted on COC)			
If applicable, was an observable VOA headspace present?			
Non Conformance Generated. (If yes see attached NCF)			

APPENDIX D

STATISTICAL DATABASE

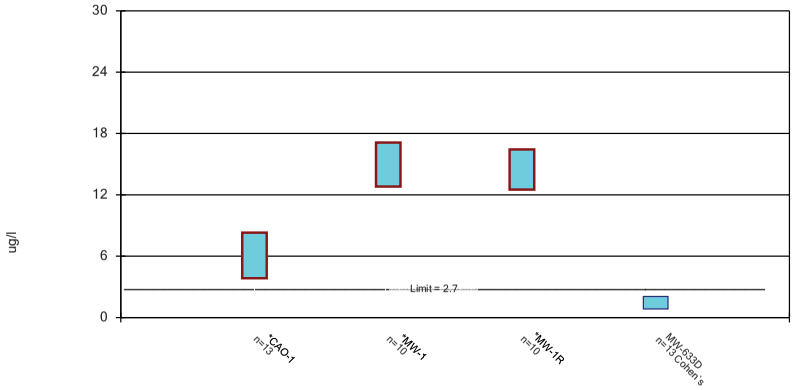
AN ELECTRONIC DATABASE FILE IS INCLUDED WITH THIS REPORT SUBMITTAL

APPENDIX E

STATISTICAL EVALUATION

Parametric Confidence Interval

Compliance limit is exceeded.* Per-well alpha = 0.01. Normality Test: Shapiro Wilk, alpha based on n.

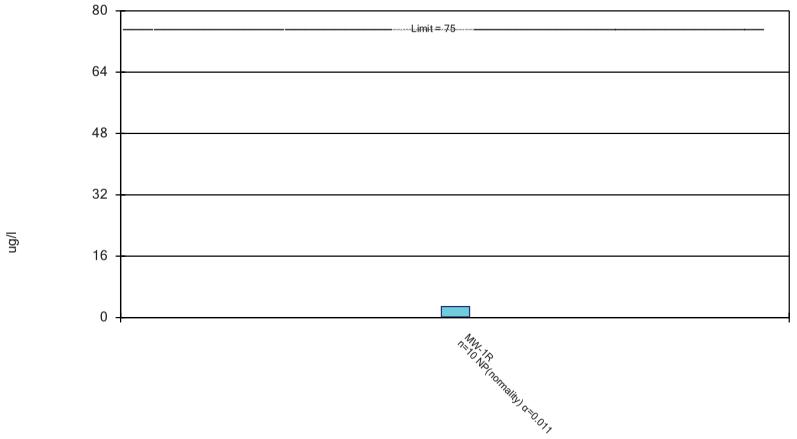


Constituent: 1,1-Dichloroethane Analysis Run 2/28/2017 10:10 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Non-Parametric Confidence Interval

Compliance Limit is not exceeded.

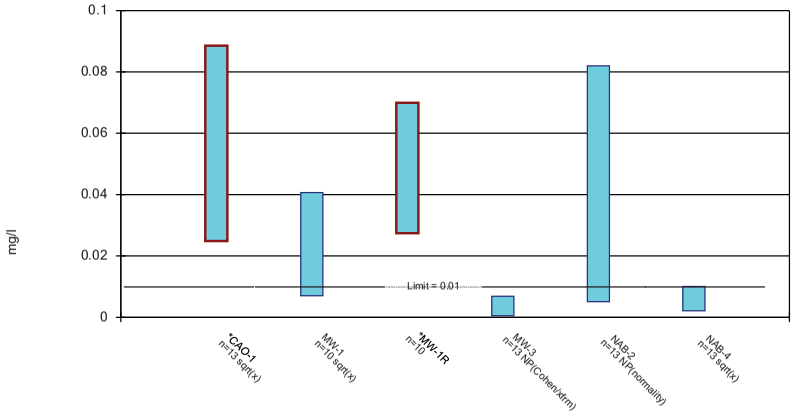


Constituent: 1,4-Dichlorobenzene Analysis Run 2/28/2017 10:10 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Parametric and Non-Parametric (NP) Confidence Interval

Compliance limit is exceeded.* Per-well alpha = 0.01. Normality Test: Shapiro Wilk, alpha based on n.

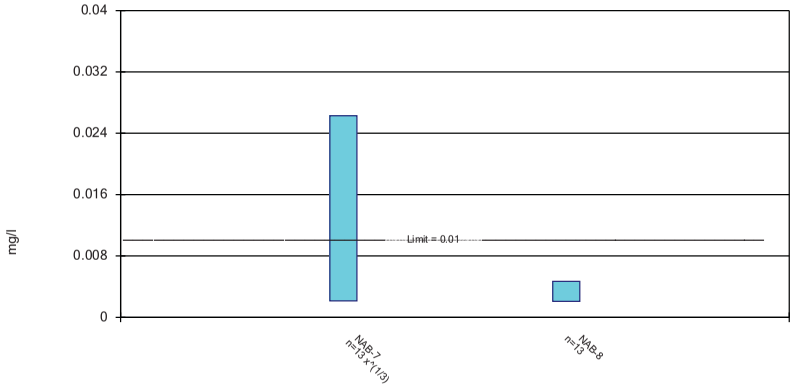


Constituent: Arsenic Analysis Run 2/28/2017 10:10 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Parametric Confidence Interval

Compliance Limit is not exceeded. Per-well alpha = 0.01. Normality Test: Shapiro Wilk, alpha based on n.

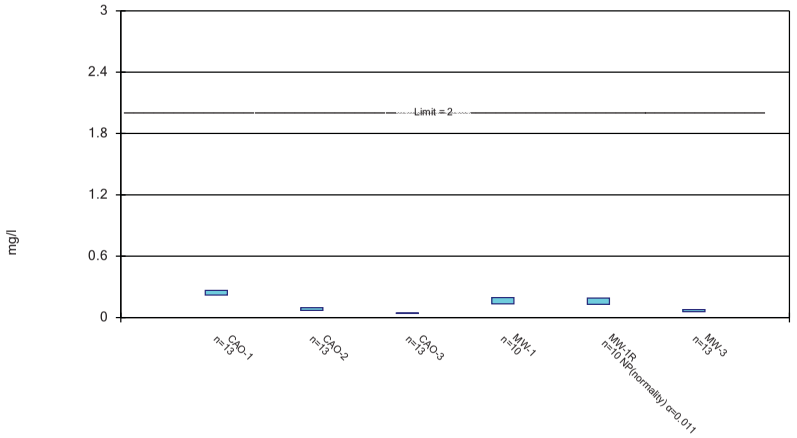


Constituent: Arsenic Analysis Run 2/28/2017 10:10 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Parametric and Non-Parametric (NP) Confidence Interval

Compliance Limit is not exceeded. Per-well alpha = 0.01 except as noted. Normality Test: Shapiro Wilk, alpha based on n.



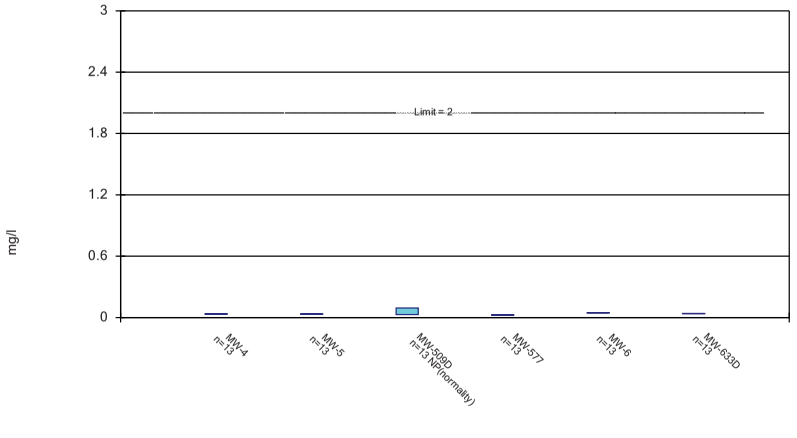
Constituent: Barium Analysis Run 2/28/2017 10:10 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Parametric and Non-Parametric (NP) Confidence Interval

Compliance Limit is not exceeded. Per-well alpha = 0.01. Normality Test: Shapiro Wilk, alpha based on n.



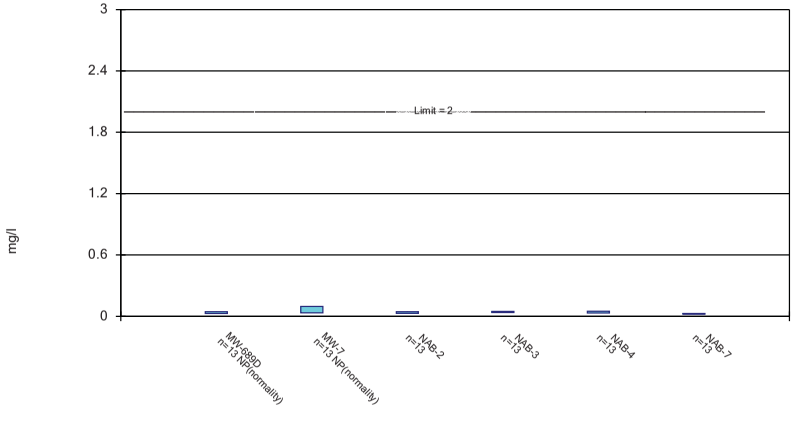
Constituent: Barium Analysis Run 2/28/2017 10:10 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Parametric and Non-Parametric (NP) Confidence Interval

Compliance Limit is not exceeded. Per-well alpha = 0.01. Normality Test: Shapiro Wilk, alpha based on n.



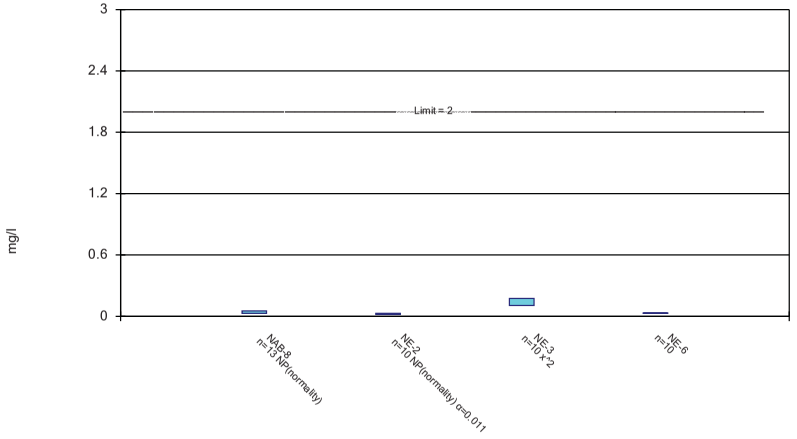
Constituent: Barium Analysis Run 2/28/2017 10:10 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Parametric and Non-Parametric (NP) Confidence Interval

Compliance Limit is not exceeded. Per-well alpha = 0.01 except as noted. Normality Test: Shapiro Wilk, alpha based on n.

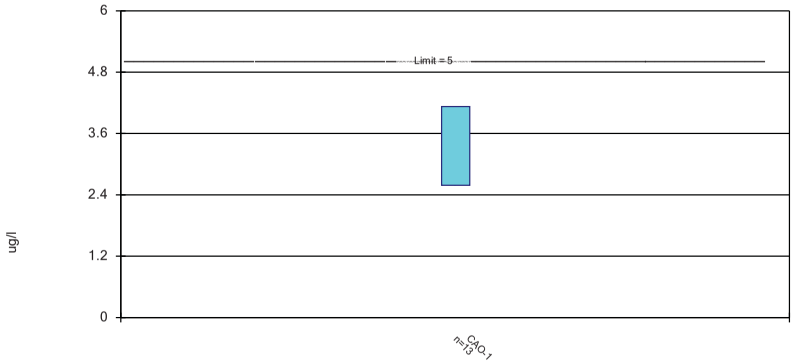


Constituent: Barium Analysis Run 2/28/2017 10:10 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Parametric Confidence Interval

Compliance Limit is not exceeded. Per-well alpha = 0.01. Normality Test: Shapiro Wilk, alpha based on n.

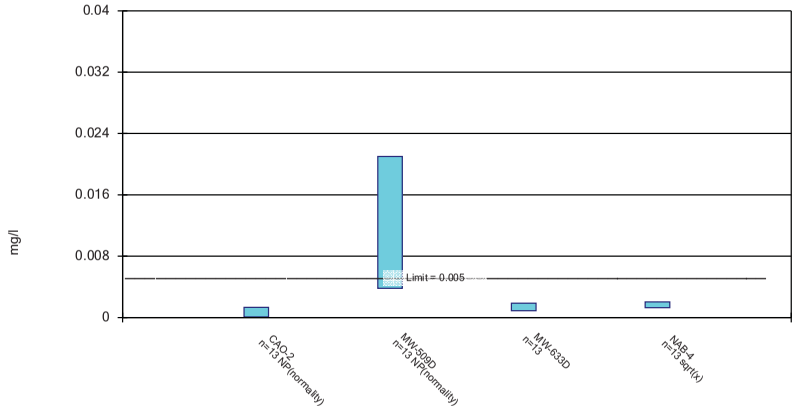


Constituent: Benzene Analysis Run 2/28/2017 10:10 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Parametric and Non-Parametric (NP) Confidence Interval

Compliance Limit is not exceeded. Per-well alpha = 0.01. Normality Test: Shapiro Wilk, alpha based on n.

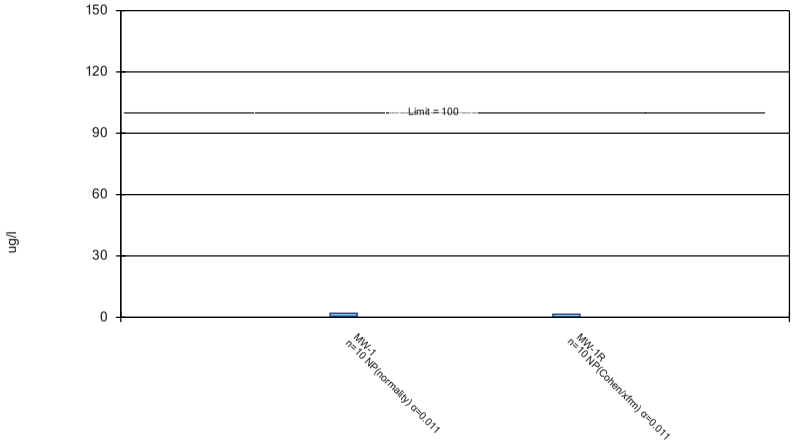


Constituent: Cadmium Analysis Run 2/28/2017 10:10 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Non-Parametric Confidence Interval

Compliance Limit is not exceeded.

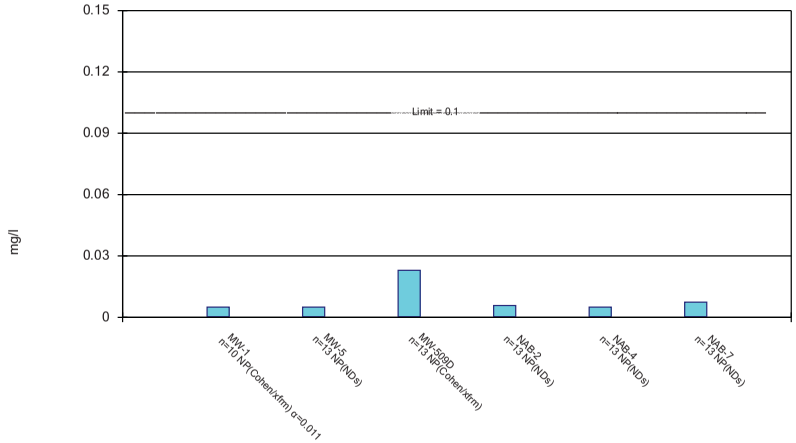


Constituent: Chlorobenzene Analysis Run 2/28/2017 10:10 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Non-Parametric Confidence Interval

Compliance Limit is not exceeded. Per-well alpha = 0.01 except as noted.

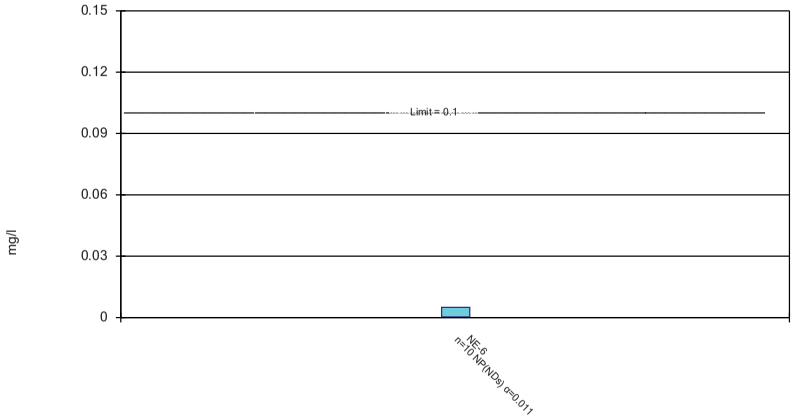


Constituent: Chromium Analysis Run 2/28/2017 10:10 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Non-Parametric Confidence Interval

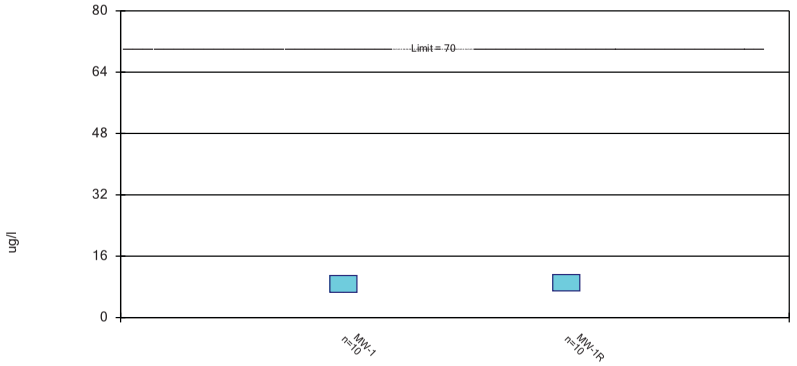
Compliance Limit is not exceeded.



Constituent: Chromium Analysis Run 2/28/2017 10:10 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Parametric Confidence Interval

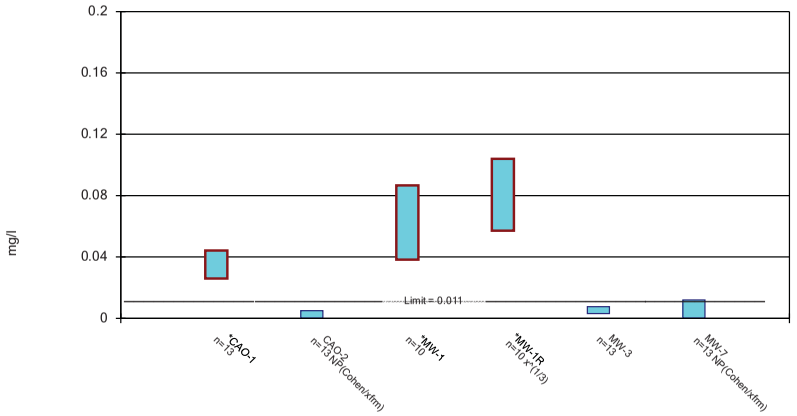
Compliance Limit is not exceeded. Per-well alpha = 0.01. Normality Test: Shapiro Wilk, alpha based on n.



Constituent: cis-1,2-Dichloroethene Analysis Run 2/28/2017 10:10 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Parametric and Non-Parametric (NP) Confidence Interval

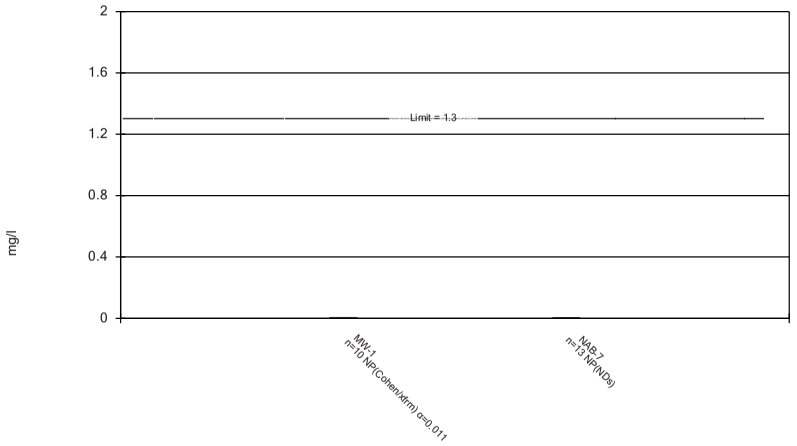
Compliance limit is exceeded.* Per-well alpha = 0.01. Normality Test: Shapiro Wilk, alpha based on n.



Constituent: Cobalt Analysis Run 2/28/2017 10:10 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Non-Parametric Confidence Interval

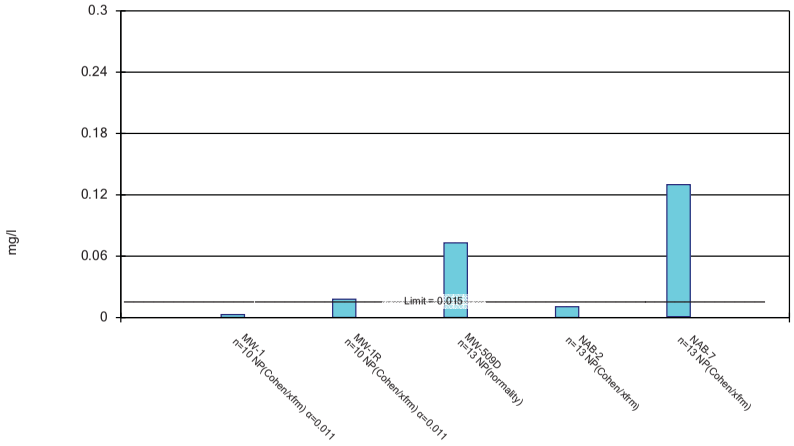
Compliance Limit is not exceeded. Per-well alpha = 0.01 except as noted.



Constituent: Copper Analysis Run 2/28/2017 10:11 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Non-Parametric Confidence Interval

Compliance Limit is not exceeded. Per-well alpha = 0.01 except as noted.



Constituent: Lead Analysis Run 2/28/2017 10:11 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

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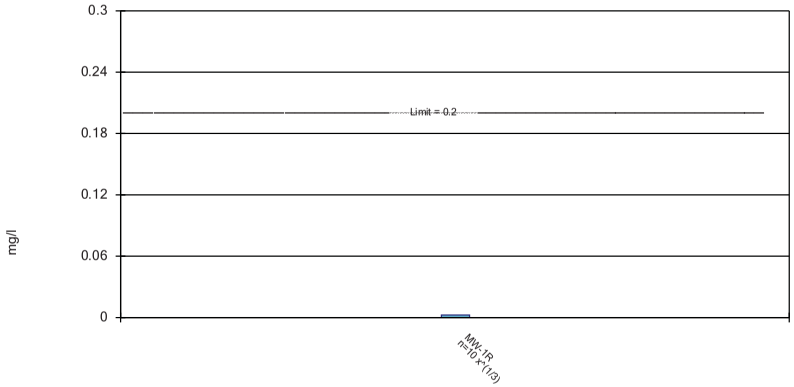
Parametric Confidence Interval

Compliance Limit is not exceeded. Per-well alpha = 0.01. Normality Test: Shapiro Wilk, alpha based on n.



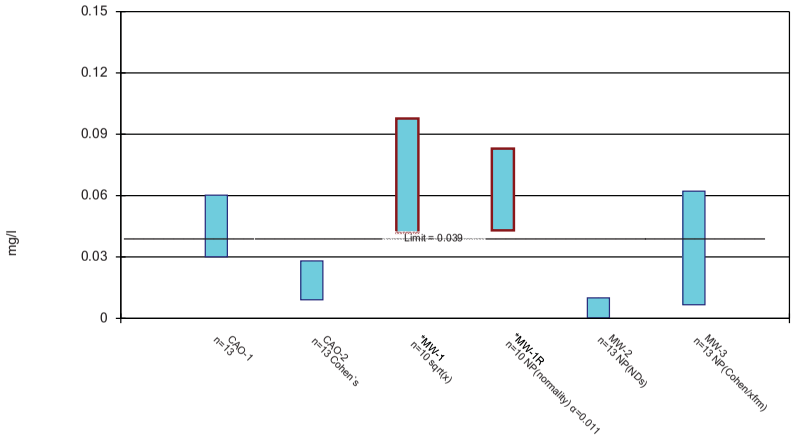
Constituent: Mercury Analysis Run 2/28/2017 10:11 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Parametric and Non-Parametric (NP) Confidence Interval

Compliance limit is exceeded.* Per-well alpha = 0.01 except as noted. Normality Test: Shapiro Wilk, alpha based on n.



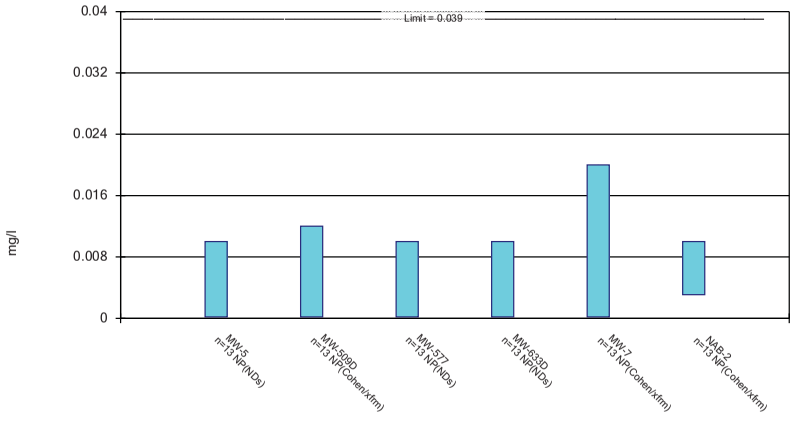
Constituent: Nickel Analysis Run 2/28/2017 10:11 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Non-Parametric Confidence Interval

Compliance Limit is not exceeded. Per-well alpha = 0.01.

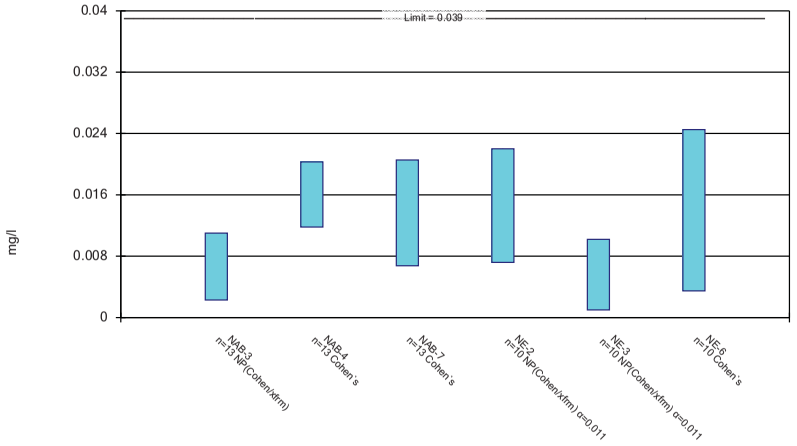


Constituent: Nickel Analysis Run 2/28/2017 10:11 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Parametric and Non-Parametric (NP) Confidence Interval

Compliance Limit is not exceeded. Per-well alpha = 0.01 except as noted. Normality Test: Shapiro Wilk, alpha based on n.



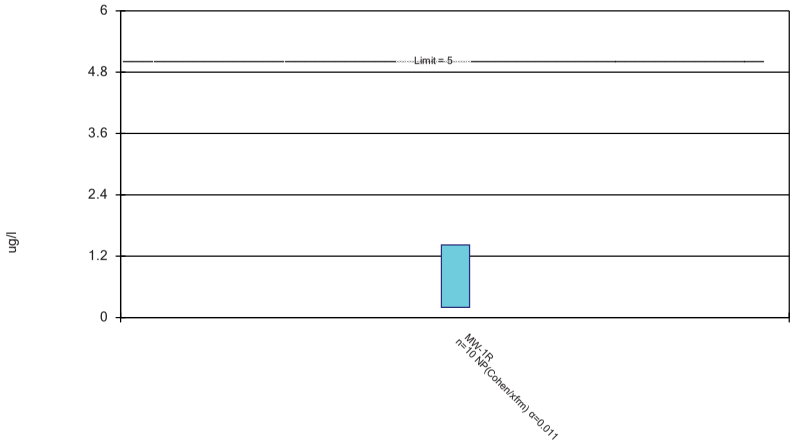
Constituent: Nickel Analysis Run 2/28/2017 10:11 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Non-Parametric Confidence Interval

Compliance Limit is not exceeded.



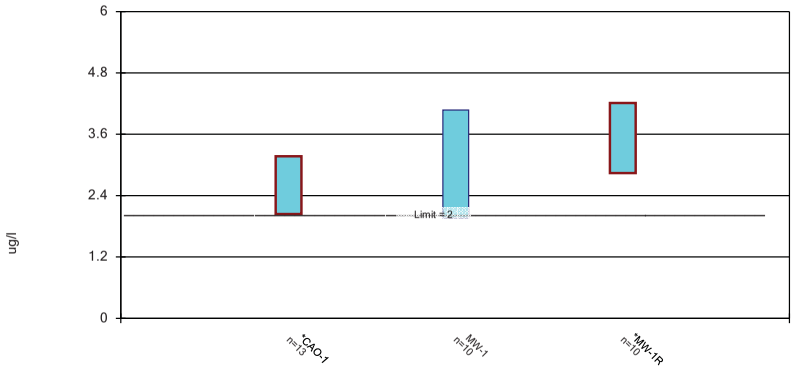
Constituent: Trichloroethene Analysis Run 2/28/2017 10:11 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Parametric Confidence Interval

Compliance limit is exceeded.* Per-well alpha = 0.01. Normality Test: Shapiro Wilk, alpha based on n.



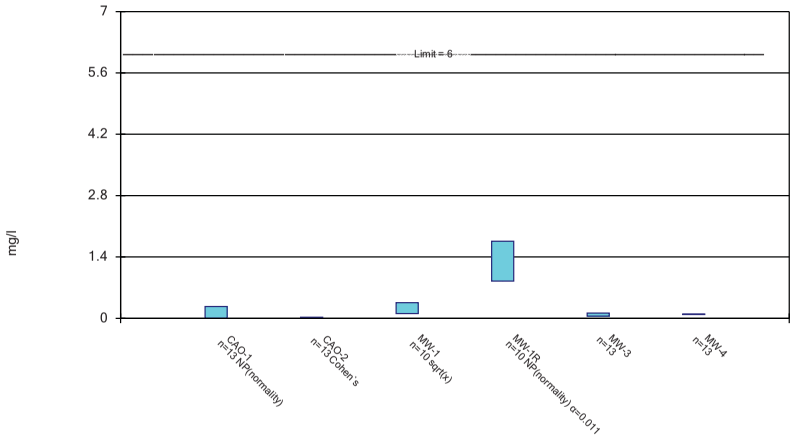
Constituent: Vinyl chloride Analysis Run 2/28/2017 10:11 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Parametric and Non-Parametric (NP) Confidence Interval

Compliance Limit is not exceeded. Per-well alpha = 0.01 except as noted. Normality Test: Shapiro Wilk, alpha based on n.

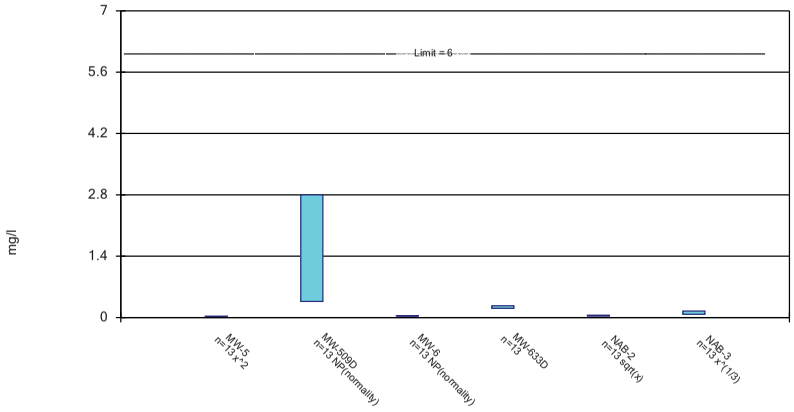


Constituent: Zinc Analysis Run 2/28/2017 10:11 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Parametric and Non-Parametric (NP) Confidence Interval

Compliance Limit is not exceeded. Per-well alpha = 0.01. Normality Test: Shapiro Wilk, alpha based on n.



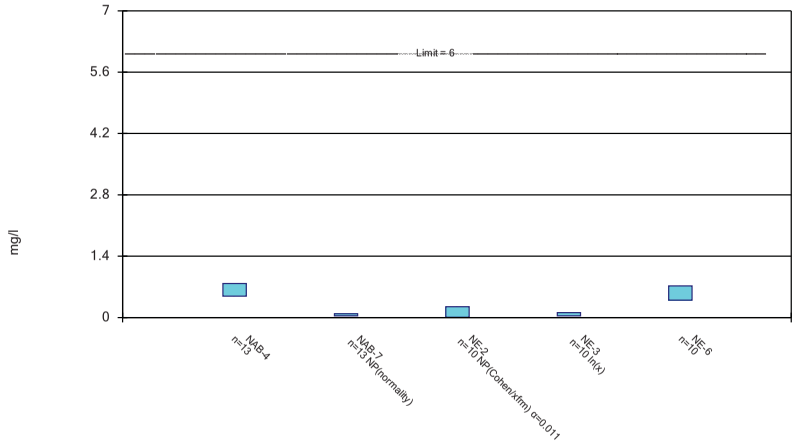
Constituent: Zinc Analysis Run 2/28/2017 10:11 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Parametric and Non-Parametric (NP) Confidence Interval

Compliance Limit is not exceeded. Per-well alpha = 0.01 except as noted. Normality Test: Shapiro Wilk, alpha based on n.

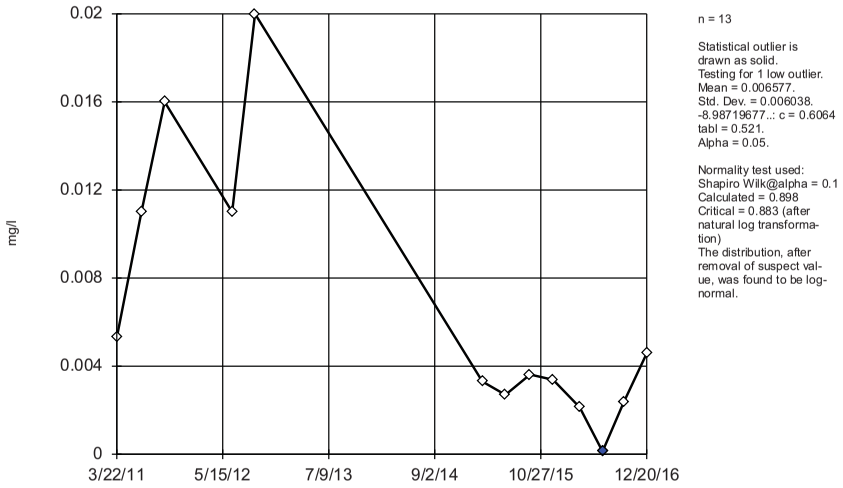


Constituent: Zinc Analysis Run 2/28/2017 10:11 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Dixon's Outlier Test

NAB-4



n = 13

Statistical outlier is drawn as solid.
Testing for 1 low outlier.
Mean = 0.006577.
Std. Dev. = 0.006038.
-8.98719677..: c = 0.6064
tabl = 0.521.
Alpha = 0.05.

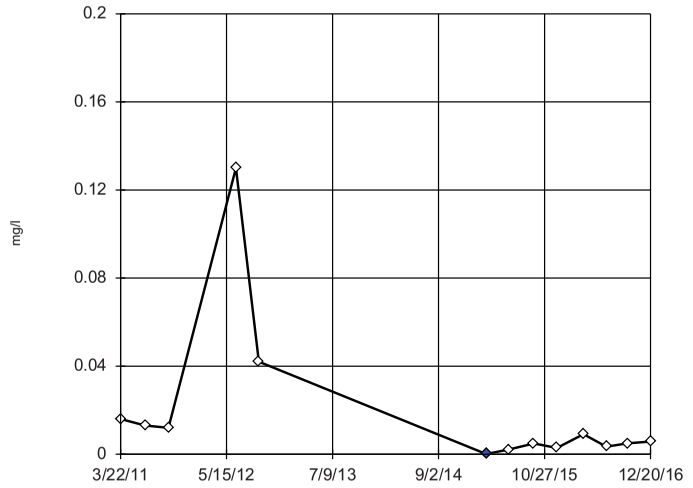
Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.898
Critical = 0.883 (after natural log transformation)
The distribution, after removal of suspect value, was found to be log-normal.

Constituent: Arsenic Analysis Run 2/28/2017 10:00 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

NAB-7



n = 13

Statistical outlier is drawn as solid.
Testing for 1 low outlier.
Mean = 0.01895.
Std. Dev. = 0.03507.
-8.98719677..: c = 0.5363
tabl = 0.521.
Alpha = 0.05.

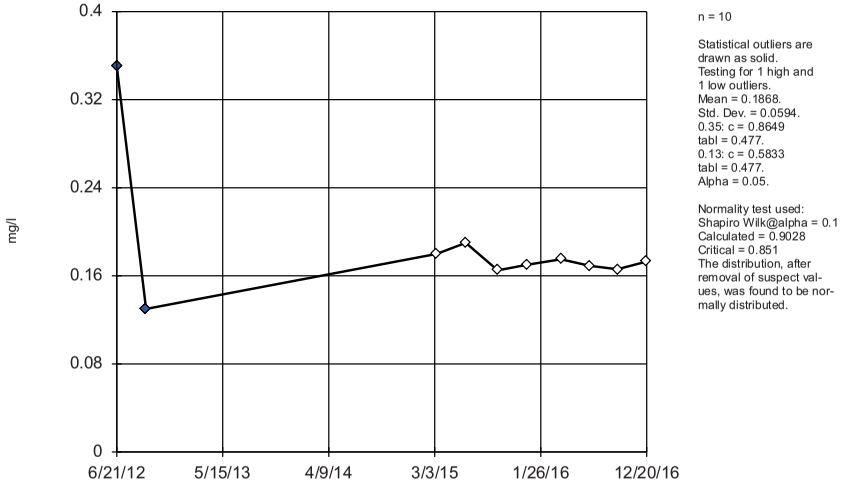
Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9195
Critical = 0.883 (after natural log transformation)
The distribution, after removal of suspect value, was found to be log-normal.

Constituent: Arsenic Analysis Run 2/28/2017 10:00 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

MW-1R



n = 10

Statistical outliers are drawn as solid.
Testing for 1 high and 1 low outliers.
Mean = 0.1868.
Std. Dev. = 0.0594.
0.35: c = 0.8649
tabl = 0.477.
0.13: c = 0.5833
tabl = 0.477.
Alpha = 0.05.

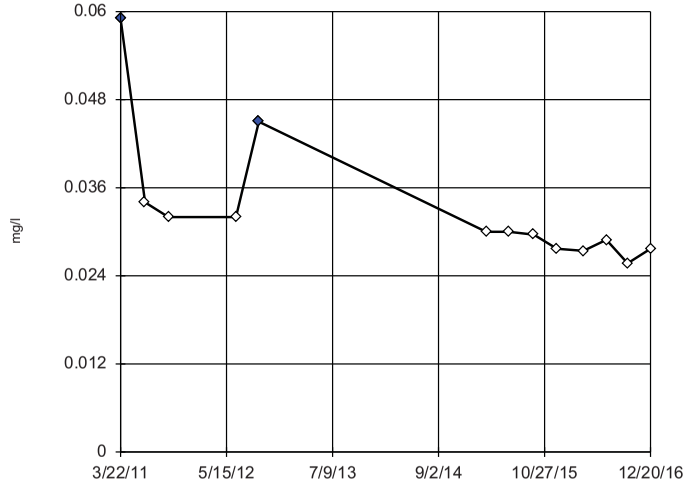
Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9028
Critical = 0.851
The distribution, after removal of suspect values, was found to be normally distributed.

Constituent: Barium Analysis Run 2/28/2017 10:00 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

MW-689D



n = 13

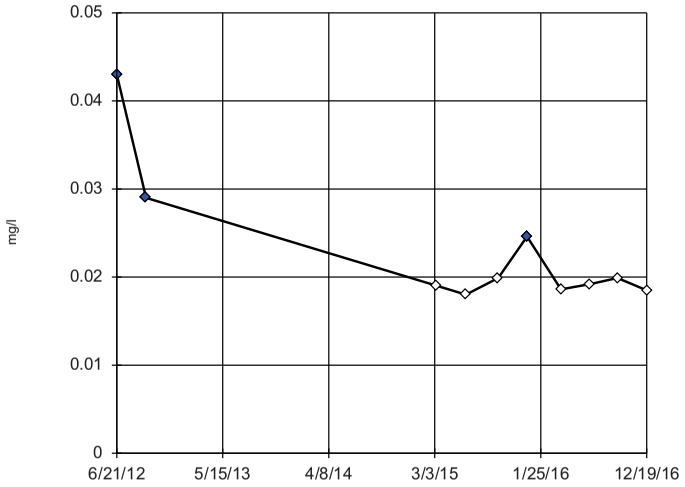
Statistical outliers are drawn as solid.
Testing for 2 high outliers.
Mean = 0.033.
Std. Dev. = 0.009175.
0.045: c = 0.7386
tabl = 0.521.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9675
Critical = 0.876
The distribution, after removal of suspect values, was found to be normally distributed.

Constituent: Barium Analysis Run 2/28/2017 10:01 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Dixon's Outlier Test

NE-2



n = 10

Statistical outliers are drawn as solid.
Testing for 3 high outliers.
Mean = 0.02297.
Std. Dev. = 0.007837.
0.0246: c = 0.7705
tabl = 0.477.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9337
Critical = 0.838
The distribution, after removal of suspect values, was found to be normally distributed.

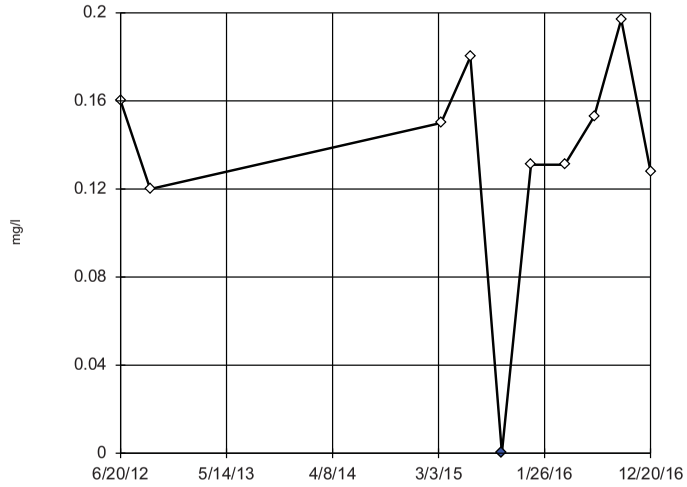
Constituent: Barium Analysis Run 2/28/2017 10:01 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

NE-3



n = 10

Statistical outlier is drawn as solid.
Testing for 1 low outlier.
Mean = 0.135.
Std. Dev. = 0.05326.
<0.00036: c = 0.6663
tabl = 0.477.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9209
Critical = 0.859
The distribution, after removal of suspect value, was found to be normally distributed.

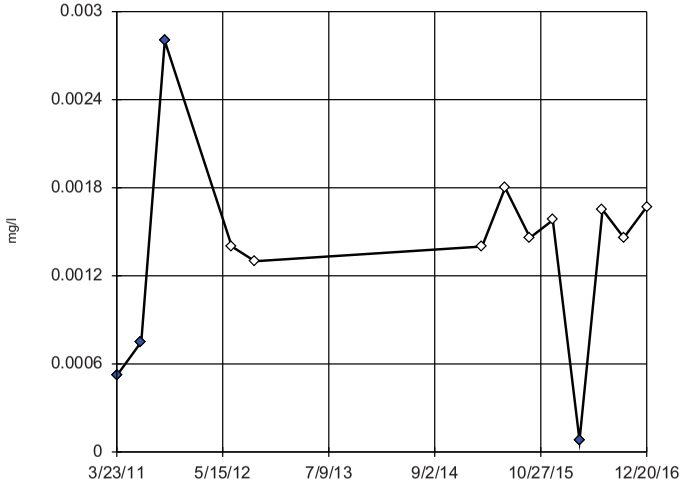
Constituent: Barium Analysis Run 2/28/2017 10:01 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

MW-633D



n = 13

Statistical outliers are drawn as solid.
Testing for 1 high and 3 low outliers.
Mean = 0.001375.
Std. Dev. = 0.0006606.
0.00167: c = 0.7533
tabl = 0.521.
0.00075: c = 0.619
tabl = 0.521.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9533
Critical = 0.859
The distribution, after removal of suspect values, was found to be normally distributed.

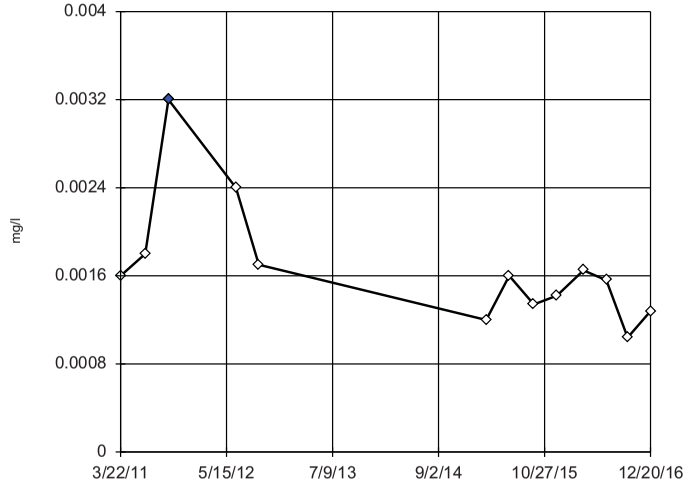
Constituent: Cadmium Analysis Run 2/28/2017 10:01 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

NAB-4



n = 13

Statistical outlier is drawn as solid.
Testing for 1 high outlier.
Mean = 0.001676.
Std. Dev. = 0.0005668.
0.0032: c = 0.7
tabl = 0.521.
Alpha = 0.05.

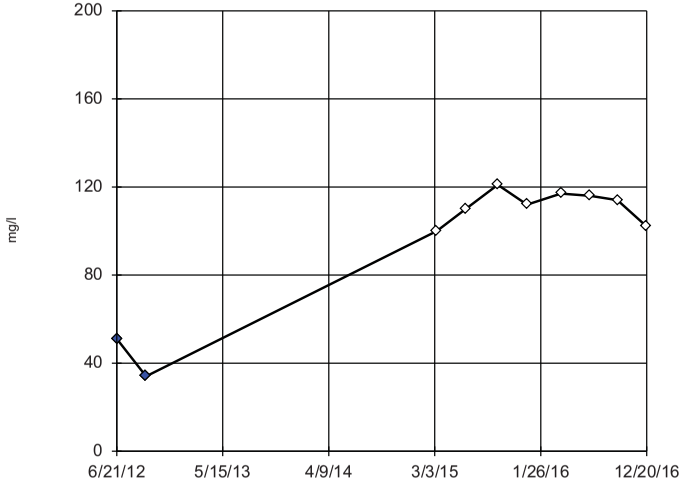
Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9195
Critical = 0.883
The distribution, after removal of suspect value, was found to be normally distributed.

Constituent: Cadmium Analysis Run 2/28/2017 10:01 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Dixon's Outlier Test

MW-1R



n = 10

Statistical outliers are drawn as solid.
Testing for 2 low outliers.
Mean = 97.7.
Std. Dev. = 30.06.
51: c = 0.7424
tabl = 0.477.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9313
Critical = 0.851
The distribution, after removal of suspect values, was found to be normally distributed.

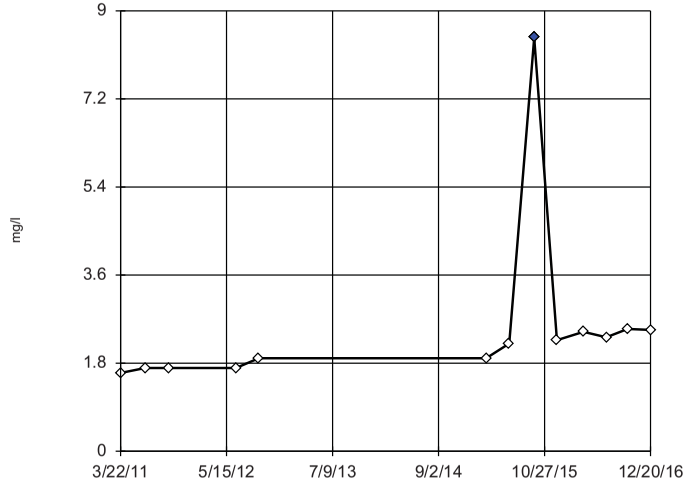
Constituent: Chloride Analysis Run 2/28/2017 10:01 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Tukey's Outlier Screening

MW-577



n = 13

Outlier is drawn as solid.
Tukey's method used in lieu of parametric test because the Shapiro Wilk normality test failed at the 0.1 alpha level.

Data were natural log transformed to achieve best W statistic (graph shown in original units).

High cutoff = 7.453, low cutoff = 0.5611, based on IQR multiplier of 3.

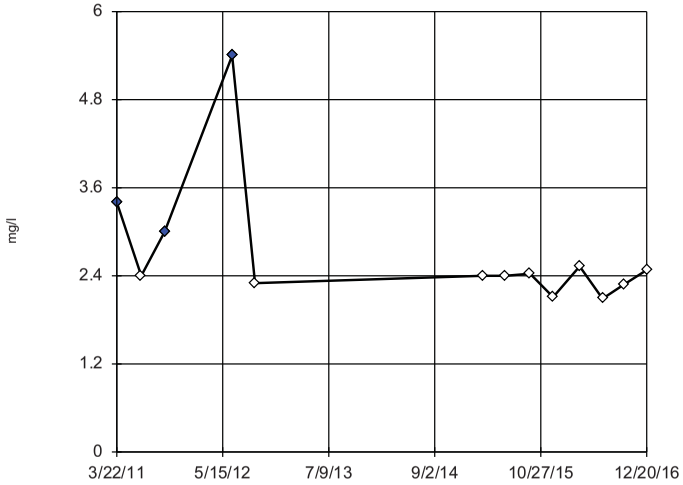
Constituent: Chloride Analysis Run 2/28/2017 10:01 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

MW-689D



n = 13

Statistical outliers are drawn as solid.
Testing for 3 high outliers.
Mean = 2.709.
Std. Dev. = 0.8825.
3: c = 0.5843
tabl = 0.521.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.898
Critical = 0.869
The distribution, after removal of suspect values, was found to be normally distributed.

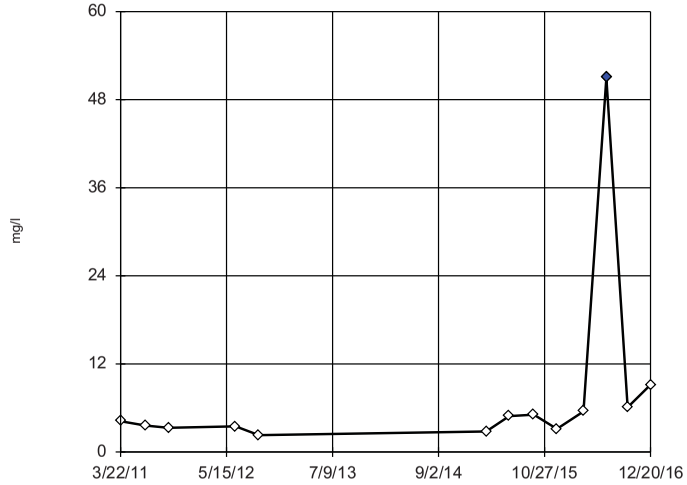
Constituent: Chloride Analysis Run 2/28/2017 10:01 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

NAB-3



n = 13

Statistical outlier is drawn as solid.
Testing for 1 high outlier.
Mean = 8.045.
Std. Dev. = 13.03.
51: c = 0.7328
tabl = 0.521.
Alpha = 0.05.

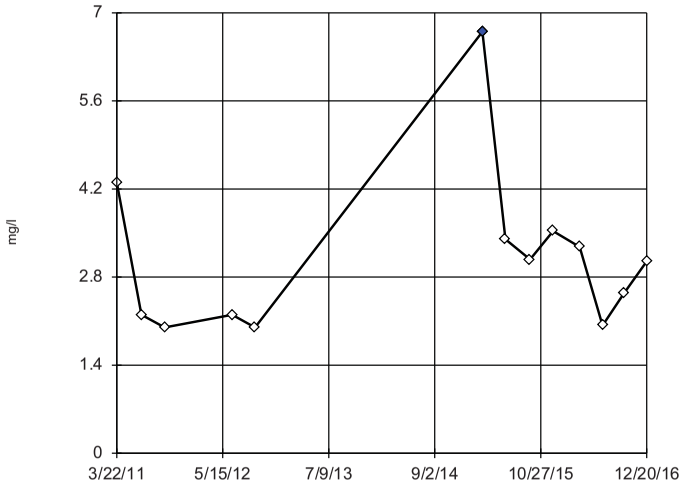
Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9747
Critical = 0.883 (after natural log transformation)
The distribution, after removal of suspect value, was found to be log-normal.

Constituent: Chloride Analysis Run 2/28/2017 10:01 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Dixon's Outlier Test

NAB-7



n = 13

Statistical outlier is drawn as solid.
Testing for 1 high outlier.
Mean = 3.101.
Std. Dev. = 1.298.
6.7: c = 0.6745
tabl = 0.521.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9
Critical = 0.883
The distribution, after removal of suspect value, was found to be normally distributed.

Constituent: Chloride Analysis Run 2/28/2017 10:01 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

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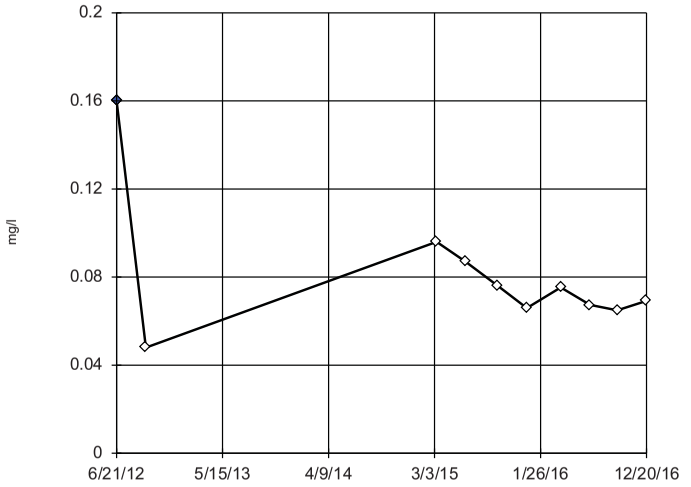
Dixon's Outlier Test

MW-1R



n = 10

Statistical outlier is drawn as solid.
Testing for 1 high outlier.
Mean = 0.08098.
Std. Dev. = 0.03066.
0.16: c = 0.673
tabl = 0.477.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9592
Critical = 0.859
The distribution, after removal of suspect value, was found to be normally distributed.

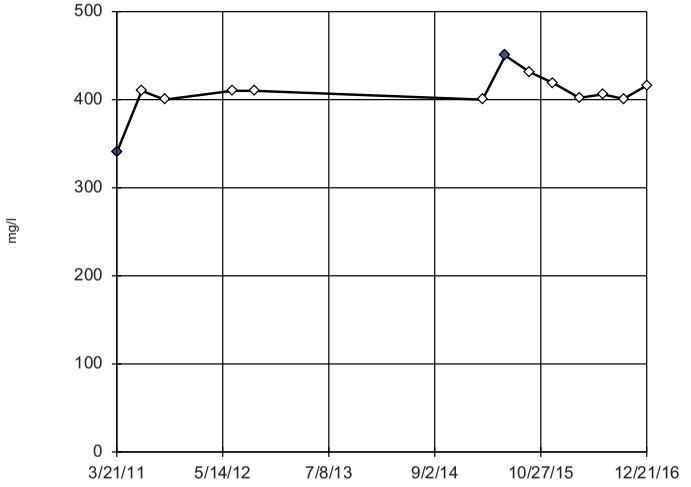
Constituent: Cobalt Analysis Run 2/28/2017 10:01 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

MW-5



n = 13

Statistical outliers are drawn as solid.
Testing for 1 high and 1 low outliers.
Mean = 407.3.
Std. Dev. = 24.71.
450: c = 0.62
tabl = 0.521.
340: c = 0.6593
tabl = 0.521.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.8843
Critical = 0.876
The distribution, after removal of suspect values, was found to be normally distributed.

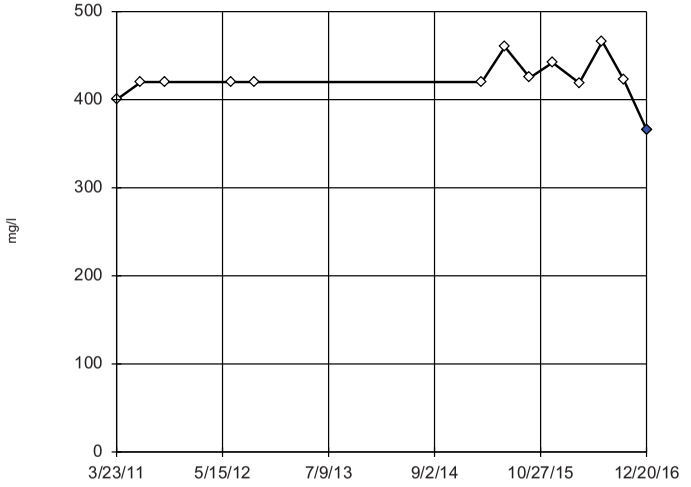
Constituent: Dissolved Solids Analysis Run 2/28/2017 10:01 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Tukey's Outlier Screening

MW-633D



n = 13

Outlier is drawn as solid.
Tukey's method used in lieu of parametric test because the Shapiro Wilk normality test failed at the 0.1 alpha level.

Data were cube transformed to achieve best W statistic (graph shown in original units).

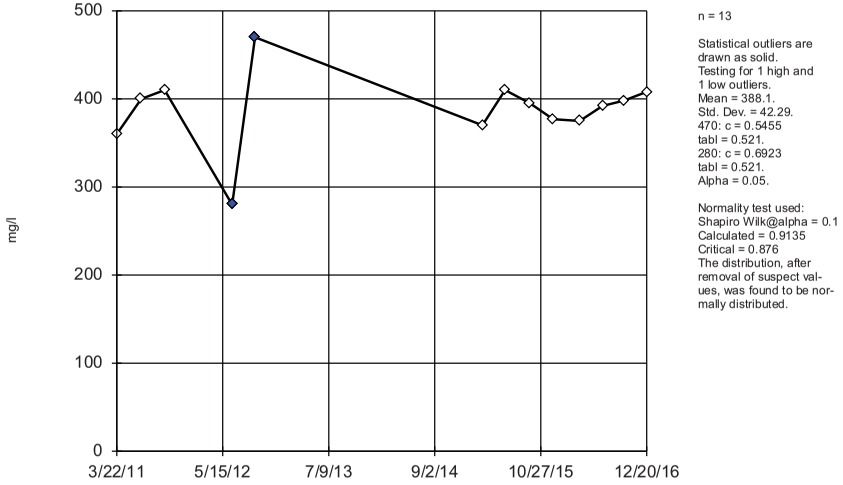
High cutoff = 471.4, low cutoff = 369.9, based on IQR multiplier of 3.

Constituent: Dissolved Solids Analysis Run 2/28/2017 10:01 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Dixon's Outlier Test

MW-689D



n = 13

Statistical outliers are drawn as solid.
Testing for 1 high and 1 low outliers.
Mean = 388.1.
Std. Dev. = 42.29.
470: c = 0.5455
tabl = 0.521.
280: c = 0.6923
tabl = 0.521.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9135
Critical = 0.876
The distribution, after removal of suspect values, was found to be normally distributed.

Constituent: Dissolved Solids Analysis Run 2/28/2017 10:01 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

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Dixon's Outlier Test

MW-7

n = 13

Statistical outlier is drawn as solid.
Testing for 1 high and 1 low outliers.
Mean = 339.2.
Std. Dev. = 26.5.
390: c = 0.4286
tabl = 0.521.
274: c = 0.6163
tabl = 0.521.
Alpha = 0.05.

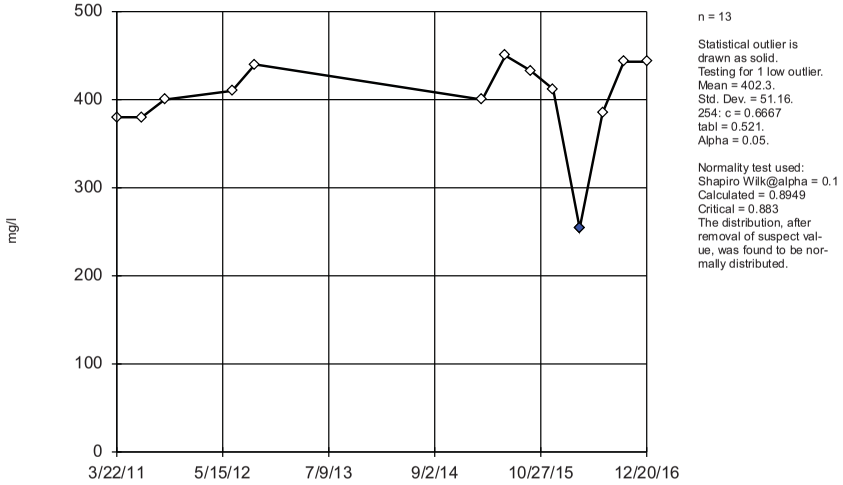
Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9401
Critical = 0.876
The distribution, after removal of suspect value, was found to be normally distributed.

Constituent: Dissolved Solids Analysis Run 2/28/2017 10:01 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

NAB-2



n = 13

Statistical outlier is drawn as solid.
Testing for 1 low outlier.
Mean = 402.3.
Std. Dev. = 51.16.
254: c = 0.6667
tabl = 0.521.
Alpha = 0.05.

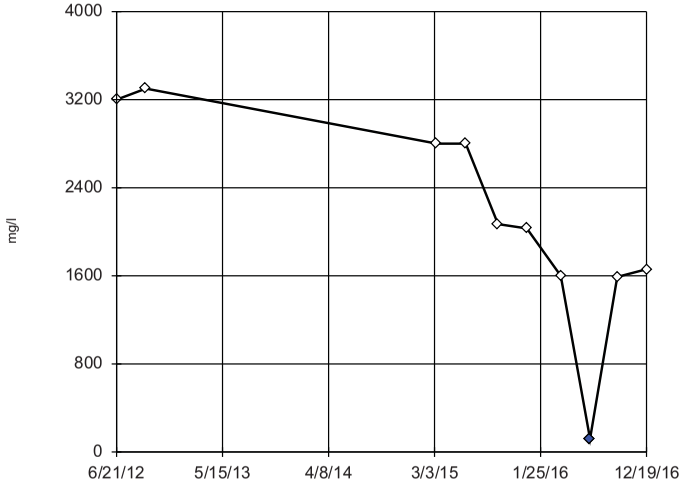
Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.8949
Critical = 0.883
The distribution, after removal of suspect value, was found to be normally distributed.

Constituent: Dissolved Solids Analysis Run 2/28/2017 10:01 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

NE-2



n = 10

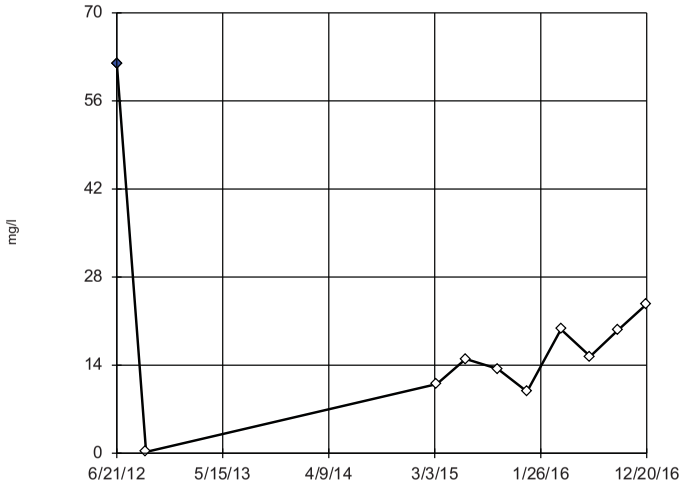
Statistical outlier is drawn as solid.
Testing for 1 low outlier.
Mean = 2117.
Std. Dev. = 957.8.
120: c = 0.4773
tabl = 0.477.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.8697
Critical = 0.859
The distribution, after removal of suspect value, was found to be normally distributed.

Constituent: Dissolved Solids Analysis Run 2/28/2017 10:01 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Dixon's Outlier Test

MW-1R



n = 10

Statistical outlier is drawn as solid.
Testing for 1 high and 1 low outliers.
Mean = 18.97.
Std. Dev. = 16.44.
62: c = 0.7339
tabl = 0.477.
0.21: c = 0.4087
tabl = 0.477.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9572
Critical = 0.851
The distribution, after removal of suspect value, was found to be normally distributed.

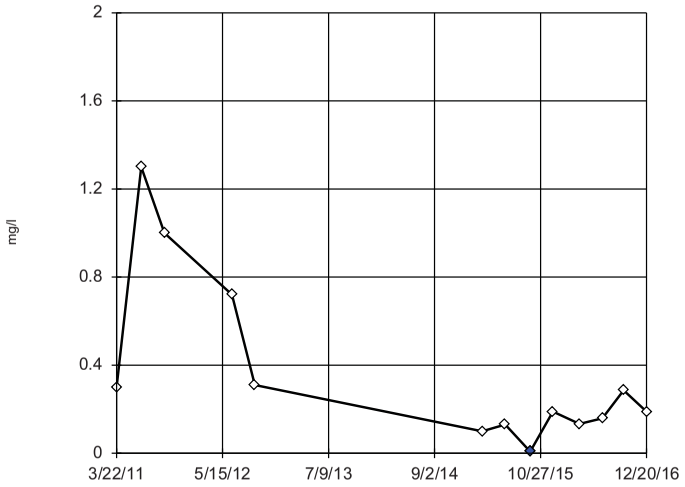
Constituent: Iron Analysis Run 2/28/2017 10:01 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

MW-577



n = 13

Statistical outlier is drawn as solid.
Testing for 1 low outlier.
Mean = 0.371.
Std. Dev. = 0.3903.
-4.89285228..: c = 0.583
tabl = 0.521.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9032
Critical = 0.883 (after natural log transformation)
The distribution, after removal of suspect value, was found to be log-normal.

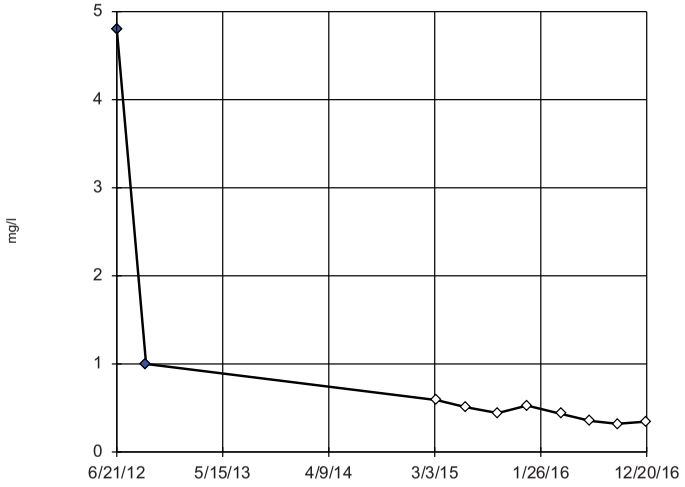
Constituent: Iron Analysis Run 2/28/2017 10:01 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

MW-1R



n = 10

Statistical outliers are drawn as solid.
Testing for 2 high outliers.
Mean = 0.9322.
Std. Dev. = 1.373.
1: c = 0.6269
tabl = 0.477.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9383
Critical = 0.851
The distribution, after removal of suspect values, was found to be normally distributed.

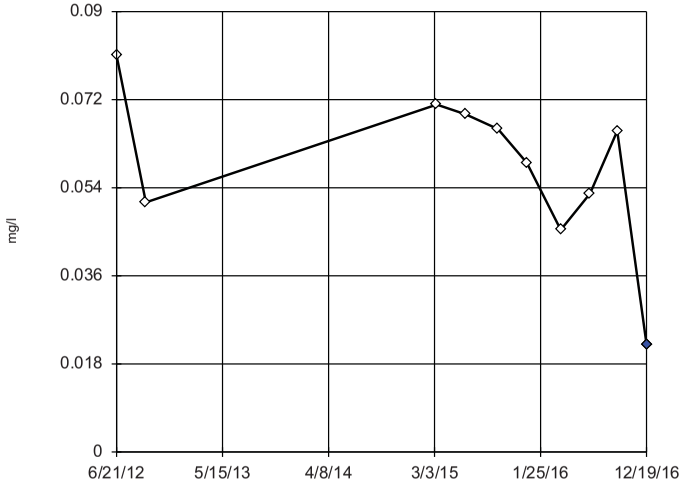
Constituent: Manganese Analysis Run 2/28/2017 10:02 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

NE-2



n = 10

Statistical outlier is drawn as solid.
Testing for 1 low outlier.
Mean = 0.05829.
Std. Dev. = 0.01657.
0.022: c = 0.4776
tabl = 0.477.
Alpha = 0.05.

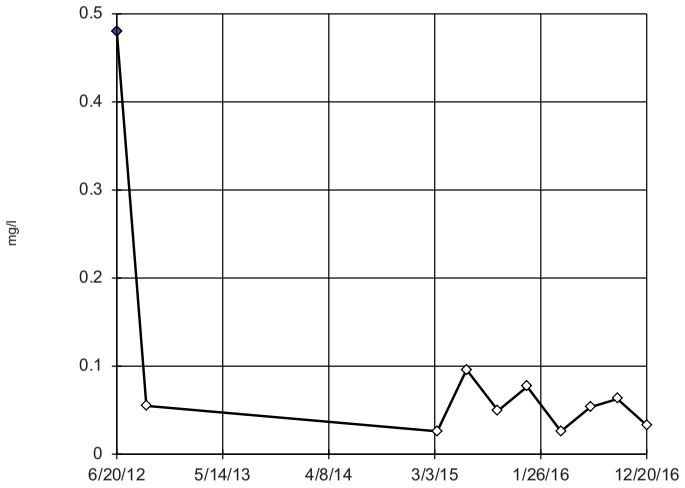
Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9715
Critical = 0.859
The distribution, after removal of suspect value, was found to be normally distributed.

Constituent: Manganese Analysis Run 2/28/2017 10:02 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Dixon's Outlier Test

NE-3



n = 10

Statistical outlier is drawn as solid.
Testing for 1 high outlier.
Mean = 0.09582.
Std. Dev. = 0.1368.
0.48: c = 0.8458
tabl = 0.477.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9367
Critical = 0.859
The distribution, after removal of suspect value, was found to be normally distributed.

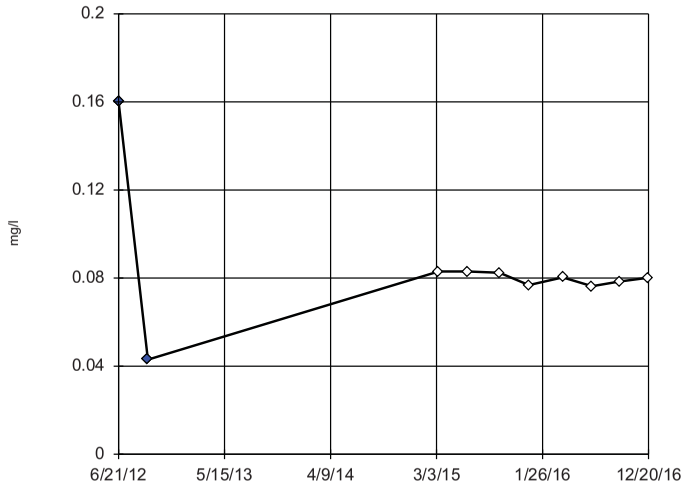
Constituent: Manganese Analysis Run 2/28/2017 10:02 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

MW-1R



n = 10

Statistical outliers are drawn as solid.
Testing for 1 high and 1 low outliers.
Mean = 0.08434.
Std. Dev. = 0.02912.
0.16: c = 0.92
tabl = 0.477.
0.043: c = 0.8325
tabl = 0.477.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.8965
Critical = 0.851
The distribution, after removal of suspect values, was found to be normally distributed.

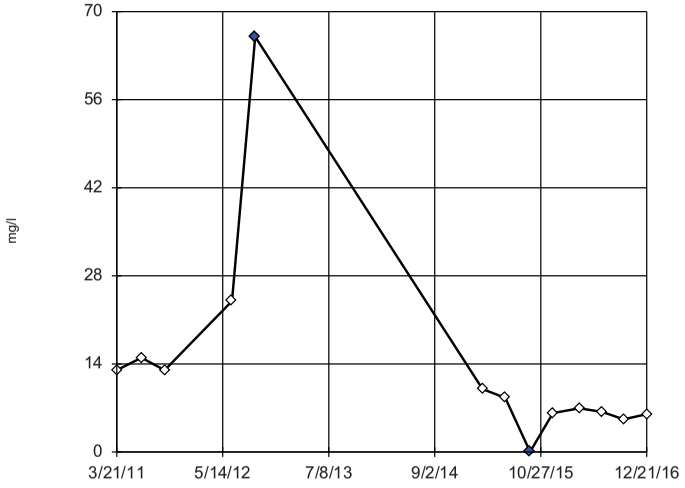
Constituent: Nickel Analysis Run 2/28/2017 10:02 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

CAO-3



n = 13

Statistical outliers are drawn as solid.
Testing for 1 high and 1 low outliers.
Mean = 13.87.
Std. Dev. = 16.73.
66: c = 0.5826
tabl = 0.521.
-3.25191569..: c = 0.7847
tabl = 0.521.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9269
Critical = 0.876 (after natural log transformation)
The distribution, after removal of suspect values, was found to be log-normal.

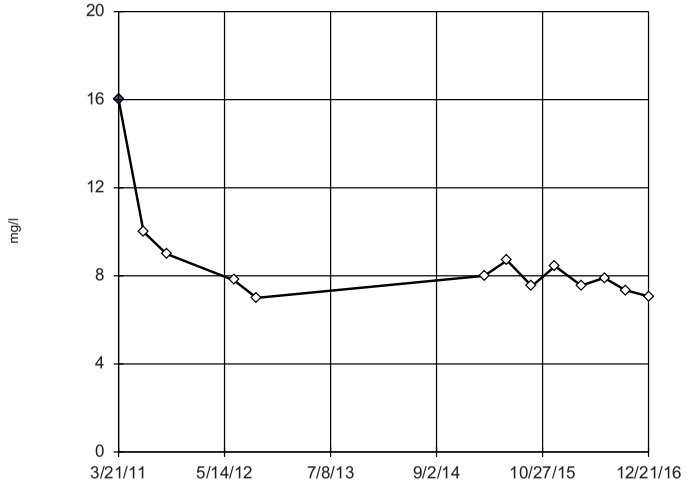
Constituent: Sulfate Analysis Run 2/28/2017 10:02 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

MW-5



n = 13

Statistical outlier is drawn as solid.
Testing for 1 high outlier.
Mean = 8.643.
Std. Dev. = 2.364.
16: c = 0.7839
tabl = 0.521.
Alpha = 0.05.

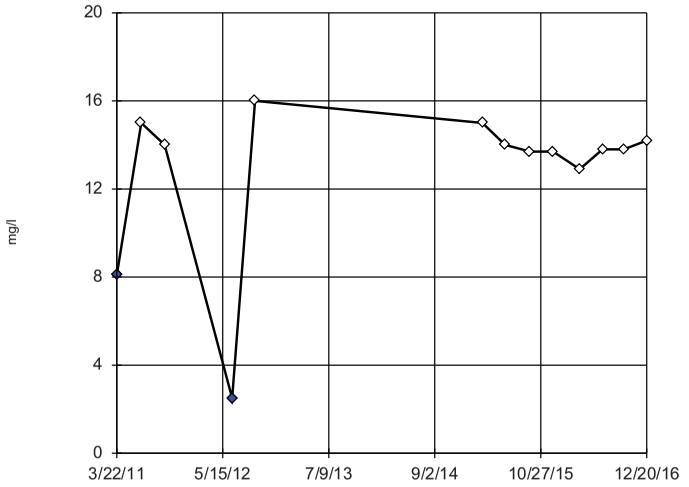
Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9237
Critical = 0.883
The distribution, after removal of suspect value, was found to be normally distributed.

Constituent: Sulfate Analysis Run 2/28/2017 10:02 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Dixon's Outlier Test

MW-689D



n = 13

Statistical outliers are drawn as solid.
Testing for 2 low outliers.
Mean = 12.82.
Std. Dev. = 3.612.
8.1: c = 0.8116
tabl = 0.521.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9019
Critical = 0.876
The distribution, after removal of suspect values, was found to be normally distributed.

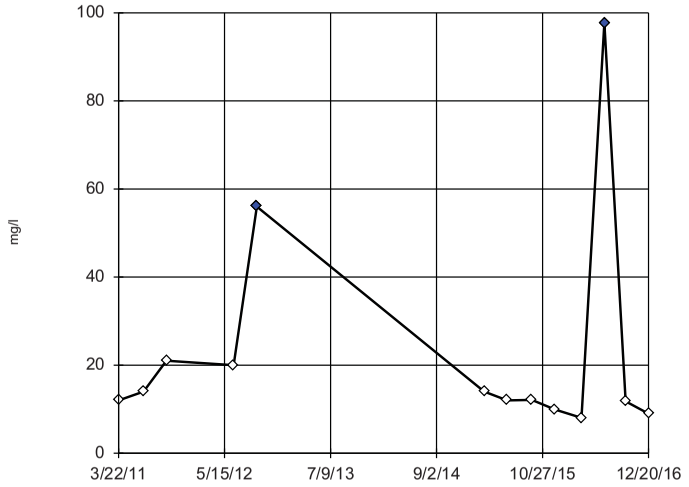
Constituent: Sulfate Analysis Run 2/28/2017 10:02 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

NAB-3



n = 13

Statistical outliers are drawn as solid.
Testing for 2 high outliers.
Mean = 22.89.
Std. Dev. = 25.69.
56: c = 0.5622
tabl = 0.521.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9322
Critical = 0.876 (after natural log transformation)
The distribution, after removal of suspect values, was found to be log-normal.

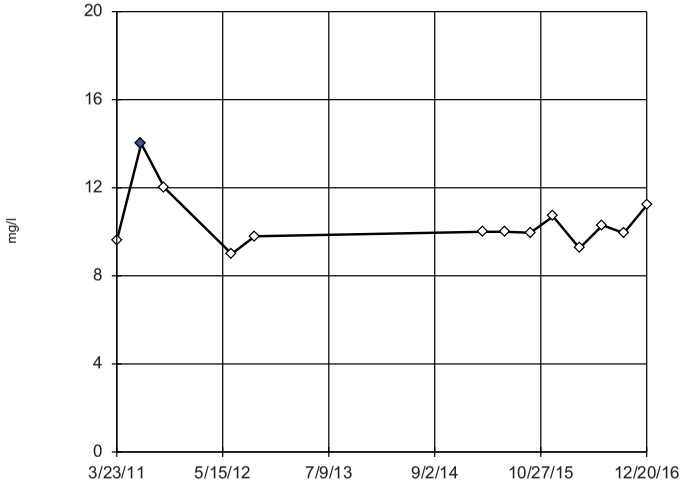
Constituent: Sulfate Analysis Run 2/28/2017 10:02 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

NAB-8



n = 13

Statistical outlier is drawn as solid.
Testing for 1 high outlier.
Mean = 10.44.
Std. Dev. = 1.331.
14: c = 0.5895
tabl = 0.521.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9218
Critical = 0.883
The distribution, after removal of suspect value, was found to be normally distributed.

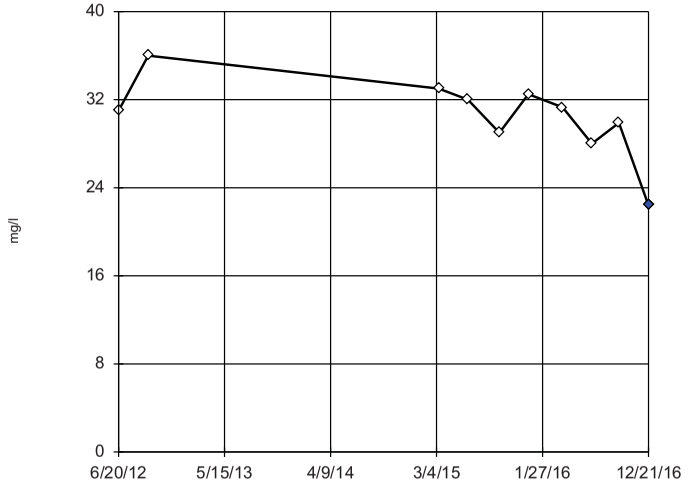
Constituent: Sulfate Analysis Run 2/28/2017 10:02 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

NE-6



n = 10

Statistical outlier is drawn as solid.
Testing for 1 low outlier.
Mean = 30.51.
Std. Dev. = 3.623.
22.4: c = 0.5283
tabl = 0.477.
Alpha = 0.05.

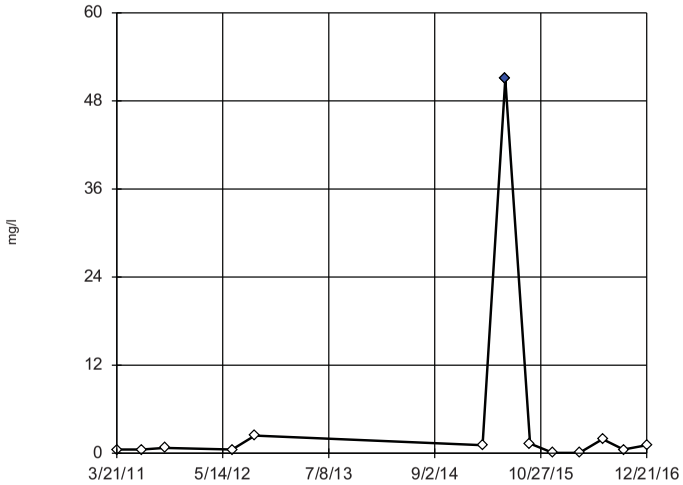
Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9722
Critical = 0.859
The distribution, after removal of suspect value, was found to be normally distributed.

Constituent: Sulfate Analysis Run 2/28/2017 10:02 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Dixon's Outlier Test

MW-4



n = 13

Statistical outlier is drawn as solid.
Testing for 1 high outlier.
Mean = 4.742.
Std. Dev. = 13.92.
51: c = 0.9631
tabl = 0.521.
Alpha = 0.05.

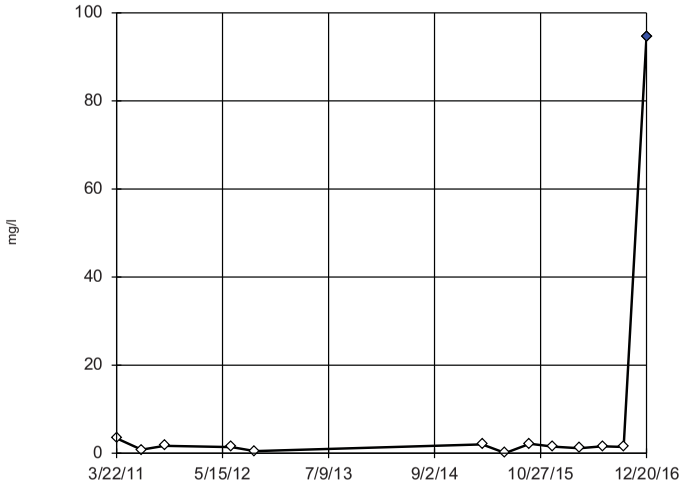
Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.899
Critical = 0.883
The distribution, after removal of suspect value, was found to be normally distributed.

Constituent: TOC [Total Organic Carbon] Analysis Run 2/28/2017 10:02 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

NAB-3



n = 13

Statistical outlier is drawn as solid.
Testing for 1 high and 1 low outliers.
Mean = 8.645.
Std. Dev. = 25.87.
94.7: c = 0.9823
tabl = 0.521.
<0.1: c = 0.2239
tabl = 0.521.
Alpha = 0.05.

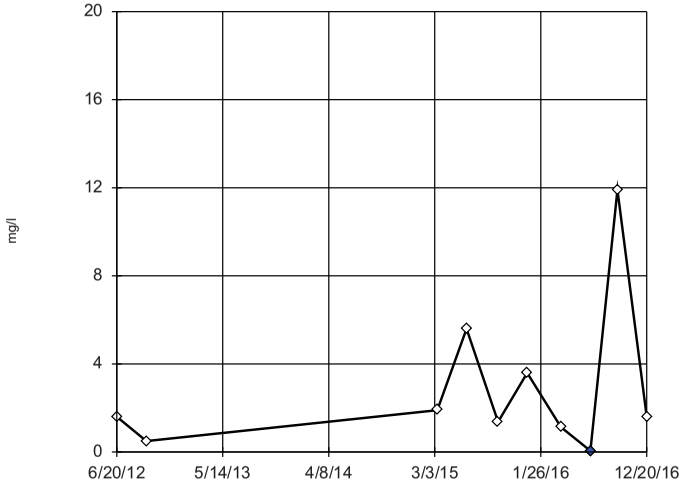
Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9212
Critical = 0.876
The distribution, after removal of suspect value, was found to be normally distributed.

Constituent: TOC [Total Organic Carbon] Analysis Run 2/28/2017 10:02 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

NE-3



n = 10

Statistical outlier is drawn as solid.
Testing for 1 low outlier.
Mean = 2.924.
Std. Dev. = 3.535.
-2.97592966..: c = 0.4858
tabl = 0.477.
Alpha = 0.05.

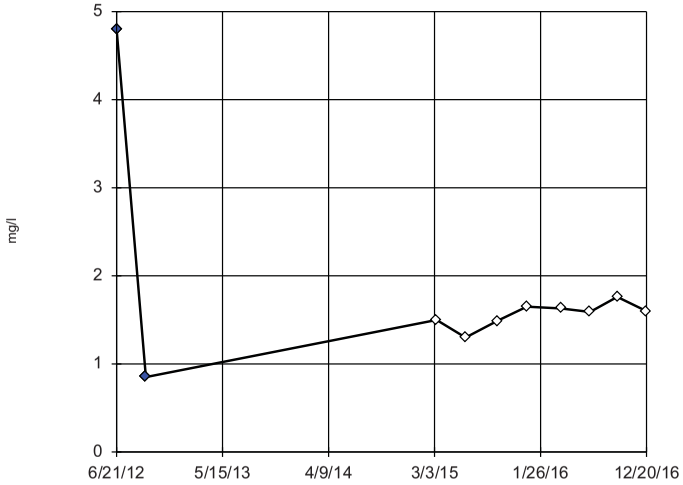
Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9544
Critical = 0.859 (after natural log transformation)
The distribution, after removal of suspect value, was found to be log-normal.

Constituent: TOC [Total Organic Carbon] Analysis Run 2/28/2017 10:02 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

MW-1R



n = 10

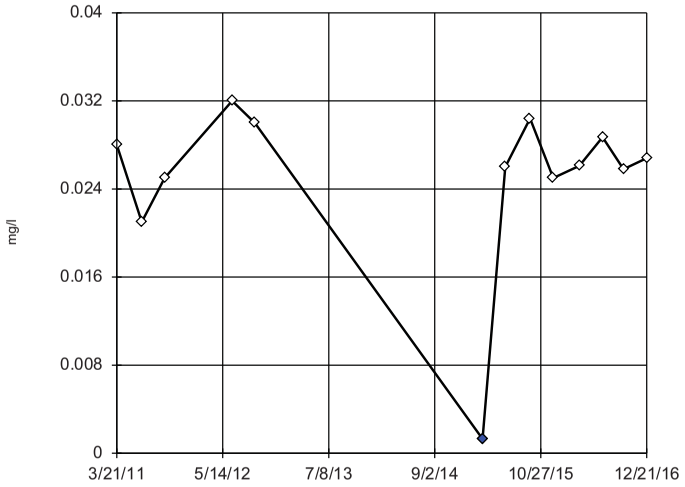
Statistical outliers are drawn as solid.
Testing for 1 high and 1 low outliers.
Mean = 1.817.
Std. Dev. = 1.079.
4.8: c = 0.8686
tabl = 0.477.
0.85: c = 0.4945
tabl = 0.477.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9442
Critical = 0.851
The distribution, after removal of suspect values, was found to be normally distributed.

Constituent: Zinc Analysis Run 2/28/2017 10:02 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Dixon's Outlier Test

MW-5



n = 13

Statistical outlier is drawn as solid.
Testing for 1 low outlier.
Mean = 0.02508.
Std. Dev. = 0.007689.
<0.0026: c = 0.8144
tabl = 0.521.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9645
Critical = 0.883
The distribution, after removal of suspect value, was found to be normally distributed.

Constituent: Zinc Analysis Run 2/28/2017 10:02 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

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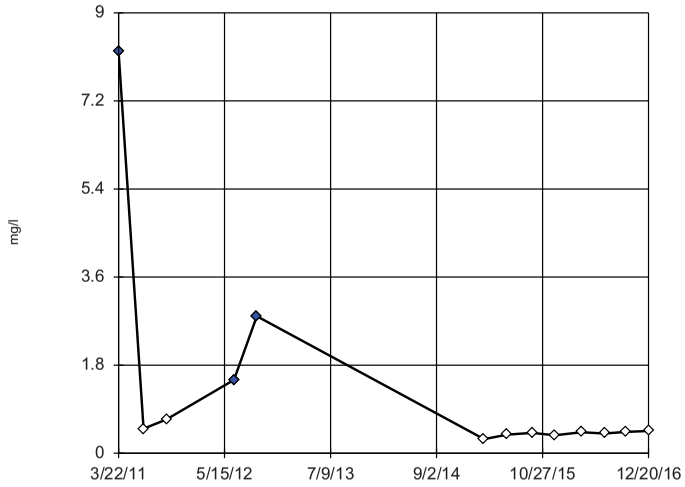
Dixon's Outlier Test

MW-509D



n = 13

Statistical outliers are drawn as solid.
Testing for 3 high outliers.
Mean = 1.297.
Std. Dev. = 2.188.
1.5: c = 0.8818
tabl = 0.521.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.8697
Critical = 0.869
The distribution, after removal of suspect values, was found to be normally distributed.

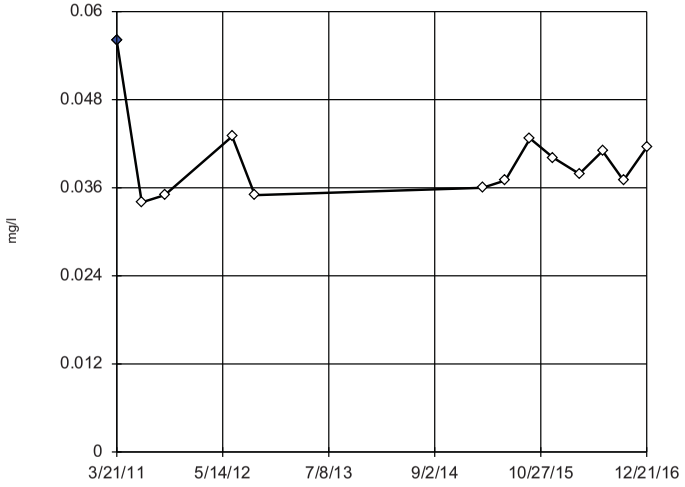
Constituent: Zinc Analysis Run 2/28/2017 10:02 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

MW-6



n = 13

Statistical outlier is drawn as solid.
Testing for 1 high outlier.
Mean = 0.03971.
Std. Dev. = 0.005765.
0.056: c = 0.6333
tabl = 0.521.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.9201
Critical = 0.883
The distribution, after removal of suspect value, was found to be normally distributed.

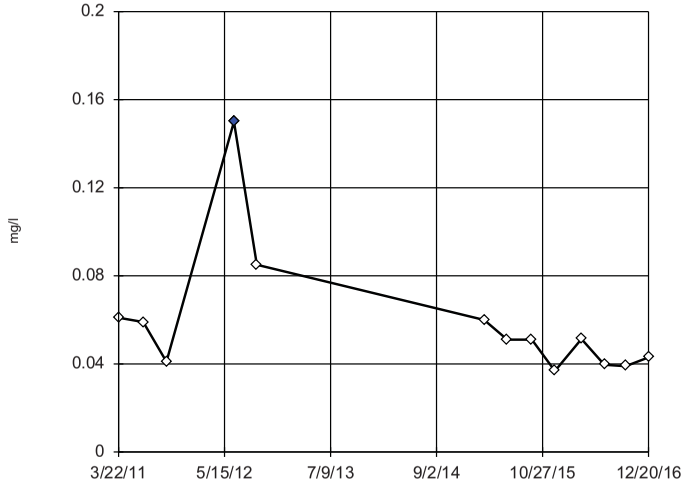
Constituent: Zinc Analysis Run 2/28/2017 10:02 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

NAB-7



n = 13

Statistical outlier is drawn as solid.
Testing for 1 high outlier.
Mean = 0.05905.
Std. Dev. = 0.03027.
0.15: c = 0.6679
tabl = 0.521.
Alpha = 0.05.

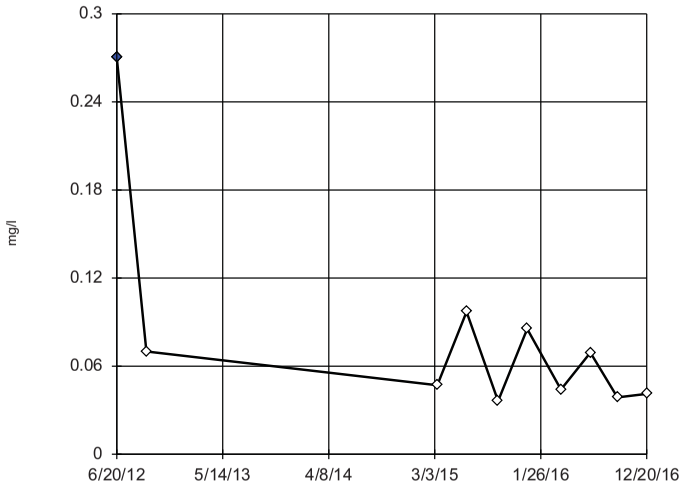
Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.925
Critical = 0.883 (after natural log transformation)
The distribution, after removal of suspect value, was found to be log-normal.

Constituent: Zinc Analysis Run 2/28/2017 10:02 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Dixon's Outlier Test

NE-3



n = 10

Statistical outlier is drawn as solid.
Testing for 1 high outlier.
Mean = 0.07988.
Std. Dev. = 0.07004.
0.27: c = 0.7479
tabl = 0.477.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.8747
Critical = 0.859
The distribution, after removal of suspect value, was found to be normally distributed.

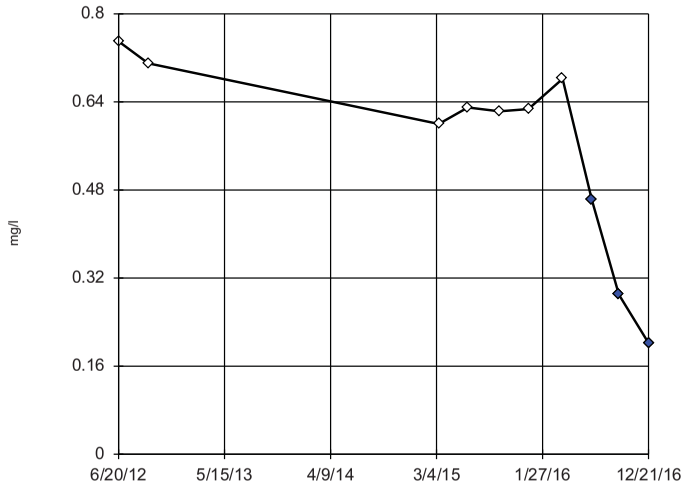
Constituent: Zinc Analysis Run 2/28/2017 10:02 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Dixon's Outlier Test

NE-6



n = 10

Statistical outliers are drawn as solid.
Testing for 3 low outliers.
Mean = 0.5576.
Std. Dev. = 0.1823.
0.462: c = 0.5565
tabl = 0.477.
Alpha = 0.05.

Normality test used:
Shapiro Wilk@alpha = 0.1
Calculated = 0.902
Critical = 0.838
The distribution, after removal of suspect values, was found to be normally distributed.

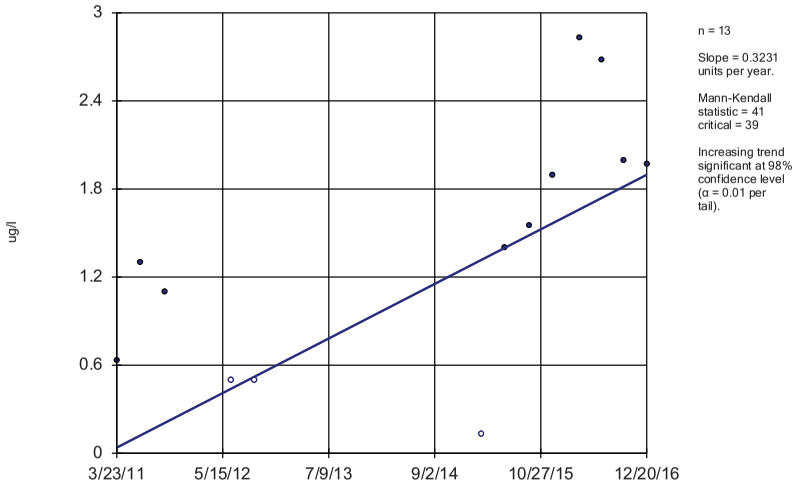
Constituent: Zinc Analysis Run 2/28/2017 10:02 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG
Hollow symbols indicate censored values.

Sen's Slope Estimator

MW-633D



n = 13

Slope = 0.3231 units per year.

Mann-Kendall statistic = 41
critical = 39

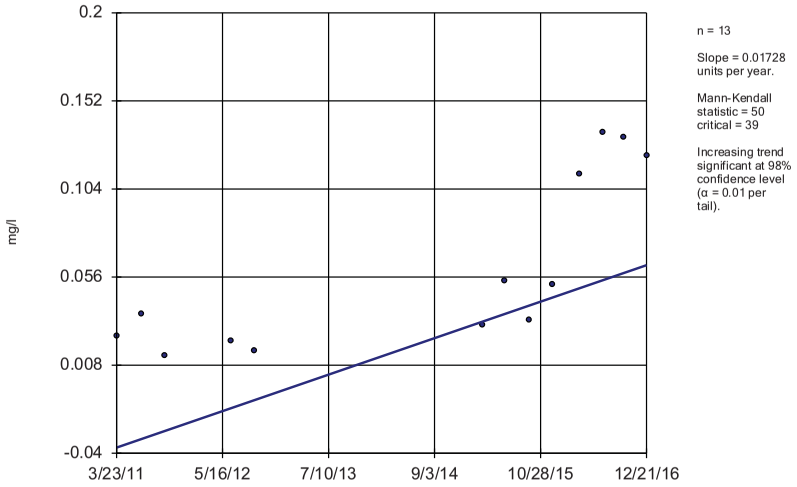
Increasing trend significant at 98% confidence level (α = 0.01 per tail).

Constituent: 1,1-Dichloroethane Analysis Run 2/28/2017 10:03 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

CAO-1



n = 13

Slope = 0.01728 units per year.

Mann-Kendall statistic = 50
critical = 39

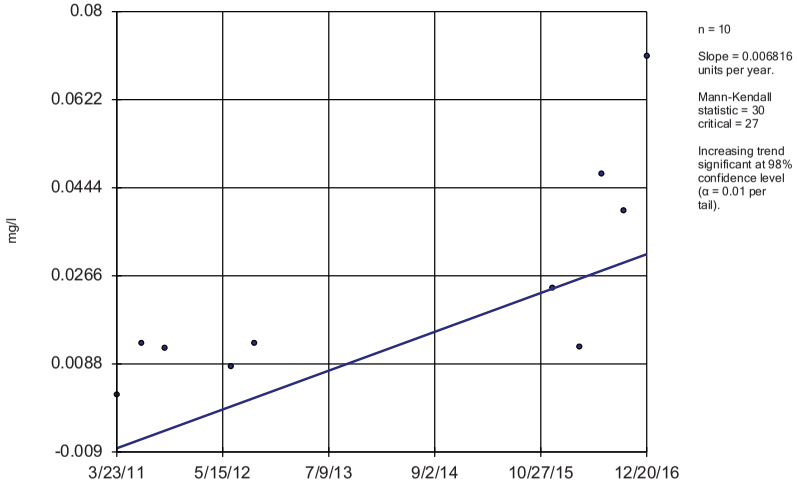
Increasing trend significant at 98% confidence level (α = 0.01 per tail).

Constituent: Arsenic Analysis Run 2/28/2017 10:03 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-1



n = 10

Slope = 0.006816 units per year.

Mann-Kendall statistic = 30
critical = 27

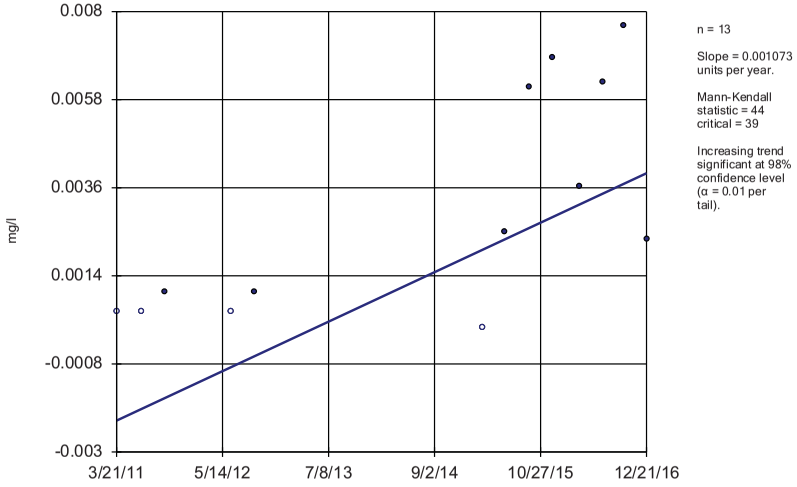
Increasing trend significant at 98% confidence level (α = 0.01 per tail).

Constituent: Arsenic Analysis Run 2/28/2017 10:03 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG
Hollow symbols indicate censored values.

Sen's Slope Estimator

MW-3



n = 13

Slope = 0.001073 units per year.

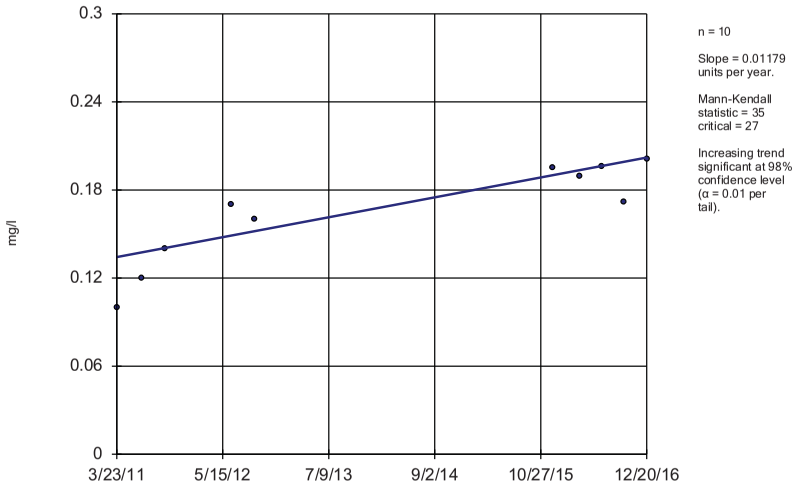
Mann-Kendall statistic = 44
critical = 39

Increasing trend significant at 98% confidence level (α = 0.01 per tail).

Constituent: Arsenic Analysis Run 2/28/2017 10:03 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sen's Slope Estimator

MW-1



n = 10

Slope = 0.01179 units per year.

Mann-Kendall statistic = 35
critical = 27

Increasing trend significant at 98% confidence level (α = 0.01 per tail).

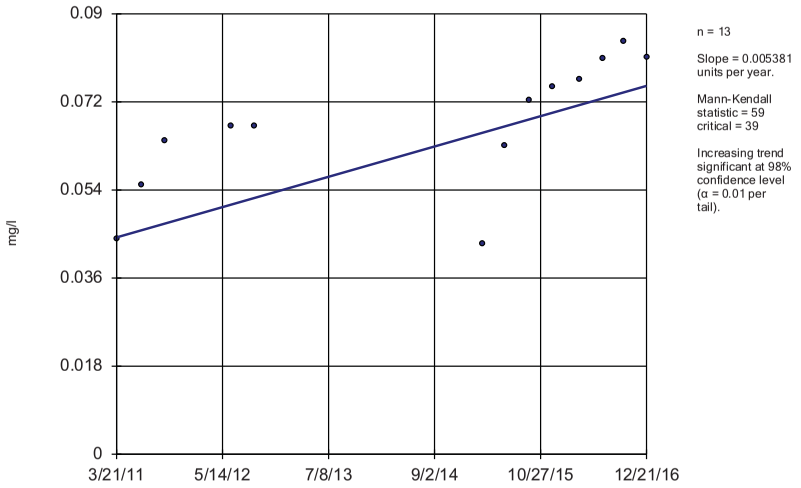
Constituent: Barium Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-3



n = 13

Slope = 0.005381 units per year.

Mann-Kendall statistic = 59
critical = 39

Increasing trend significant at 98% confidence level (α = 0.01 per tail).

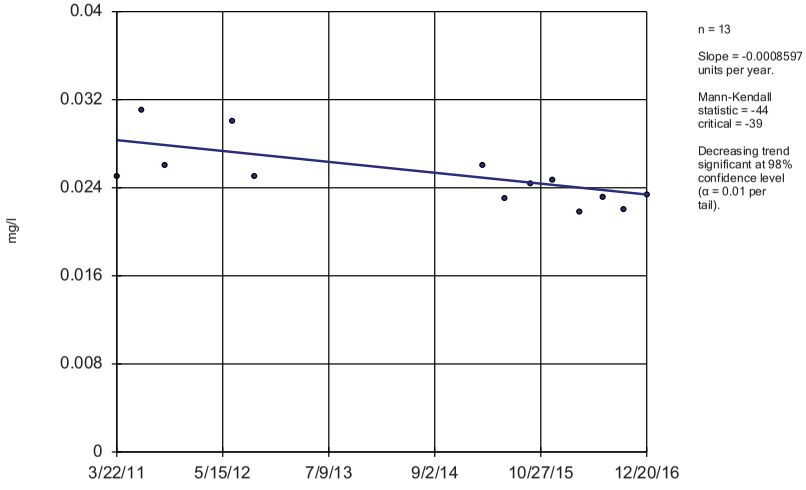
Constituent: Barium Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-577



n = 13

Slope = -0.0008597 units per year.

Mann-Kendall statistic = -44
critical = -39

Decreasing trend significant at 98% confidence level (α = 0.01 per tail).

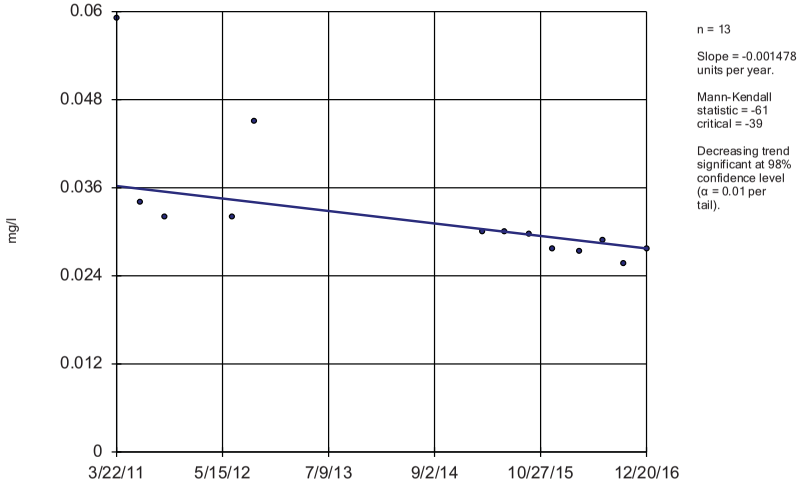
Constituent: Barium Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-689D



n = 13

Slope = -0.001478 units per year.

Mann-Kendall statistic = -61
critical = -39

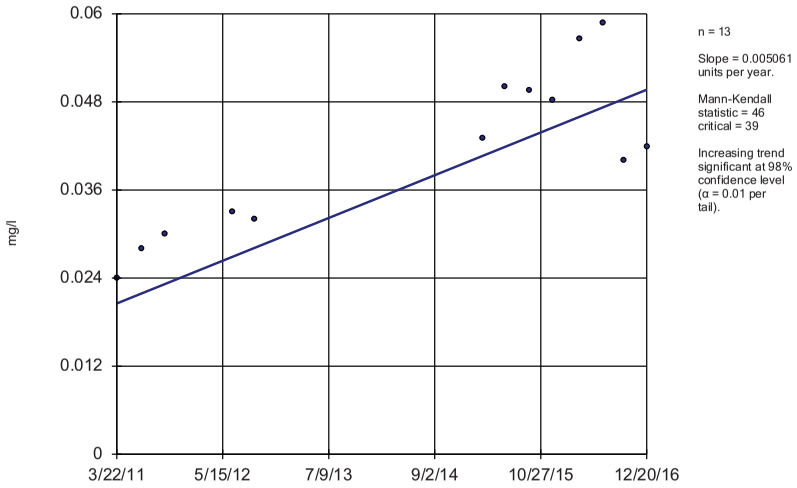
Decreasing trend significant at 98% confidence level (α = 0.01 per tail).

Constituent: Barium Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sen's Slope Estimator

NAB-4



n = 13

Slope = 0.005061 units per year.

Mann-Kendall statistic = 46
critical = 39

Increasing trend significant at 98% confidence level (α = 0.01 per tail).

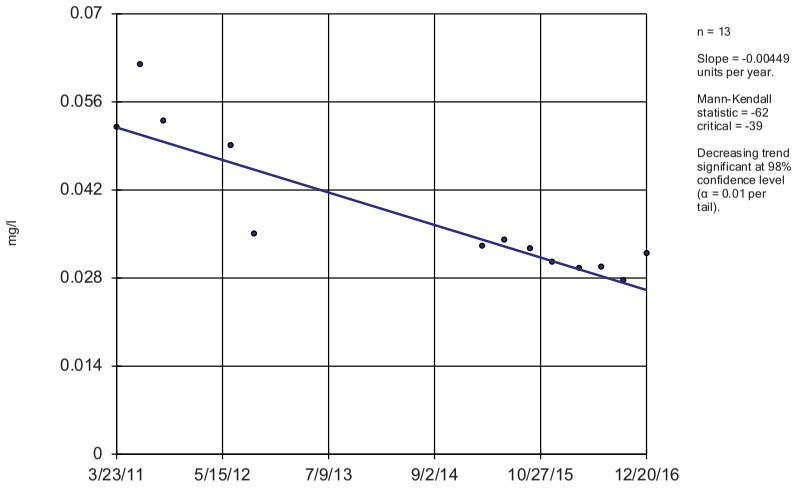
Constituent: Barium Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

NAB-8



n = 13

Slope = -0.00449 units per year.

Mann-Kendall statistic = -62
critical = -39

Decreasing trend significant at 98% confidence level (α = 0.01 per tail).

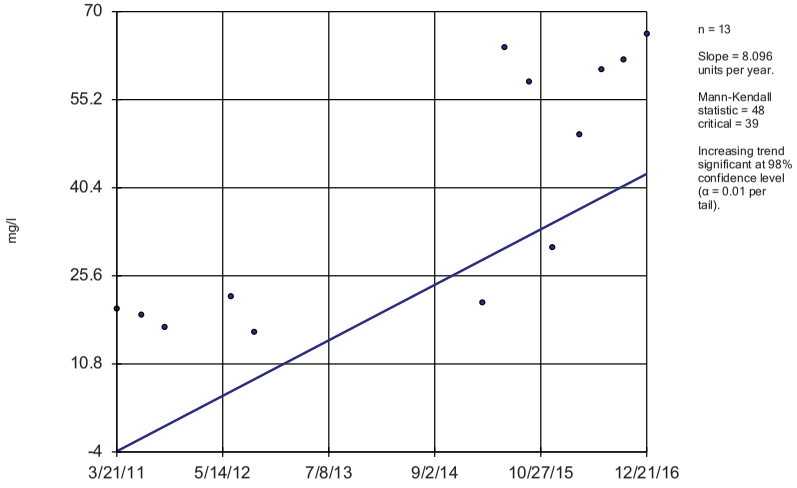
Constituent: Barium Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

CAO-2



n = 13

Slope = 8.096 units per year.

Mann-Kendall statistic = 48
critical = 39

Increasing trend significant at 98% confidence level (α = 0.01 per tail).

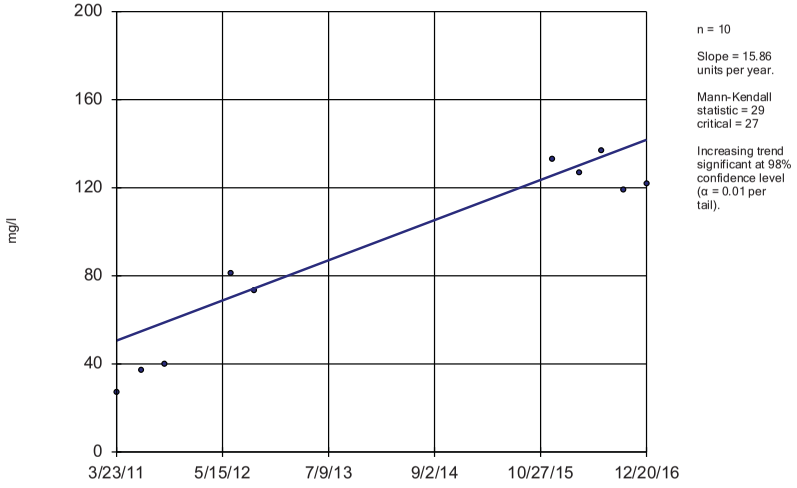
Constituent: Chloride Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-1



n = 10

Slope = 15.86 units per year.

Mann-Kendall statistic = 29
critical = 27

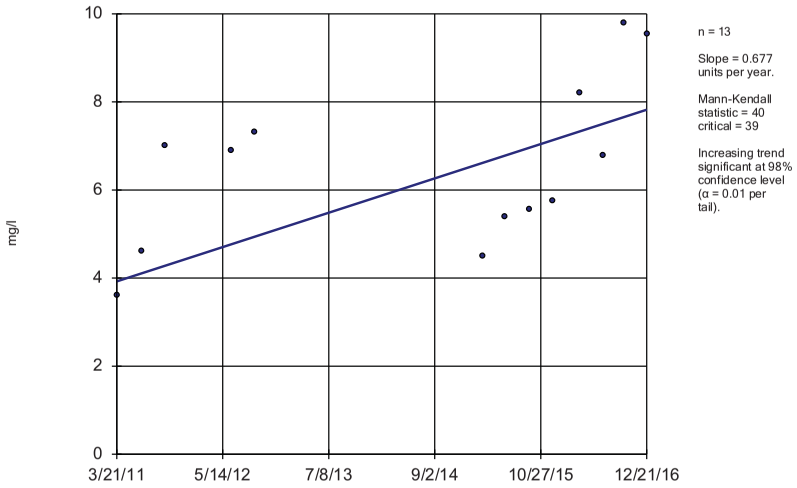
Increasing trend significant at 98% confidence level (α = 0.01 per tail).

Constituent: Chloride Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sen's Slope Estimator

MW-3



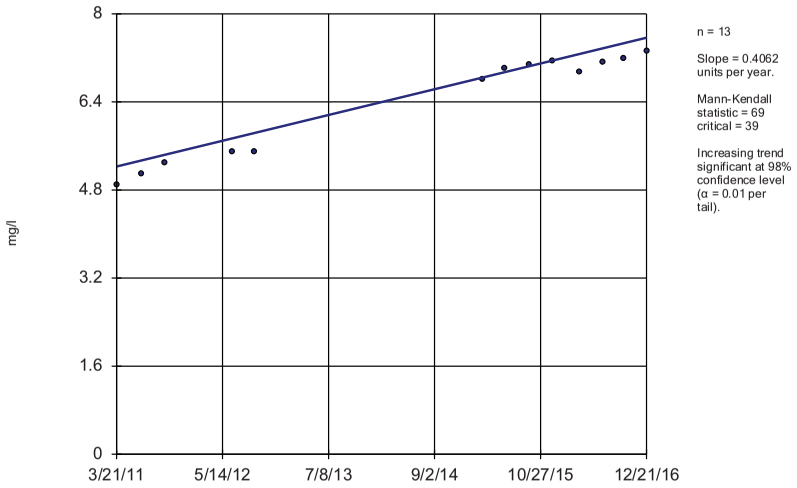
Constituent: Chloride Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-4



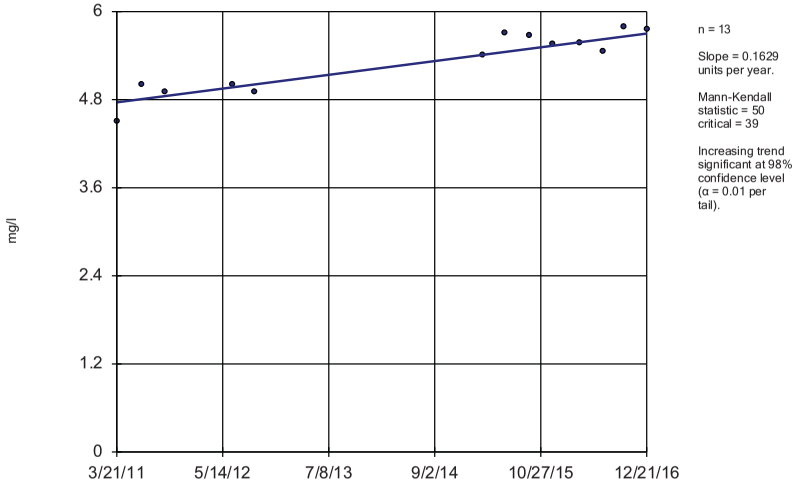
Constituent: Chloride Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-5



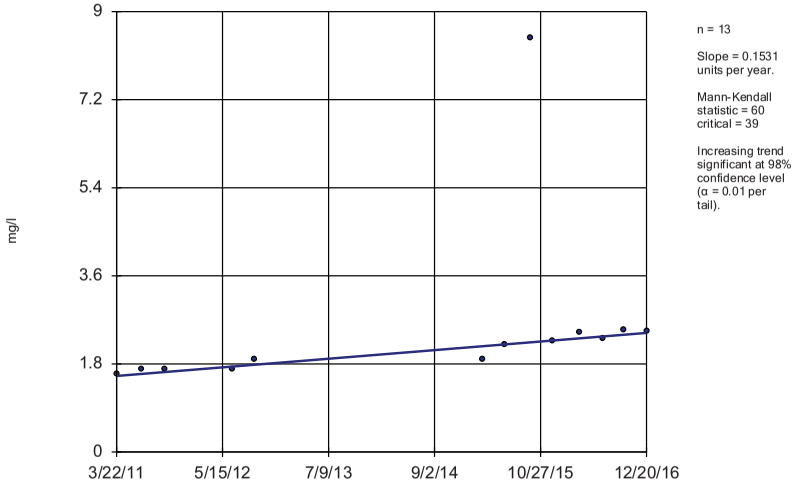
Constituent: Chloride Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-577

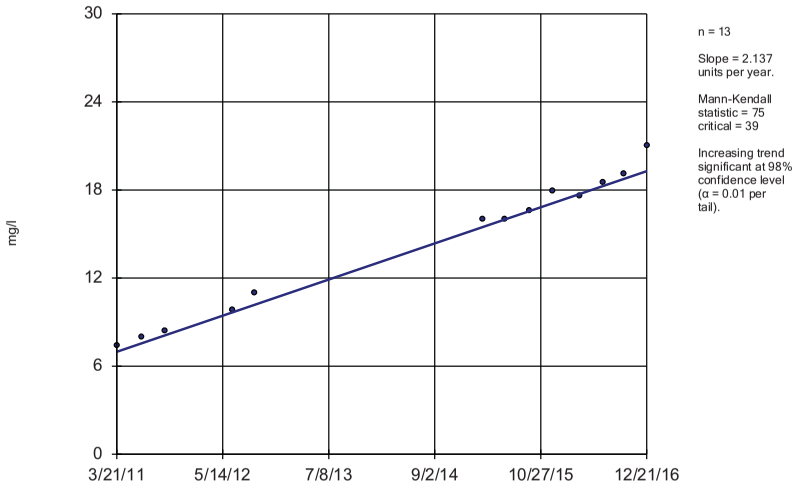


Constituent: Chloride Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sen's Slope Estimator

MW-6



n = 13

Slope = 2.137 units per year.

Mann-Kendall statistic = 75
critical = 39

Increasing trend significant at 98% confidence level (α = 0.01 per tail).

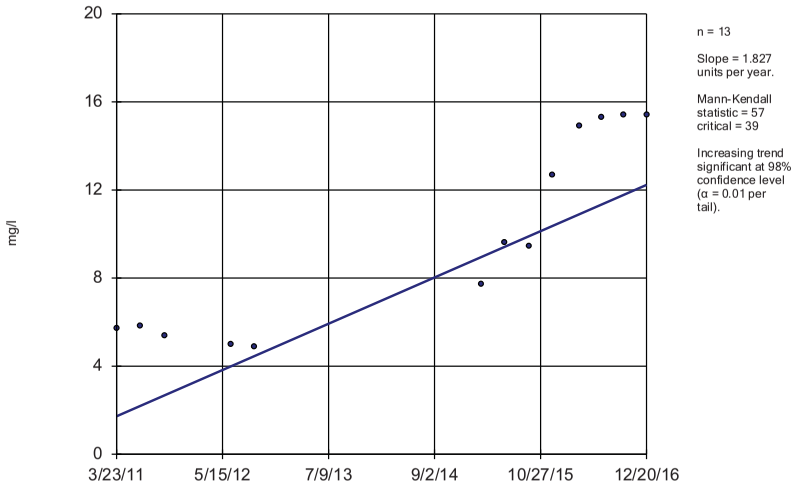
Constituent: Chloride Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-633D



n = 13

Slope = 1.827 units per year.

Mann-Kendall statistic = 57
critical = 39

Increasing trend significant at 98% confidence level (α = 0.01 per tail).

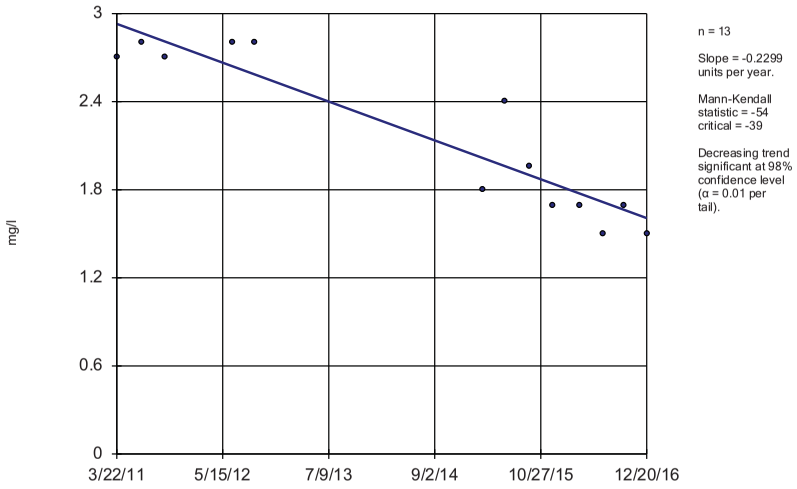
Constituent: Chloride Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-7



n = 13

Slope = -0.2299 units per year.

Mann-Kendall statistic = -54
critical = -39

Decreasing trend significant at 98% confidence level (α = 0.01 per tail).

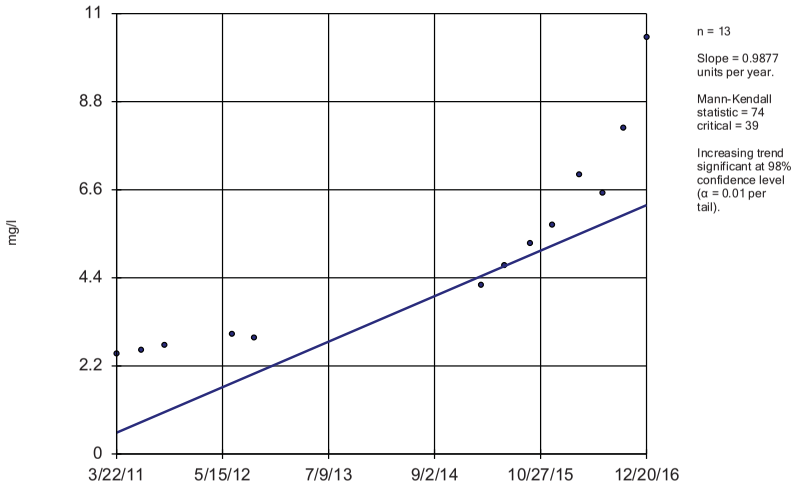
Constituent: Chloride Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

NAB-2



n = 13

Slope = 0.9877 units per year.

Mann-Kendall statistic = 74
critical = 39

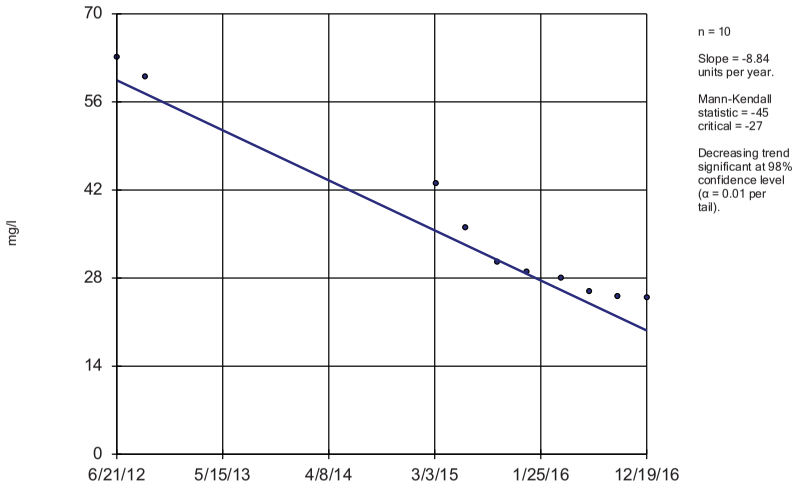
Increasing trend significant at 98% confidence level (α = 0.01 per tail).

Constituent: Chloride Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sen's Slope Estimator

NE-2



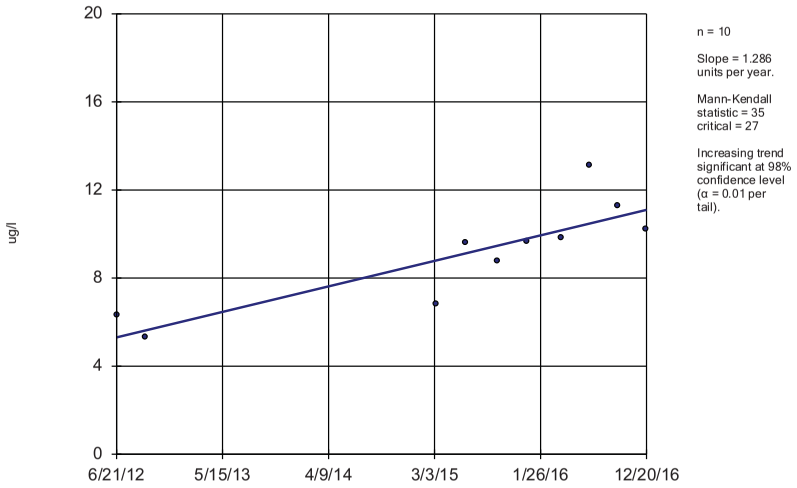
Constituent: Chloride Analysis Run 2/28/2017 10:03 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-1R



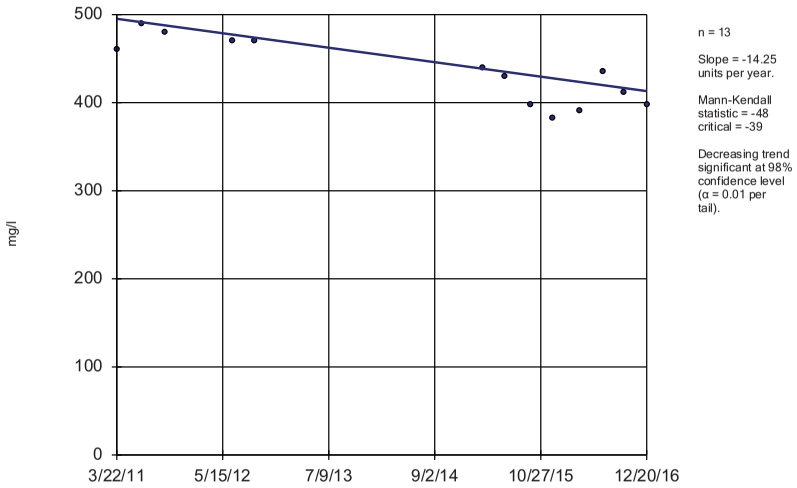
Constituent: cis-1,2-Dichloroethene Analysis Run 2/28/2017 10:04 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-577



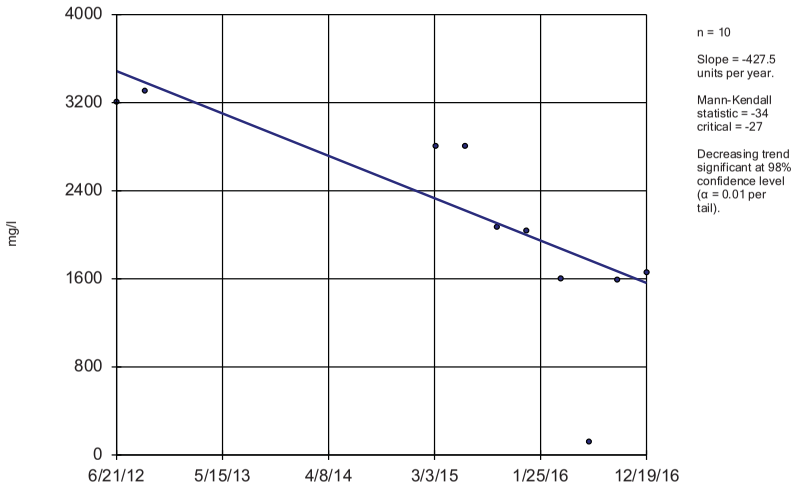
Constituent: Dissolved Solids Analysis Run 2/28/2017 10:04 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

NE-2



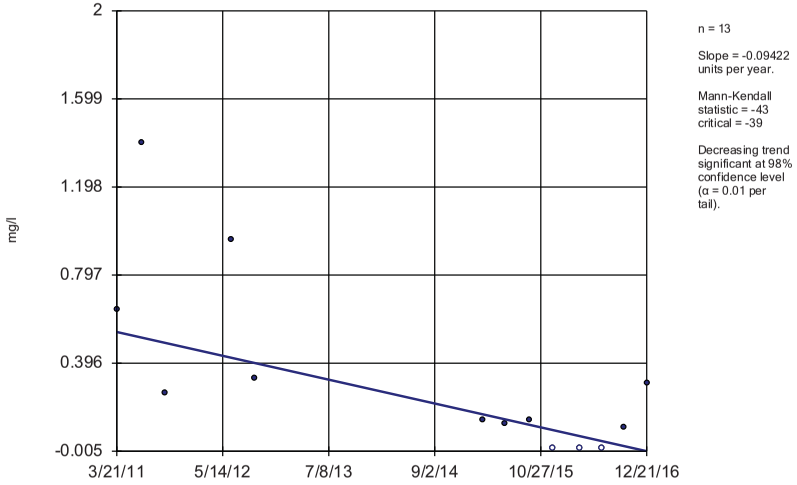
Constituent: Dissolved Solids Analysis Run 2/28/2017 10:04 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG
Hollow symbols indicate censored values.

Sen's Slope Estimator

CAO-2



n = 13

Slope = -0.09422 units per year.

Mann-Kendall statistic = -43
critical = -39

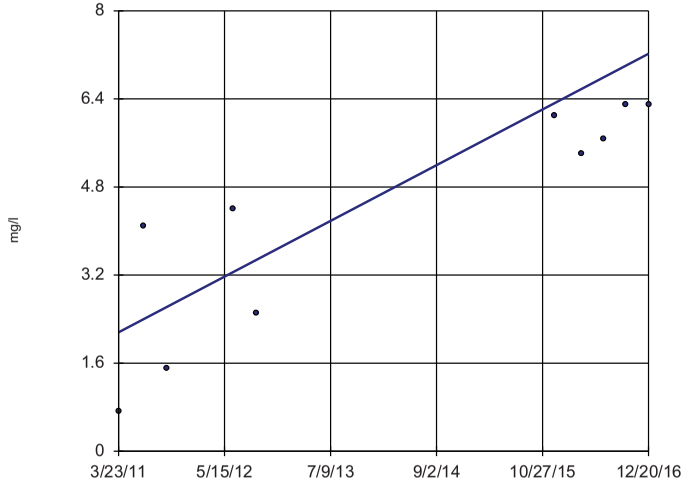
Decreasing trend significant at 98% confidence level (α = 0.01 per tail).

Constituent: Iron Analysis Run 2/28/2017 10:04 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-1



n = 10

Slope = 0.8799 units per year.

Mann-Kendall statistic = 34
critical = 27

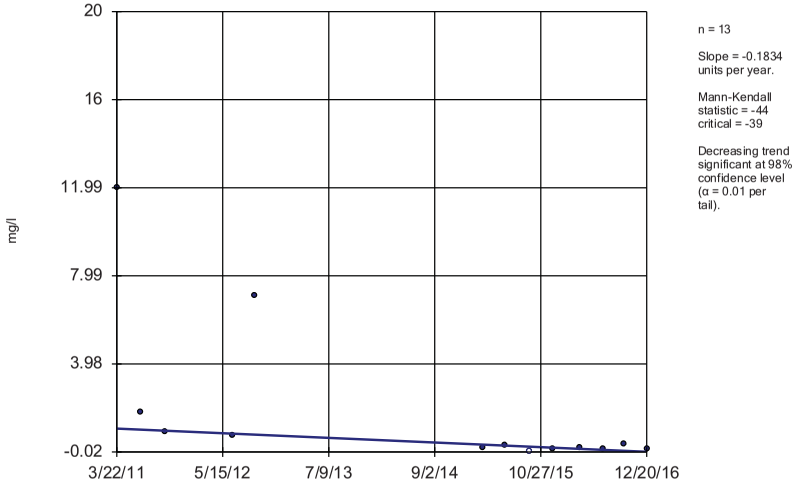
Increasing trend significant at 98% confidence level (α = 0.01 per tail).

Constituent: Iron Analysis Run 2/28/2017 10:04 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG
Hollow symbols indicate censored values.

Sen's Slope Estimator

MW-689D



n = 13

Slope = -0.1834 units per year.

Mann-Kendall statistic = -44
critical = -39

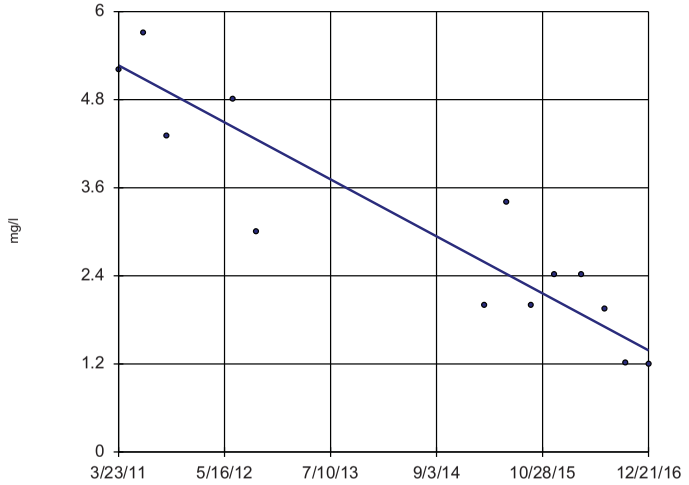
Decreasing trend significant at 98% confidence level (α = 0.01 per tail).

Constituent: Iron Analysis Run 2/28/2017 10:04 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

CAO-1



n = 13

Slope = -0.6746 units per year.

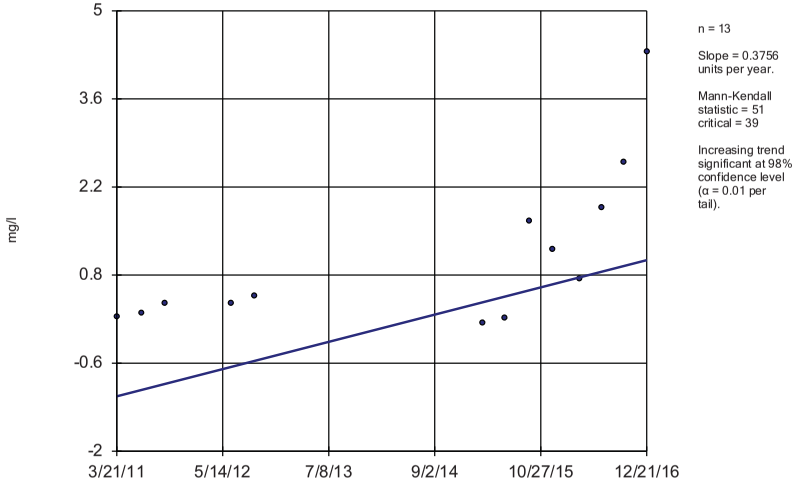
Mann-Kendall statistic = -60
critical = -39

Decreasing trend significant at 98% confidence level (α = 0.01 per tail).

Constituent: Manganese Analysis Run 2/28/2017 10:04 AM
NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sen's Slope Estimator

CAO-2



n = 13

Slope = 0.3756 units per year.

Mann-Kendall statistic = 51
critical = 39

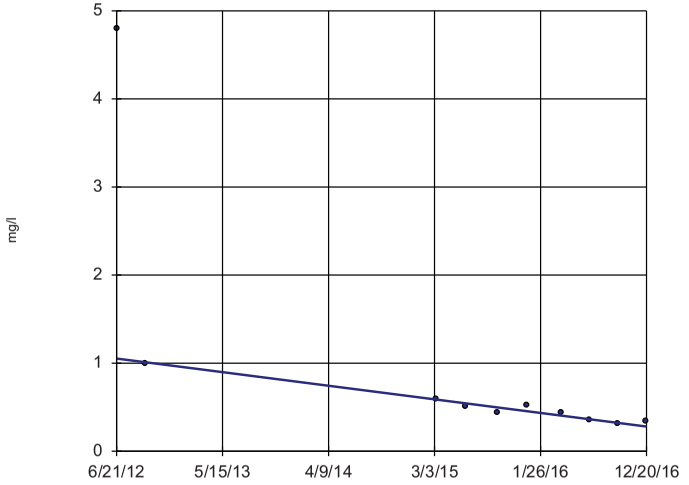
Increasing trend significant at 98% confidence level (α = 0.01 per tail).

Constituent: Manganese Analysis Run 2/28/2017 10:04 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sen's Slope Estimator

MW-1R



n = 10

Slope = -0.1715 units per year.

Mann-Kendall statistic = -39
critical = -27

Decreasing trend significant at 98% confidence level (α = 0.01 per tail).

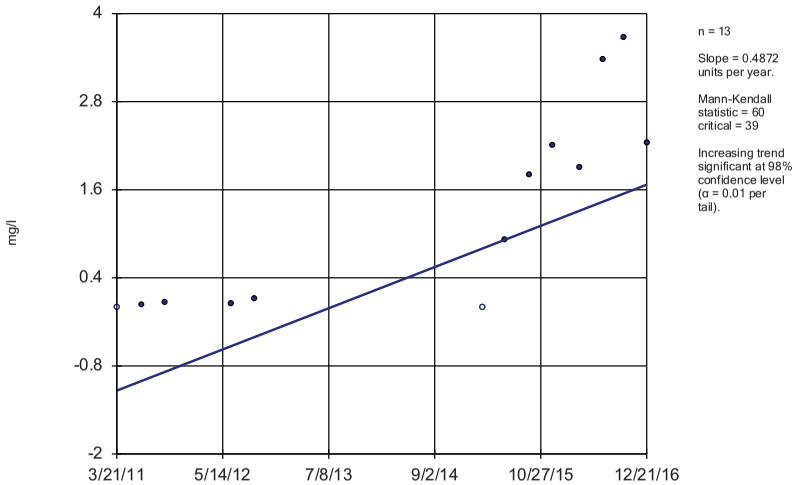
Constituent: Manganese Analysis Run 2/28/2017 10:04 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Hollow symbols indicate censored values.

Sen's Slope Estimator

MW-3



n = 13

Slope = 0.4872 units per year.

Mann-Kendall statistic = 60
critical = 39

Increasing trend significant at 98% confidence level (α = 0.01 per tail).

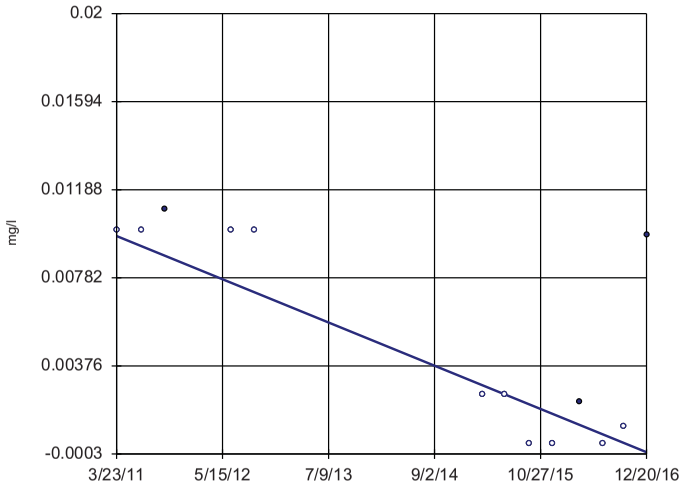
Constituent: Manganese Analysis Run 2/28/2017 10:04 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Hollow symbols indicate censored values.

Sen's Slope Estimator

MW-2



n = 13

Slope = -0.001732 units per year.

Mann-Kendall statistic = -40
critical = -39

Decreasing trend significant at 98% confidence level (α = 0.01 per tail).

Constituent: Nickel Analysis Run 2/28/2017 10:04 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

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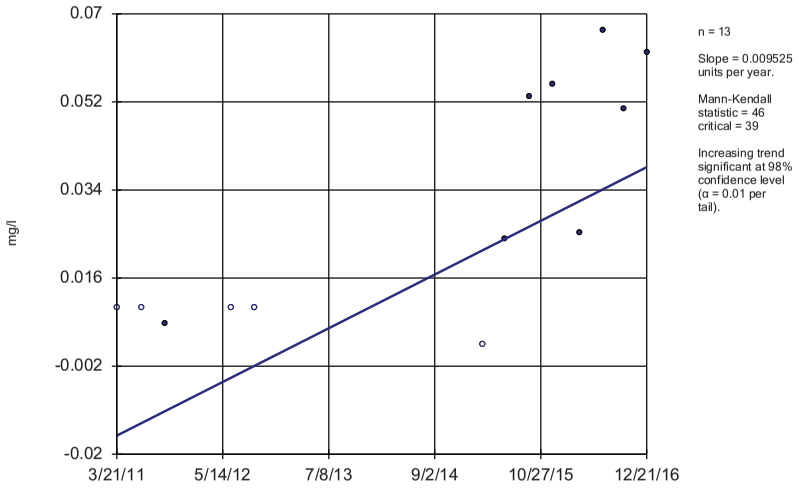
Hollow symbols indicate censored values.

Sen's Slope Estimator

MW-3



n = 13

Slope = 0.009525 units per year.

Mann-Kendall statistic = 46
critical = 39

Increasing trend significant at 98% confidence level (α = 0.01 per tail).

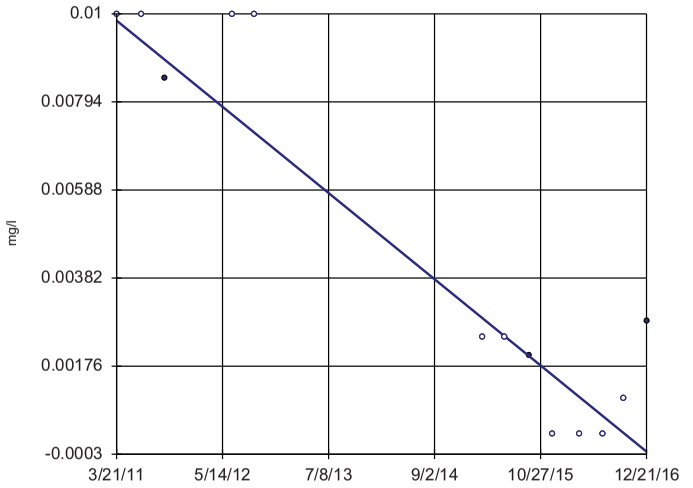
Constituent: Nickel Analysis Run 2/28/2017 10:04 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG
Hollow symbols indicate censored values.

Sen's Slope Estimator

MW-5



n = 13

Slope = -0.00175 units per year.

Mann-Kendall statistic = -44
critical = -39

Decreasing trend significant at 98% confidence level (α = 0.01 per tail).

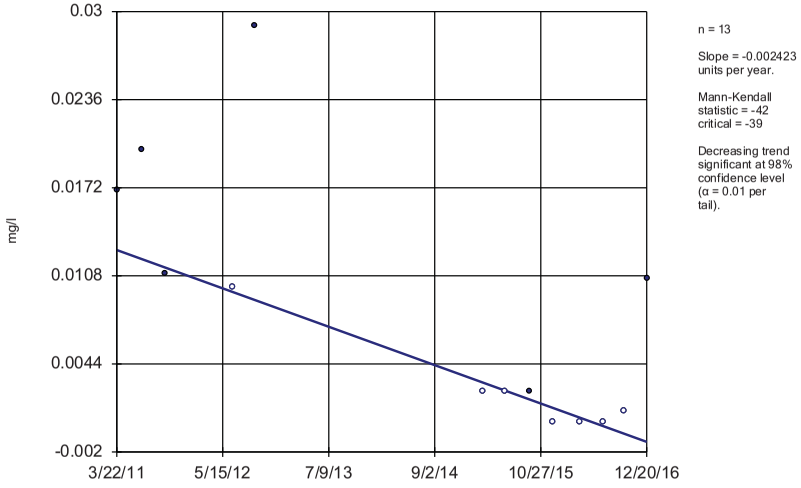
Constituent: Nickel Analysis Run 2/28/2017 10:04 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG
Hollow symbols indicate censored values.

Sen's Slope Estimator

MW-7



n = 13

Slope = -0.002423 units per year.

Mann-Kendall statistic = -42
critical = -39

Decreasing trend significant at 98% confidence level (α = 0.01 per tail).

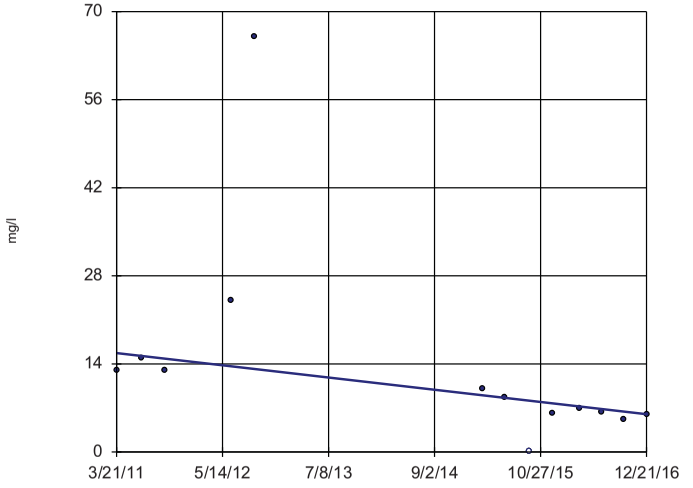
Constituent: Nickel Analysis Run 2/28/2017 10:04 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG
Hollow symbols indicate censored values.

Sen's Slope Estimator

CAO-3



n = 13

Slope = -1.687 units per year.

Mann-Kendall statistic = -45
critical = -39

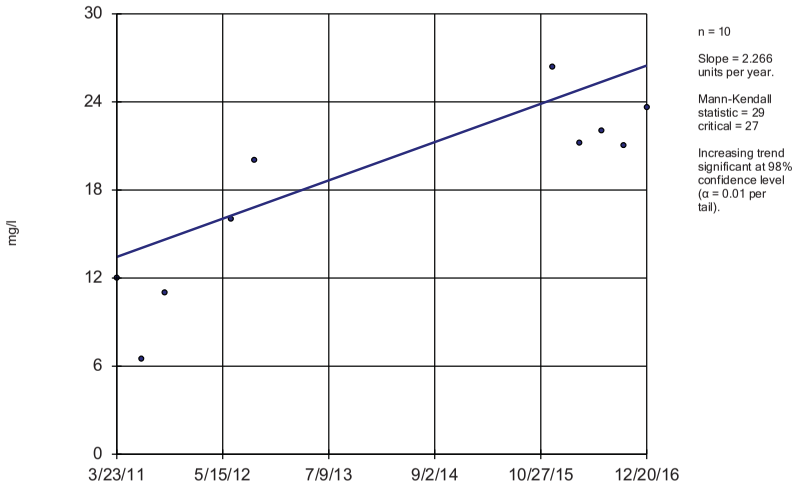
Decreasing trend significant at 98% confidence level (α = 0.01 per tail).

Constituent: Sulfate Analysis Run 2/28/2017 10:04 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sen's Slope Estimator

MW-1



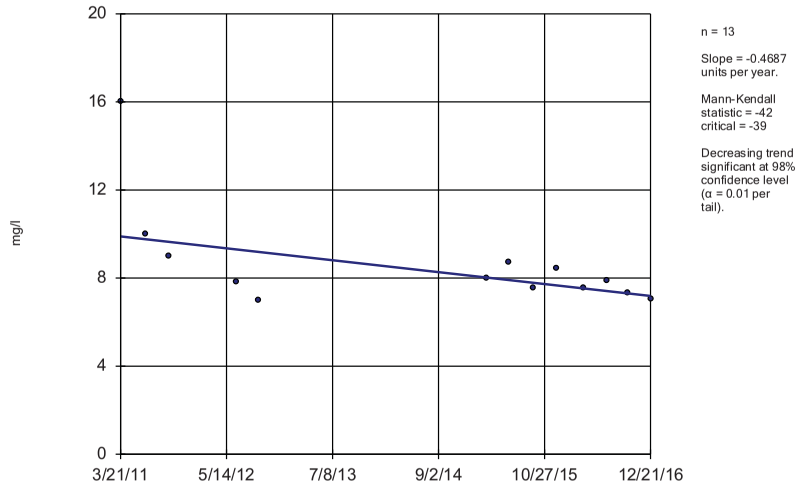
Constituent: Sulfate Analysis Run 2/28/2017 10:04 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-5



Constituent: Sulfate Analysis Run 2/28/2017 10:04 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

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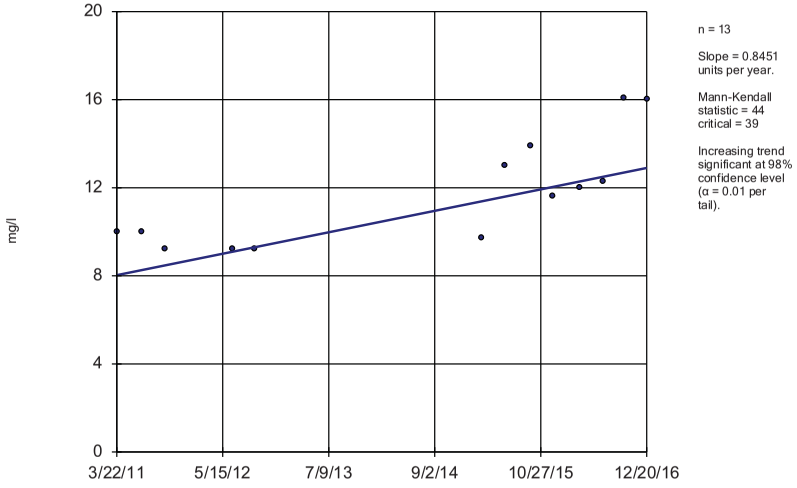
Sen's Slope Estimator

MW-509D



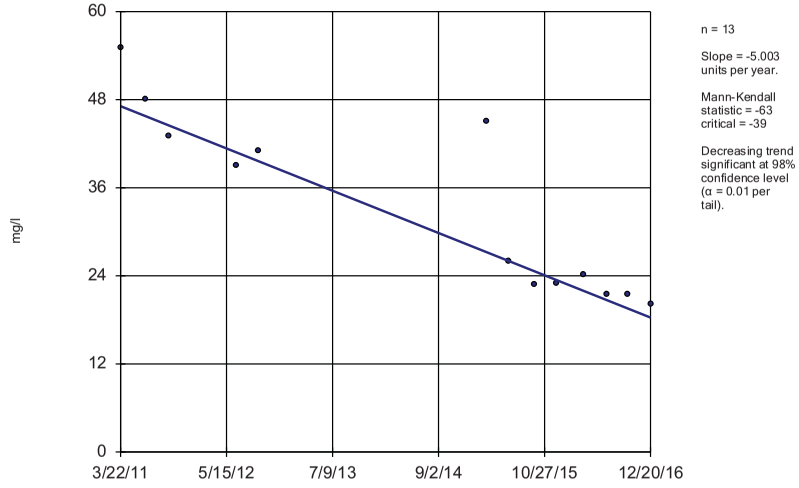
Constituent: Sulfate Analysis Run 2/28/2017 10:05 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-577

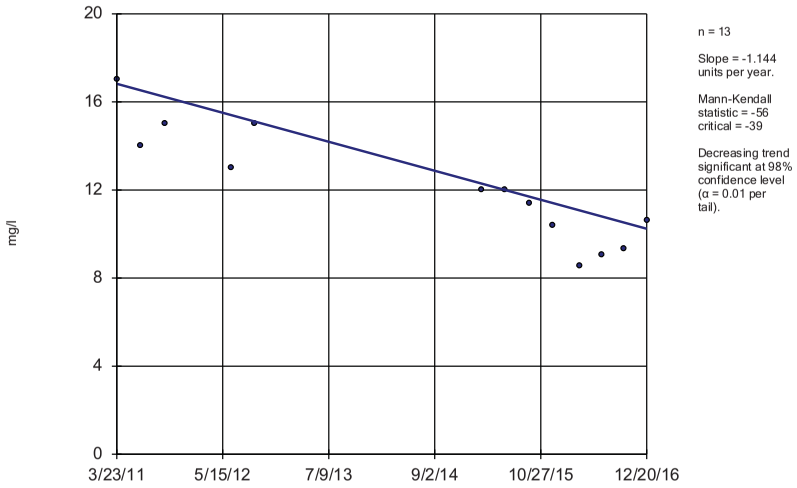


Constituent: Sulfate Analysis Run 2/28/2017 10:05 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sen's Slope Estimator

MW-633D



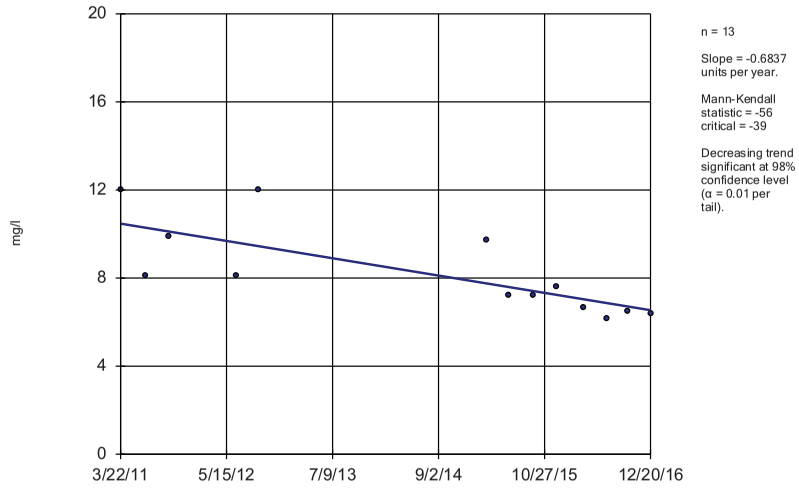
Constituent: Sulfate Analysis Run 2/28/2017 10:05 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-7



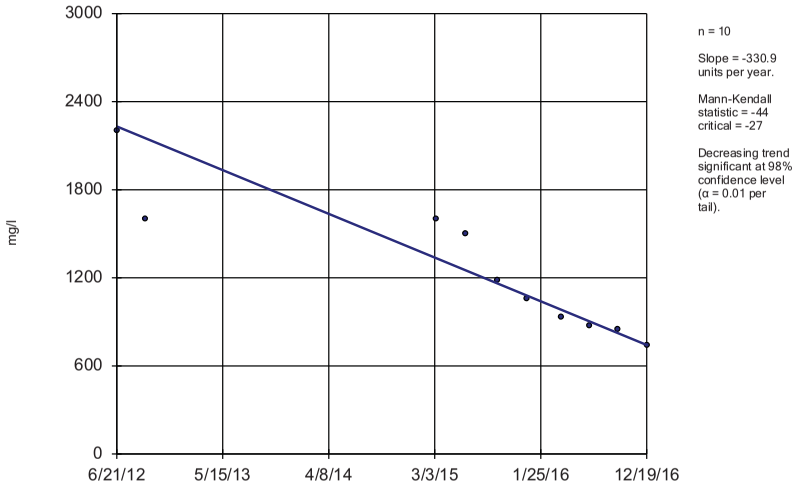
Constituent: Sulfate Analysis Run 2/28/2017 10:05 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

NE-2



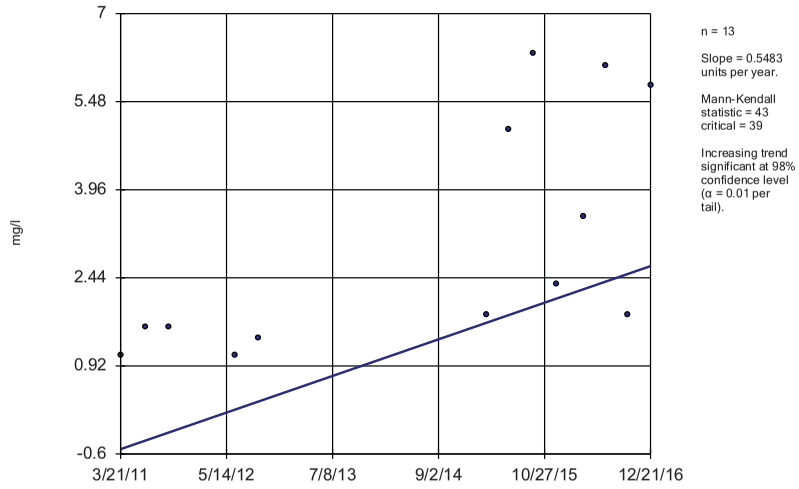
Constituent: Sulfate Analysis Run 2/28/2017 10:05 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

CAO-2

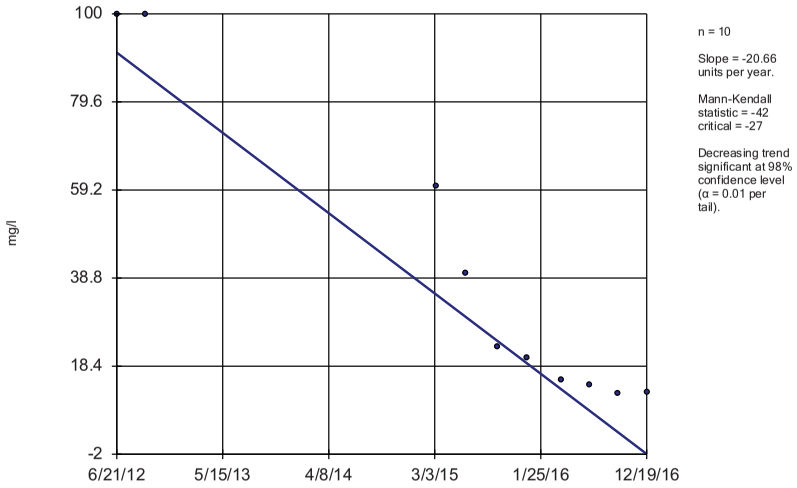


Constituent: TOC [Total Organic Carbon] Analysis Run 2/28/2017 10:05 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sen's Slope Estimator

NE-2



n = 10

Slope = -20.66 units per year.

Mann-Kendall statistic = -42
critical = -27

Decreasing trend significant at 98% confidence level (α = 0.01 per tail).

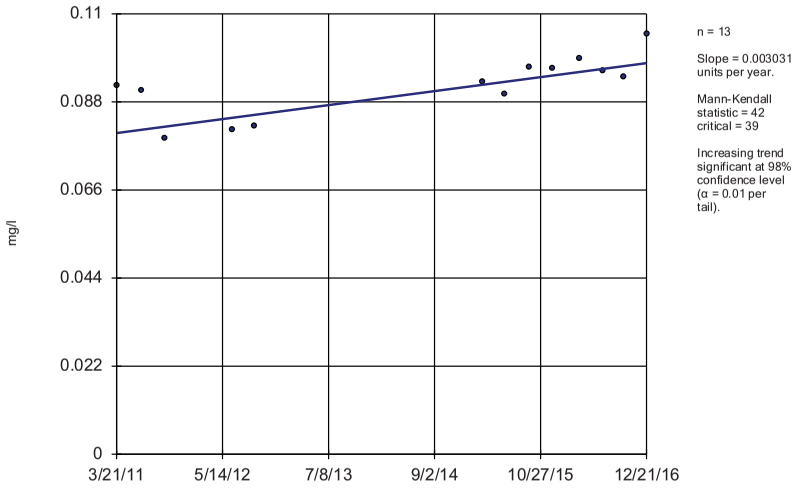
Constituent: TOC [Total Organic Carbon] Analysis Run 2/28/2017 10:05 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-4



n = 13

Slope = 0.003031 units per year.

Mann-Kendall statistic = 42
critical = 39

Increasing trend significant at 98% confidence level (α = 0.01 per tail).

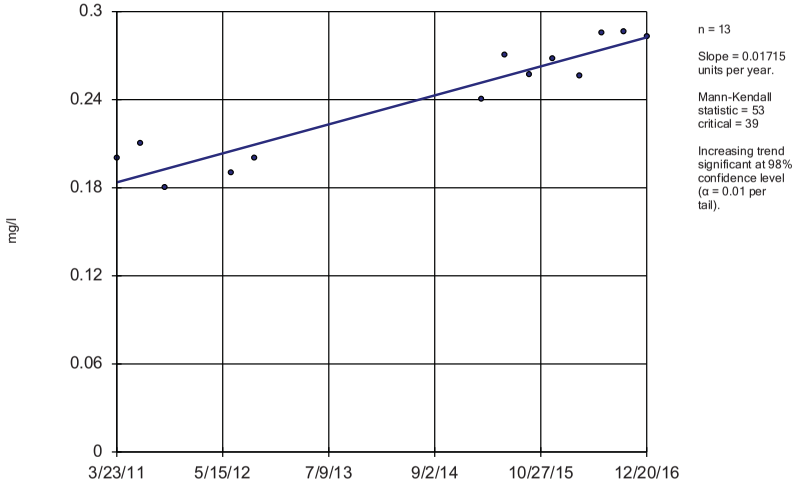
Constituent: Zinc Analysis Run 2/28/2017 10:05 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

MW-633D



n = 13

Slope = 0.01715 units per year.

Mann-Kendall statistic = 53
critical = 39

Increasing trend significant at 98% confidence level (α = 0.01 per tail).

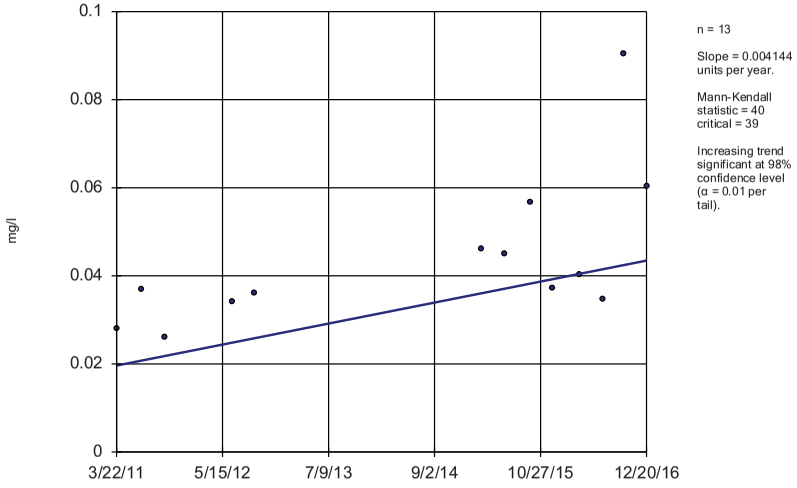
Constituent: Zinc Analysis Run 2/28/2017 10:05 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sanitas™ v.9.5.30 Software licensed to SCS Engineers. UG

Sen's Slope Estimator

NAB-2



n = 13

Slope = 0.004144 units per year.

Mann-Kendall statistic = 40
critical = 39

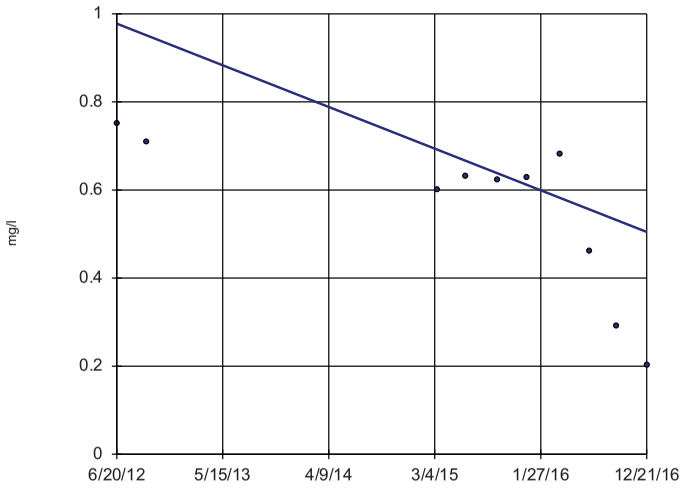
Increasing trend significant at 98% confidence level (α = 0.01 per tail).

Constituent: Zinc Analysis Run 2/28/2017 10:05 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix

Sen's Slope Estimator

NE-6



n = 10

Slope = -0.1049 units per year.

Mann-Kendall statistic = -29
critical = -27

Decreasing trend significant at 98% confidence level (α = 0.01 per tail).

Constituent: Zinc Analysis Run 2/28/2017 10:05 AM

NABORS Client: SCS Engineers Data: NABORS_DATABASE_SanitasMatrix