3R – 467 2014 GWMR 04 / 16 / 2015

John F. (Rick) Greiner, CPG, P.G.

ConocoPhillips Company Risk Management & Remediation Program Manager/Director Corp. Waste Management Program 600 N. Dairy Ashford, MA 1004 Houston, TX 77079 Phone: 281-293-3264 E-mail: Rick.Greiner@conocophillips.com



Mr. Glenn von Gonten New Mexico Oil Conservation Division 1220 South St. Francis DR Santa Fe, NM 87505

April 16, 2015

Re: API No. 30-045-10923 2014 Annual Groundwater Monitoring Report

Dear Mr. von Gonten:

Enclosed is the 2014 Annual Groundwater Monitoring Report for the Marcotte No. 1 site. This report, prepared by Conestoga-Rovers & Associates (CRA), contains the results of groundwater monitoring activities conducted during 2014.

Please let me know if you have any questions.

Sinters Rick Greiner

Enc



www.CRAworld.com



Final Report

2014 GROUNDWATER SAMPLING REPORT

ConocoPhillips Company Marcotte No. 1 San Juan County, New Mexico API# 30-045-10923

Prepared for: ConocoPhillips Company

Conestoga-Rovers & Associates

6121 Indian School Road, NE Suite 200 Albuquerque, New Mexico 87110



September 2014 • 085692 • Report No. 1

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Section 1.0 Introduction

Conestoga-Rovers & Associates (CRA) conducted groundwater monitoring at the ConocoPhillips Company (ConocoPhillips) Marcotte No. 1 natural gas well (Site) on April 2, 2014. The Site is located within Unit Letter G, Section 8, Township 31N, Range 10W, off County Road 2391, in San Juan County, New Mexico (latitude: 36.915560° N; longitude: -107.901902° W) (**Figure 1**). This report summarizes the monitoring activities and groundwater data that were collected during this single monitoring event.

1.1 Site History

The Site is located on private land that is leased by ConocoPhillips. The natural gas well is currently operated by Burlington Resources Oil and Gas Company LP, a wholly owned subsidiary of ConocoPhillips. A Site detail map is included as **Figure 2**.

Hydrocarbon impacted soil was discovered at the Site in 2003 during excavation work to reset the production equipment. Approximately 3,000 cubic yards of impacted soil was removed from the former pit in September 2003. Soil was landfarmed on the adjacent Marcotte No. 2 site with approval from both the New Mexico Oil Conservation Division (NMOCD) and U.S. Bureau of Land Management.

Impacted soils were excavated to approximately 30 feet below ground surface (ft bgs). Groundwater was encountered at a depth of 30 ft bgs. Soil impacts were noted at the time to end in a black stained gravel layer about 6 inches above the groundwater. The horizontal extent of contamination was determined by the limits of the open excavation on the north, northeast and east and by four soil borings located northwest, west and southwest.

Observation of the water in the bottom of the open excavation showed minor free phase hydrocarbons. Water and oil was removed from the excavation using a pump truck over a period of 2 months. Prior to backfilling the excavation no free phase hydrocarbons were visible on the water surface.

Two of the original soil borings were converted into groundwater monitoring wells (MW-2 and MW-3) in 2003 and an additional monitor well, MW-1, installed through the center of the excavation, was installed in September 2004 (**Figure 2**). Monitor Wells MW-2 and MW-3 were completed with above-ground (3 ft tall) well shrouds and MW-1 was installed with an at-grade concrete monument. Monitor Wells MW-2 and MW-3 were first gauged for depth to water (**Table 1**) and sampled in October 2003. Monitor Well MW-1 was first gauged and sampled in September 2004. Samples collected from Site monitor wells were analyzed for benzene, toluene, ethylbenzene and xylenes (BTEX) and for dissolved metals, anions and general chemistry parameters. Concentrations of BTEX constituents were below the NM Water Quality Control Commission (NMWQCC) standards for these compounds. Concentrations of dissolved manganese were found at a concentration above NMWQCC standards in groundwater of MW-2 and sulfates were above standards in all wells from the 2003/2004 background samples.



Monitor Wells MW-2 and MW-3 were sampled again in December 2003 and quarterly in 2004. It is uncertain why MW-1 was not sampled. It is possible that MW-1 was not able to be found due to its location in an area of brush and it's at-grade surface completion.

Site wells were not sampled again until December 9, 2010. Monitor Wells MW-2 and MW-3 were gauged and sample collected for analysis of BTEX, dissolved metals, anions and general chemistry parameters. MW-1 was noted as not located. BTEX constituents were not detected at any concentrations above laboratory detection limits. Total dissolved solids (TDS) and chlorides were detected in MW-2 and MW-3 during the December 2010 sampling event at concentrations above the NMWQCC standards.

1.2 Site Setting

The Site is located in San Juan County, New Mexico, on privately owned land. The elevation at the Site is approximately 5,850 feet above mean sea level. An ephemeral wash borders the location approximately 30 yards to the north. A seasonal irrigation ditch is located approximately 100 yards to the west. Subsurface soils at the location are mainly fine to coarse sands with trace amounts of cobbles and boulders. Groundwater flows generally parallel to the adjacent wash in a west-southwest direction towards the Animas River located ½ mile west.

Subsurface soils at the Site consist of very fine to medium grained sands with trace gravels and cobbles and with minor silty-clay lenses. Groundwater was measured at a depth of 31.85 ft-below the top of PVC casing (TOC) during the April 2014 event. The groundwater gradient is presumed essentially parallel to the adjacent wash in a westerly direction, towards the Animas River located.

Section 2.0 Groundwater Monitoring Summary

A groundwater sampling event was conducted at the Site on April 2, 2014. Monitor well MW-1 was found and sampled during this event (see **Figure 2**). Prior to sampling, depth to groundwater was measured in each well using an oil/water interface probe (**Table 1**). Monitor Well MW-2 was the only Site well with measurable groundwater. Groundwater was measured in MW-2 at a depth of 31.85 ft below TOC. The measured dry depth of MW-1 (23.20 ft below TOC), is lower than the measured total depths of MW-2 and MW-3 (37.40 and 38.45 ft below TOC, respectively). CRA believes this indicates that MW-1 has an obstruction inside the PVC casing. A groundwater potentiometric surface map for the site was not able to be generated due to the single groundwater elevation measurement

The groundwater flow at the site is presumed to be aligned with the flow of the adjacent Miller Canyon Wash, towards the Animas River, located ½ mile to the west.



2.1 Groundwater Monitoring Methodology

During monitoring events, at least three well volumes were purged from Site Monitor Well MW-2 with a hand-held disposable bailer. Field parameters of pH, temperature, conductivity, oxidation/reduction potential, and dissolved oxygen were collected during purging of groundwater prior to sampling. The field parameter readings are presented in **Table 2**. Purge water generated during purging of Site monitor wells was placed in the on-Site produced water tank. The groundwater sample was placed in laboratory prepared bottles, packed on ice, and shipped under chain-of-custody documentation to Pace Analytical Services, Inc. of Lenexa, KS.

The MW-2 groundwater sample was analyzed for the presence of BTEX by EPA method 8260, semivolatile compounds by EPA method 8270, gasoline and diesel-range organics by EPA 8015, for dissolved metals by EPA method 6010, for dissolved mercury by EPA method 7470, and for general chemistry parameters by various EPA methods. A summary of analytical results is presented in **Table 3**. Groundwater laboratory analytical results are presented in **Appendix A**.

2.2 Groundwater Monitoring Analytical Results

The NMWQCC mandates that groundwater quality in New Mexico be protected, and has issued groundwater quality standards in Title 20, Chapter 6, Part 2, Section 3103 of the New Mexico Administrative Code (20.6.2.3103 NMAC). Groundwater quality standards have been set for the protection of human health, domestic water supply, and irrigation use.

A groundwater inorganic concentration map is included as **Figure 3**. Groundwater analytical results are discussed below.

<u>April 2014</u>

- **BTEX:** BTEX constituents were not detected in concentrations above the laboratory detection limits.
- **Dissolved Manganese:** The NMWQCC domestic water supply groundwater quality standard for dissolved manganese is 0.2 mg/L. Monitoring Well MW-2 exceeded this standard with an analytical result of 0.853 mg/L.
- **Sulfate:** The NMWQCC domestic water supply groundwater quality standard for sulfate is 600 mg/L. Monitoring Well MW-2 exceeded this standard with an analytical result of 2,360mg/L.
- **TDS:** The NMWQCC domestic water supply groundwater quality standard for TDS is 1,000 mg/L. Monitoring Well MW-2 exceeded this standard with an analytical result of 3,030 mg/L.



Section 3.0 Water Well Database Search

A list of domestic water well users and well coordinates was generated from the New Mexico Office of the State Engineer Water Rights Reporting System database. The list of wells, including the owner and well coordinates, is presented in **Appendix B.** The wells have been plotted on a map and are presented as **Figure 4**.

Section 4.0 Conclusions and Recommendations

BTEX constituents in groundwater at the Site have not been detected in any concentrations above laboratory detection limits since groundwater monitoring began. Chloride was detected at concentrations above the NMWQCC standard in the past in Monitor Wells MW-2 and MW-3; however, the April 2, 2014 sampling event showed that chloride in MW-2 was present at a concentration well below the standard. Sulfate has been detected in the past at concentrations above the NMWQCC standard in Site wells and remains above the standard in Monitor Well MW-2. TDS has also been detected in MW-2 and MW-3 at concentrations above the NMWQCC standard.

CRA recommends:

- The plugging and abandonment of Monitor Wells MW-1 and MW-3;
- Reinstalling MW-1 and MW-3 deeper to intercept the Site groundwater; and
- Install an up-gradient well to assess background concentrations.

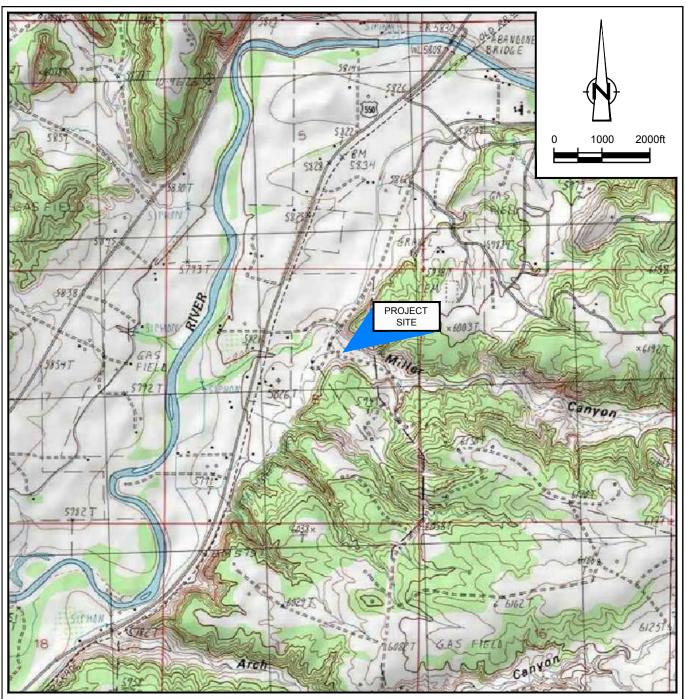
Once these wells are reinstalled and a top of well casing survey completed, a Site-specific groundwater flow direction and gradient can be established.

Until a better understanding of Site groundwater conditions are enabled by the installation and sampling of replacement wells, CRA does not recommend further groundwater monitoring.



Figures





SOURCE: USGS 7.5 MINUTE QUAD "CEDAR HILL, NEW MEXICO"

LAT/LONG: 36.9155° NORTH, 107.9019° WEST COORDINATE: NAD83 DATUM, U.S. FOOT STATE PLANE ZONE - NEW MEXICO WEST

Figure 1

SITE VICINITY MAP MARCOTTE No.1 NATURAL GAS WELL SITE SECTION 8, T31N-R10W, SAN JUAN COUNTY, NEW MEXICO *ConocoPhillips Company*

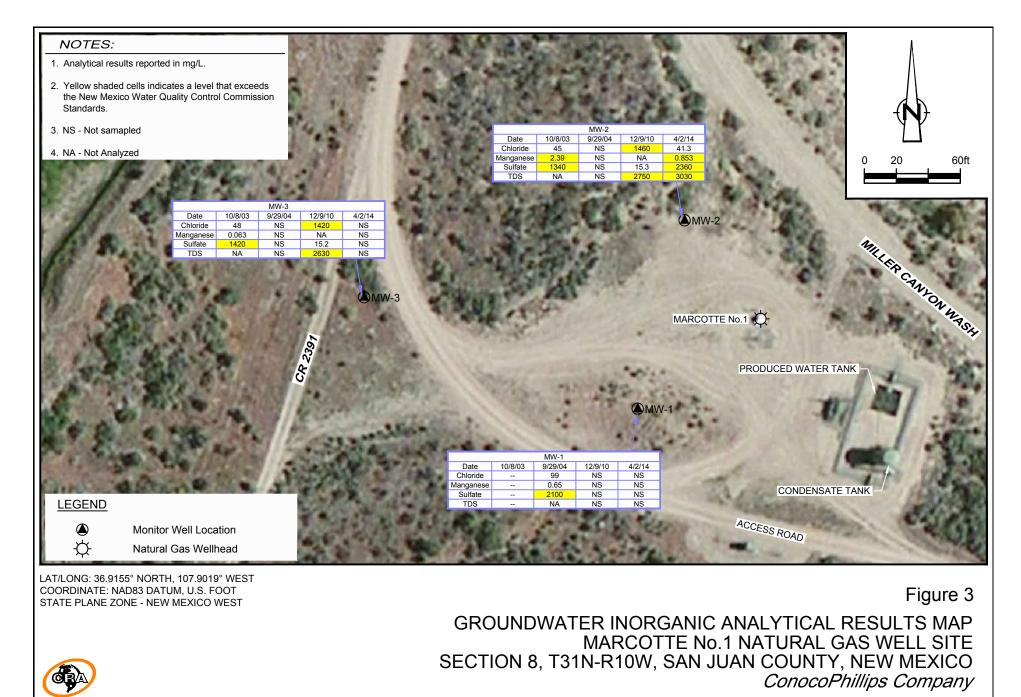


Figure 2

SITE PLAN MARCOTTE No.1 NATURAL GAS WELL SITE SECTION 8, T31N-R10W, SAN JUAN COUNTY, NEW MEXICO ConocoPhillips Company

085692-00(000)GN-DL001 JUL 16/2014

STATE PLANE ZONE - NEW MEXICO WEST



085602 00/000\CN DL001 UU

085692-00(000)GN-DL001 JUL 30/2014



LAT/LONG: 36.9155° NORTH, 107.9019° WEST COORDINATE: NAD83 DATUM, U.S. FOOT STATE PLANE ZONE - NEW MEXICO WEST

Figure 4

VICINITY WATER WELL LOCATION MAP MARCOTTE No.1 NATURAL GAS WELL SITE SECTION 8, T31N-R10W, SAN JUAN COUNTY, NEW MEXICO *ConocoPhillips Company*

085692-00(000)GN-DL001 JUL 30/2014



Monitor Well Specifications and Groundwater Elevations ConocoPhillips Company Marcotte No. 1 San Juan County, New Mexico

Well ID	Total Depth (ft below TOC)	Screen Interval (ft bgs)	Date Measured	Depth to Groundwater (ft below TOC)
			9/29/2004	23.20
N 4147 1	22.20*	TT1	12/13/2004	23.67
MW-1	23.20*	Unknown	4/2/2014	DRY
			10/6/2003	29.71
			12/16/2003	30.09
			3/15/2004	30.62
			6/21/2004	30.05
MW-2	37.40	22-37 ft bgs	9/29/2004	
			12/13/2004	29.88
			12/9/2010	29.78
			4/2/2014	31.85
			10/6/2003	30.74
			12/16/2003	34.14
			3/15/2004	
			6/21/2004	36.62
MW-3	38.45	23-38 ft bgs	9/29/2004	28.72
			12/13/2004	32.35
			12/9/2010	35.51
			4/2/2014	DRY

Notes:

bgs = Below ground surface

ft = Feet

TOC = Top of casing

*Total depth measured 4/2/2014-may represent an obstruction; well completion data unavailable.

Field Parameters ConocoPhillips Company Marcotte No. 1 San Juan County, NM

Well ID	Sample Date	Temperature (°C)	рН	TDS (g/L)	Conductivity (μS/cm)	DO (mg/L)	ORP (mV)	Volume (gallons)
	4/2/2014	14.37	6.84	2.189	3367	1.36	70.4	1.5
	4/2/2014	14.32	6.90	2.183	3361	1.39	55.0	2.5
MW-2	4/2/2014	14.41	6.91	2.184	3359	1.56	49.3	2.75
	4/2/2014	14.42	6.92	2.180	3354	1.58	44.2	3.0

Groundwater Analytical Results Summary ConocoPhillips Company Marcotte No. 1 San Juan County, New Mexico

Well ID	Sample ID	Date	Sample Type	Benzene (mg/L)	Toluene (mg/L)	Ethylbenzene (mg/L)	Xylenes (mg/L)	Total Hardness, dissolved (mg/L)	Arsenic, dissolved (mg/L)	Barium, dissolved (mg/L)	Cadmium, dissolved (mg/L)	Calcium, dissolved (mg/L)	Chromium, dissolved (mg/L)	Copper, dissolved (mg/L)	Iron, dissolved (mg/L)
	NMWQCC Groundwater Qua	ality Standards		0.01	0.75	0.75	0.62		0.1	1.0	0.01		0.05	1.0	1.0
MW-1	MW-1	9/29/2004	0	< 0.0003	< 0.0002	0.038	0.0369		< 0.001	0.017	0.0009	286	0.0003	0.001	0.19
10100-1	MW-1	12/13/2004	(orig)	0.0004	0.0007	0.0007	0.0202								
	M P Unit 1 MW-2	10/8/2003	(orig)	< 0.0003	< 0.0002	< 0.0002	< 0.0002		0.0036	0.047	< 0.0001	266	0.0008	0.0021	0.98
	MW-2	12/16/2003	(orig)	0.0004	< 0.0002	< 0.0002	< 0.0002								
	MW-2	3/15/2004	(orig)	0.0004	0.0003	< 0.0002	0.0002								
	MW-2	6/21/2004	(orig)	< 0.0003	< 0.0002	< 0.0002	< 0.0002								
MW-2	MW-2	9/29/2004	(orig)	< 0.0003	0.0003	0.0003	0.0007								
	MW-2	12/13/2004	(orig)	0.0003	0.0013	0.0003	0.0112								
	MW-2	12/9/2010	(orig)	< 0.001	< 0.001	< 0.001	< 0.001	1100	0.003	0.009	< 0.001	360	< 0.001		0.042
	GW-085692-040214-CM-MW-2	4/2/2014	(orig)	< 0.001	< 0.001	< 0.001	< 0.003	1180	0.0011	0.0128	< 0.50		<1.0	<1.0	
	M P Unit 1 MW-3	10/8/2003	(orig)	< 0.0003	0.0002	< 0.0002	< 0.002		0.0012	0.037	< 0.0001	262	0.0012	0.0017	0.47
	MW-3	12/16/2003	(orig)	0.0005	< 0.0002	< 0.0002	< 0.0002								
	MW-3	6/21/2004	(orig)	< 0.0003	< 0.002	< 0.0002	< 0.002								
MW-3	MW-3	9/29/2004	(orig)	< 0.0003	< 0.002	< 0.0002	< 0.002								
	MW-3	12/13/2004	(orig)	< 0.0003	0.0003	< 0.0002	0.0016								
	MW-3	12/9/2010	(orig)	< 0.001	< 0.001	< 0.001	< 0.001	1130	0.002	0.009	< 0.001	367	< 0.001		0.009

Notes:

BDL = below detection limit (actual laboratory detection limit not available)

mg/L = milligrams per liter (parts per million)

NA = Not Analyzed

NMWQCC = New Mexico Water Quality Control Commission

Groundwater Analytical Results Summary ConocoPhillips Company Marcotte No. 1 San Juan County, New Mexico

Well ID	Sample ID	Date	Sample Type	Magnesium, dissolved (mg/L)	Manganese, dissolved (mg/L)	Molybdenum, dissolved (mg/L)	Potassium, dissolved (mg/L)	Selenium (mg/L)	Silver (mg/L)	Sodium, dissolved (mg/L)	Zinc, dissolved (mg/L)	Alkalinity, total as CaCO3 (mg/L)	TDS (mg/L)	Chloride (mg/L)
	NMWQCC Groundwater Qua	ality Standards			0.2	1.0		0.05	0.05		10		1000	250
MW-1	MW-1	9/29/2004	0	39.9	0.65		2.5			727	< 0.02	318		99
10100-1	MW-1	12/13/2004	(orig)											
	M P Unit 1 MW-2	10/8/2003	(orig)	34.9	2.39		1.6			419	0.02	302		45
	MW-2	12/16/2003	(orig)											
	MW-2	3/15/2004	(orig)											
	MW-2	6/21/2004	(orig)											
MW-2	MW-2	9/29/2004	(orig)											
	MW-2	12/13/2004	(orig)											
	MW-2	12/9/2010	(orig)	50			6.56	0.005	0.031	603		410	2750	1460
	GW-085692-040214-CM-MW-2	4/2/2014	(orig)		0.853	0.0039						290	3030	41.3
	M P Unit 1 MW-3	10/8/2003	(orig)	34.5	0.063		1.6			409	< 0.01	291		48
	MW-3	12/16/2003	(orig)											
	MW-3	6/21/2004	(orig)											
MW-3	MW-3	9/29/2004	(orig)											
	MW-3	12/13/2004	(orig)											
	MW-3	12/9/2010	(orig)	50.9			4.28	0.027	0.031	550		370	2630	1420

Notes:

BDL = below detection limit (actual laboratory detection lin

mg/L = milligrams per liter (parts per million)

NA = Not Analyzed

NMWQCC = New Mexico Water Quality Control Commiss

Groundwater Analytical Results Summary ConocoPhillips Company Marcotte No. 1 San Juan County, New Mexico

Well ID	Sample ID	Date	Sample Type	Fluoride (mg/L)	Sulfate (mg/L)	Nitrate, NO3 as N (mg/L)	Orthophosphate, as P (mg/L)	Cyanide (mg/L)	pН
10	NMWQCC Groundwater Qua		турс						
		,		1.6	600	10		0.2	6 - 9
MW-1	MW-1	9/29/2004	0		2100				7.1
10100 1	MW-1	12/13/2004	(orig)						
	M P Unit 1 MW-2	10/8/2003	(orig)		1340				7.9
	MW-2	12/16/2003	(orig)						
	MW-2	3/15/2004	(orig)						
	MW-2	6/21/2004	(orig)						
MW-2	MW-2	9/29/2004	(orig)						
	MW-2	12/13/2004	(orig)						
	MW-2	12/9/2010	(orig)	BDL	15.3	6.36		0.003	6.71
	GW-085692-040214-CM-MW-2	4/2/2014	(orig)	0.68	2360	<0.10	0.10		7.3
	M P Unit 1 MW-3	10/8/2003	(orig)		1420				7.9
	MW-3	12/16/2003	(orig)						
	MW-3	6/21/2004	(orig)						
MW-3	MW-3	9/29/2004	(orig)						
	MW-3	12/13/2004	(orig)						
	MW-3	12/9/2010	(orig)	1.14	15.2	<0.10		0.002	6.92

Notes:

BDL = below detection limit (actual laboratory detection lin

mg/L = milligrams per liter (parts per million)

NA = Not Analyzed

NMWQCC = New Mexico Water Quality Control Commise

Appendix A

Groundwater Laboratory Analytical Report





Pace Analytical Services, Inc. 9608 Loiret Blvd. Lenexa, KS 66219 (913)599-5665

April 18, 2014

Jeff Walker COP Conestoga-Rovers & Associa 6121 Indian School Rd. NE Ste 200 Albuquerque, NM 87110

RE: Project: 085692 MARCOTTE NO 1 Pace Project No.: 60166332

Dear Jeff Walker:

Enclosed are the analytical results for sample(s) received by the laboratory on April 04, 2014. The results relate only to the samples included in this report. Results reported herein conform to the most current TNI standards and the laboratory's Quality Assurance Manual, where applicable, unless otherwise noted in the body of the report.

If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Alice Flanagan

Alice Flanagan alice.flanagan@pacelabs.com Project Manager

Enclosures

cc: Angela Bown, COP Conestoga-Rovers & Associa Christine Matthews, CRA





CERTIFICATIONS

Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Kansas Certification IDs

9608 Loiret Boulevard, Lenexa, KS 66219 WY STR Certification #: 2456.01 Arkansas Certification #: 13-012-0 Illinois Certification #: 003097 Iowa Certification #: 118 Kansas/NELAP Certification #: E-10116 Louisiana Certification #: 03055 Nevada Certification #: KS000212008A Oklahoma Certification #: 9205/9935 Texas Certification #: T104704407-13-4 Utah Certification #: KS000212013-3 Illinois Certification #: 003097



SAMPLE SUMMARY

Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Lab ID	Sample ID	Matrix	Date Collected	Date Received
60166332001	GW-085692-040214-CM-MW-2	Water	04/02/14 13:45	04/04/14 08:35
60166332002	TRIP BLANK	Water	04/02/14 13:45	04/04/14 08:35



SAMPLE ANALYTE COUNT

 Project:
 085692 MARCOTTE NO 1

 Pace Project No.:
 60166332

Lab ID Sample ID Method Analysts	Analytes Reported
60166332001 GW-085692-040214-CM-MW-2 EPA 8015B JDE	3
EPA 5030B/8015B JTK	3
EPA 6010 NDJ	1
EPA 6020 JGP	13
EPA 7470 TDS	1
EPA 8270 JMT	73
EPA 5030B/8260 PRG	69
EPA 120.1 NDL	1
SM 2320B JMC1	1
SM 2540C RAH	1
SM 4500-H+B NDL	1
EPA 300.0 OL	4
EPA 353.2 AJM	1



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Method: EPA 8015B

Description:8015B Diesel Range OrganicsClient:COP Conestoga-Rovers & Associates, Inc. NMDate:April 18, 2014

General Information:

1 sample was analyzed for EPA 8015B. All samples were received in acceptable condition with any exceptions noted below.

Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

Sample Preparation:

The samples were prepared in accordance with EPA 3510C with any exceptions noted below.

Initial Calibrations (including MS Tune as applicable):

All criteria were within method requirements with any exceptions noted below.

Continuing Calibration:

All criteria were within method requirements with any exceptions noted below.

Surrogates:

All surrogates were within QC limits with any exceptions noted below.

Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

QC Batch: GCSV/16583

A matrix spike/matrix spike duplicate was not performed due to insufficient sample volume.

Additional Comments:



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Method: EPA 5030B/8015B

Description:Gasoline Range OrganicsClient:COP Conestoga-Rovers & Associates, Inc. NMDate:April 18, 2014

General Information:

1 sample was analyzed for EPA 5030B/8015B. All samples were received in acceptable condition with any exceptions noted below.

Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

Initial Calibrations (including MS Tune as applicable): All criteria were within method requirements with any exceptions noted below.

Continuing Calibration:

All criteria were within method requirements with any exceptions noted below.

Internal Standards:

All internal standards were within QC limits with any exceptions noted below.

Surrogates:

All surrogates were within QC limits with any exceptions noted below.

Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

QC Batch: GCV/4730

A matrix spike/matrix spike duplicate was not performed due to insufficient sample volume.

Additional Comments:



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Method: EPA 6010

Description:6010 MET ICP, DissolvedClient:COP Conestoga-Rovers & Associates, Inc. NMDate:April 18, 2014

General Information:

1 sample was analyzed for EPA 6010. All samples were received in acceptable condition with any exceptions noted below.

Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

Sample Preparation:

The samples were prepared in accordance with EPA 3010 with any exceptions noted below.

Initial Calibrations (including MS Tune as applicable):

All criteria were within method requirements with any exceptions noted below.

Continuing Calibration:

All criteria were within method requirements with any exceptions noted below.

Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

Additional Comments:



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Method: EPA 6020

Description:6020 MET ICPMS, DissolvedClient:COP Conestoga-Rovers & Associates, Inc. NMDate:April 18, 2014

General Information:

1 sample was analyzed for EPA 6020. All samples were received in acceptable condition with any exceptions noted below.

Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

Sample Preparation:

The samples were prepared in accordance with EPA 3010 with any exceptions noted below.

Initial Calibrations (including MS Tune as applicable):

All criteria were within method requirements with any exceptions noted below.

Continuing Calibration:

All criteria were within method requirements with any exceptions noted below.

Internal Standards:

All internal standards were within QC limits with any exceptions noted below.

Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

QC Batch: MPRP/26827

A matrix spike and/or matrix spike duplicate (MS/MSD) were performed on the following sample(s): 60166332001

M1: Matrix spike recovery exceeded QC limits. Batch accepted based on laboratory control sample (LCS) recovery.

- MS (Lab ID: 1360510)
 - Manganese, Dissolved
 - Molybdenum, Dissolved

Additional Comments:



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Method: EPA 7470

Description:7470 Mercury, DissolvedClient:COP Conestoga-Rovers & Associates, Inc. NMDate:April 18, 2014

General Information:

1 sample was analyzed for EPA 7470. All samples were received in acceptable condition with any exceptions noted below.

Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

Sample Preparation:

The samples were prepared in accordance with EPA 7470 with any exceptions noted below.

Initial Calibrations (including MS Tune as applicable):

All criteria were within method requirements with any exceptions noted below.

Continuing Calibration:

All criteria were within method requirements with any exceptions noted below.

Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

Additional Comments:



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Method: EPA 8270

Description:8270 MSSV Semivolatile OrganicClient:COP Conestoga-Rovers & Associates, Inc. NMDate:April 18, 2014

General Information:

1 sample was analyzed for EPA 8270. All samples were received in acceptable condition with any exceptions noted below.

Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

Sample Preparation:

The samples were prepared in accordance with EPA 3510 with any exceptions noted below.

Initial Calibrations (including MS Tune as applicable):

All criteria were within method requirements with any exceptions noted below.

Continuing Calibration:

All criteria were within method requirements with any exceptions noted below.

Internal Standards:

All internal standards were within QC limits with any exceptions noted below.

Surrogates:

All surrogates were within QC limits with any exceptions noted below.

Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

QC Batch: OEXT/43560

L0: Analyte recovery in the laboratory control sample (LCS) was outside QC limits.

- LCS (Lab ID: 1357257)
- Benzoic acid

Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

QC Batch: MSSV/13901

A matrix spike/matrix spike duplicate was not performed due to insufficient sample volume.

Additional Comments:



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Method: EPA 5030B/8260

Description:8260 MSVClient:COP Conestoga-Rovers & Associates, Inc. NMDate:April 18, 2014

General Information:

1 sample was analyzed for EPA 5030B/8260. All samples were received in acceptable condition with any exceptions noted below.

Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

Initial Calibrations (including MS Tune as applicable): All criteria were within method requirements with any exceptions noted below.

Continuing Calibration:

All criteria were within method requirements with any exceptions noted below.

Internal Standards:

All internal standards were within QC limits with any exceptions noted below.

Surrogates:

All surrogates were within QC limits with any exceptions noted below.

Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

QC Batch: MSV/60740

A matrix spike/matrix spike duplicate was not performed due to insufficient sample volume.

Additional Comments:



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Method: EPA 120.1

Description:120.1 Specific ConductanceClient:COP Conestoga-Rovers & Associates, Inc. NMDate:April 18, 2014

General Information:

1 sample was analyzed for EPA 120.1. All samples were received in acceptable condition with any exceptions noted below.

Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

Duplicate Sample:

All duplicate sample results were within method acceptance criteria with any exceptions noted below.

Additional Comments:



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Method: SM 2320B

Description:2320B AlkalinityClient:COP Conestoga-Rovers & Associates, Inc. NMDate:April 18, 2014

General Information:

1 sample was analyzed for SM 2320B. All samples were received in acceptable condition with any exceptions noted below.

Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

Duplicate Sample:

All duplicate sample results were within method acceptance criteria with any exceptions noted below.

Additional Comments:



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Method: SM 2540C

Description:2540C Total Dissolved SolidsClient:COP Conestoga-Rovers & Associates, Inc. NMDate:April 18, 2014

General Information:

1 sample was analyzed for SM 2540C. All samples were received in acceptable condition with any exceptions noted below.

Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

Duplicate Sample:

All duplicate sample results were within method acceptance criteria with any exceptions noted below.

Additional Comments:



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Method: SM 4500-H+B

Description:4500H+ pH, ElectrometricClient:COP Conestoga-Rovers & Associates, Inc. NMDate:April 18, 2014

General Information:

1 sample was analyzed for SM 4500-H+B. All samples were received in acceptable condition with any exceptions noted below.

Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

- H6: Analysis initiated outside of the 15 minute EPA recommended holding time.
 - GW-085692-040214-CM-MW-2 (Lab ID: 60166332001)

Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

Duplicate Sample:

All duplicate sample results were within method acceptance criteria with any exceptions noted below.

Additional Comments:



PROJECT NARRATIVE

Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Method: EPA 300.0

Description:300.0 IC Anions 28 DaysClient:COP Conestoga-Rovers & Associates, Inc. NMDate:April 18, 2014

General Information:

1 sample was analyzed for EPA 300.0. All samples were received in acceptable condition with any exceptions noted below.

Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

Initial Calibrations (including MS Tune as applicable):

All criteria were within method requirements with any exceptions noted below.

Continuing Calibration:

All criteria were within method requirements with any exceptions noted below.

Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

QC Batch: WETA/29042

A matrix spike and/or matrix spike duplicate (MS/MSD) were performed on the following sample(s): 60166574001,60166574005

M1: Matrix spike recovery exceeded QC limits. Batch accepted based on laboratory control sample (LCS) recovery.

- MSD (Lab ID: 1361974)
 - Sulfate

Additional Comments:



PROJECT NARRATIVE

Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Method: EPA 353.2

Description:353.2 Nitrogen, NO2/NO3 pres.Client:COP Conestoga-Rovers & Associates, Inc. NMDate:April 18, 2014

General Information:

1 sample was analyzed for EPA 353.2. All samples were received in acceptable condition with any exceptions noted below.

Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

Initial Calibrations (including MS Tune as applicable):

All criteria were within method requirements with any exceptions noted below.

Continuing Calibration:

All criteria were within method requirements with any exceptions noted below.

Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

Duplicate Sample:

All duplicate sample results were within method acceptance criteria with any exceptions noted below.

Additional Comments:



PROJECT NARRATIVE

Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Method: EPA 365.1

Description:365.1 Orthophosphate as PClient:COP Conestoga-Rovers & Associates, Inc. NMDate:April 18, 2014

General Information:

1 sample was analyzed for EPA 365.1. All samples were received in acceptable condition with any exceptions noted below.

Hold Time:

The samples were analyzed within the method required hold times with any exceptions noted below.

Initial Calibrations (including MS Tune as applicable):

All criteria were within method requirements with any exceptions noted below.

Continuing Calibration:

All criteria were within method requirements with any exceptions noted below.

Method Blank:

All analytes were below the report limit in the method blank, where applicable, with any exceptions noted below.

Laboratory Control Spike:

All laboratory control spike compounds were within QC limits with any exceptions noted below.

Matrix Spikes:

All percent recoveries and relative percent differences (RPDs) were within acceptance criteria with any exceptions noted below.

Duplicate Sample:

All duplicate sample results were within method acceptance criteria with any exceptions noted below.

Additional Comments:

This data package has been reviewed for quality and completeness and is approved for release.



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Sample: GW-085692-040214-CM- MW-2	Lab ID: 601	66332001	Collected: 04/02/1	4 13:45	5 Received: 04	1/04/14 08:35 N	latrix: Water	
Parameters	Results	Units	Report Limit	DF	Prepared	Analyzed	CAS No.	Qua
8015B Diesel Range Organics	Analytical Met	hod: EPA 80	15B Preparation Me	ethod: E	PA 3510C			
TPH-DRO	ND m	g/L	0.50	1	04/04/14 00:00	04/07/14 16:29		
Surrogates								
p-Terphenyl (S)	67 %		28-127	1		04/07/14 16:29		
n-Tetracosane (S)	49 %		22-121	1	04/04/14 00:00	04/07/14 16:29	646-31-1	
Gasoline Range Organics	Analytical Met	hod: EPA 50	30B/8015B					
TPH-GRO	ND m	g/L	0.50	1		04/09/14 21:48		
Surrogates								
4-Bromofluorobenzene (S)	94 %		56-137	1		04/09/14 21:48	460-00-4	
Preservation pH	1.0		0.10	1		04/09/14 21:48		
6010 MET ICP, Dissolved	Analytical Met	hod: EPA 60	10 Preparation Met	nod: EP	A 3010			
Total Hardness by 2340B, Dissolved	1180000 ug	g/L	500	1	04/15/14 10:30	04/15/14 15:22		
6020 MET ICPMS, Dissolved	Analytical Met	hod: EPA 60	20 Preparation Met	nod: EP	A 3010			
Arsenic, Dissolved	1.1 ug	g/L	1.0	1	04/14/14 15:50	04/16/14 11:05	7440-38-2	
Barium, Dissolved	12.8 ug		1.0	1	04/14/14 15:50	04/16/14 11:05	7440-39-3	
Cadmium, Dissolved	ND ug	,	0.50	1		04/16/14 11:05		
Chromium, Dissolved	ND ug		1.0	1		04/16/14 11:05		
Cobalt, Dissolved	ND ug		1.0	1		04/16/14 11:05		
Copper, Dissolved	ND ug		1.0	1		04/16/14 11:05		
Lead, Dissolved	ND ug		1.0	1		04/16/14 11:05		
Manganese, Dissolved	853 ug		1.0	1		04/16/14 11:05		M1
Maliganese, Dissolved Molybdenum, Dissolved	3.9 ug	,	1.0	1		04/16/14 11:05		M1
	3.9 ug ND ug		1.0			04/16/14 11:05		
Nickel, Dissolved				1				
Selenium, Dissolved	ND ug		1.0	1		04/16/14 11:05		
Silver, Dissolved	ND uç		0.50	1		04/16/14 11:05		
Zinc, Dissolved	ND uç	g/L	10.0	1	04/14/14 15:50	04/16/14 11:05	7440-66-6	
7470 Mercury, Dissolved	Analytical Met	hod: EPA 74	70 Preparation Met	nod: EP	A 7470			
Mercury, Dissolved	ND uç	g/L	0.20	1	04/07/14 10:00	04/07/14 15:14	7439-97-6	
8270 MSSV Semivolatile Organic	Analytical Met	hod: EPA 82	70 Preparation Met	nod: EP	A 3510			
Acenaphthene	ND uç	g/L	10.0	1	04/09/14 00:00	04/10/14 22:19	83-32-9	
Acenaphthylene	ND ug	g/L	10.0	1	04/09/14 00:00	04/10/14 22:19	208-96-8	
Anthracene	ND uç	g/L	10.0	1	04/09/14 00:00	04/10/14 22:19	120-12-7	
Benzo(a)anthracene	ND ug	g/L	10.0	1	04/09/14 00:00	04/10/14 22:19	56-55-3	
Benzo(a)pyrene	ND ug		10.0	1	04/09/14 00:00	04/10/14 22:19	50-32-8	
Benzo(b)fluoranthene	ND ug		10.0	1	04/09/14 00:00	04/10/14 22:19	205-99-2	
Benzo(g,h,i)perylene	ND ug		10.0	1		04/10/14 22:19		
Benzo(k)fluoranthene	ND ug		10.0	1		04/10/14 22:19		
Benzoic acid	ND ug		50.0	1		04/10/14 22:19		L2
Benzyl alcohol	ND ug		20.0	1		04/10/14 22:19		
4-Bromophenylphenyl ether	ND ug		10.0	1		04/10/14 22:19		
Butylbenzylphthalate	ND ug		10.0	1	04/09/14 00:00			

REPORT OF LABORATORY ANALYSIS

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Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Sample: GW-085692-040214-CM- MW-2	Lab ID: 60166	332001 Collected:	04/02/1	14 13:45	Received: 04	/04/14 08:35	Matrix: Water	
Parameters	Results	Units Repo	rt Limit	DF	Prepared	Analyzed	CAS No.	Qual
8270 MSSV Semivolatile Organic	Analytical Metho	d: EPA 8270 Prepara	tion Met	hod: EPA	3510			
Carbazole	ND ug/L		10.0	1	04/09/14 00:00	04/10/14 22:19	86-74-8	
4-Chloro-3-methylphenol	ND ug/L		20.0	1	04/09/14 00:00	04/10/14 22:19	59-50-7	
4-Chloroaniline	ND ug/L		20.0	1	04/09/14 00:00	04/10/14 22:19	106-47-8	
bis(2-Chloroethoxy)methane	ND ug/L		10.0	1	04/09/14 00:00	04/10/14 22:19	111-91-1	
bis(2-Chloroethyl) ether	ND ug/L		10.0	1	04/09/14 00:00	04/10/14 22:19	111-44-4	
bis(2-Chloroisopropyl) ether	ND ug/L		10.0	1	04/09/14 00:00	04/10/14 22:19	39638-32-9	
2-Chloronaphthalene	ND ug/L		10.0	1	04/09/14 00:00	04/10/14 22:19	91-58-7	
2-Chlorophenol	ND ug/L		10.0	1		04/10/14 22:19		
4-Chlorophenylphenyl ether	ND ug/L		10.0	1		04/10/14 22:19		
Chrysene	ND ug/L		10.0	1		04/10/14 22:19		
Dibenz(a,h)anthracene	ND ug/L		10.0	1		04/10/14 22:19		
Dibenzofuran	ND ug/L		10.0	1		04/10/14 22:19		
1,2-Dichlorobenzene	ND ug/L		10.0	1		04/10/14 22:19		
1,3-Dichlorobenzene	ND ug/L		10.0	1		04/10/14 22:19		
1,4-Dichlorobenzene	ND ug/L		10.0	1		04/10/14 22:19		
3,3'-Dichlorobenzidine	ND ug/L		20.0	1		04/10/14 22:19		
2,4-Dichlorophenol	ND ug/L		10.0	1		04/10/14 22:19		
Diethylphthalate	ND ug/L		10.0	1		04/10/14 22:19		
2,4-Dimethylphenol	ND ug/L		10.0	1		04/10/14 22:19		
Dimethylphthalate	ND ug/L		10.0	1		04/10/14 22:19		
	-		10.0	1		04/10/14 22:19		
Di-n-butylphthalate	ND ug/L		50.0	1		04/10/14 22:19		
4,6-Dinitro-2-methylphenol	ND ug/L							
2,4-Dinitrophenol	ND ug/L		50.0 10.0	1 1		04/10/14 22:19 04/10/14 22:19		
2,4-Dinitrotoluene	ND ug/L							
2,6-Dinitrotoluene	ND ug/L		10.0	1		04/10/14 22:19		
Di-n-octylphthalate	ND ug/L		10.0	1		04/10/14 22:19		
bis(2-Ethylhexyl)phthalate	ND ug/L		10.0	1		04/10/14 22:19		
Fluoranthene	ND ug/L		10.0	1		04/10/14 22:19		
Fluorene	ND ug/L		10.0	1		04/10/14 22:19		
Hexachloro-1,3-butadiene	ND ug/L		10.0	1		04/10/14 22:19		
Hexachlorobenzene	ND ug/L		10.0	1		04/10/14 22:19		
Hexachlorocyclopentadiene	ND ug/L		10.0	1		04/10/14 22:19		
Hexachloroethane	ND ug/L		10.0	1		04/10/14 22:19	-	
Indeno(1,2,3-cd)pyrene	ND ug/L		10.0	1		04/10/14 22:19		
Isophorone	ND ug/L		10.0	1		04/10/14 22:19		
2-Methylnaphthalene	ND ug/L		10.0	1		04/10/14 22:19		
2-Methylphenol(o-Cresol)	ND ug/L		10.0	1		04/10/14 22:19		
3&4-Methylphenol(m&p Cresol)	ND ug/L		10.0	1		04/10/14 22:19		
Naphthalene	ND ug/L		10.0	1		04/10/14 22:19		
2-Nitroaniline	ND ug/L		50.0	1		04/10/14 22:19		
3-Nitroaniline	ND ug/L		50.0	1		04/10/14 22:19		
4-Nitroaniline	ND ug/L		50.0	1		04/10/14 22:19		
Nitrobenzene	ND ug/L		10.0	1		04/10/14 22:19		
2-Nitrophenol	ND ug/L		10.0	1	04/09/14 00:00	04/10/14 22:19	88-75-5	
4-Nitrophenol	ND ug/L		50.0	1	04/09/14 00:00	04/10/14 22:19	100-02-7	
N-Nitroso-di-n-propylamine	ND ug/L		10.0	1	04/09/14 00:00	04/10/14 22:19	621-64-7	

REPORT OF LABORATORY ANALYSIS

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Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Sample: GW-085692-040214-CM- MW-2	Lab ID: 6016	6332001	Collected:	04/02/1	4 13:45	Received: 0	4/04/14 08:35	Matrix: Water	
Parameters	Results	Units	Repor	t Limit	DF	Prepared	Analyzed	CAS No.	Qual
8270 MSSV Semivolatile Organic	Analytical Meth	od: EPA 82	270 Preparat	tion Meth	nod: EPA	3510			
N-Nitrosodiphenylamine	ND ug/	L		10.0	1	04/09/14 00:00	0 04/10/14 22:1	9 86-30-6	
Pentachlorophenol	ND ug/	L		50.0	1	04/09/14 00:00	0 04/10/14 22:1	9 87-86-5	
Phenanthrene	ND ug/	L		10.0	1	04/09/14 00:00	0 04/10/14 22:1	9 85-01-8	
Phenol	ND ug/	L		10.0	1	04/09/14 00:00	0 04/10/14 22:1	9 108-95-2	
Pyrene	ND ug/	L		10.0	1	04/09/14 00:00	0 04/10/14 22:1	9 129-00-0	
Pyridine	ND ug/	L		10.0	1	04/09/14 00:00	0 04/10/14 22:1	9 110-86-1	
1,2,4-Trichlorobenzene	ND ug/			10.0	1	04/09/14 00:00	0 04/10/14 22:1	9 120-82-1	
2,4,5-Trichlorophenol	ND ug/			50.0	1	04/09/14 00:00	0 04/10/14 22:1	9 95-95-4	
2,4,6-Trichlorophenol	ND ug/			10.0	1	04/09/14 00:00	0 04/10/14 22:1	9 88-06-2	
Surrogates	0								
Nitrobenzene-d5 (S)	85 %			10-135	1	04/09/14 00:00	0 04/10/14 22:1	9 4165-60-0	
2-Fluorobiphenyl (S)	89 %			19-124	1	04/09/14 00:00	0 04/10/14 22:1	9 321-60-8	
Terphenyl-d14 (S)	87 %			24-131	1	04/09/14 00:00	0 04/10/14 22:1	9 1718-51-0	
Phenol-d6 (S)	26 %			10-120	1	04/09/14 00:00	0 04/10/14 22:1	9 13127-88-3	
2-Fluorophenol (S)	40 %			13-120	1	04/09/14 00:00	0 04/10/14 22:1	9 367-12-4	
2,4,6-Tribromophenol (S)	84 %			29-121	1	04/09/14 00:00	0 04/10/14 22:1	9 118-79-6	
8260 MSV	Analytical Meth	od: EPA 50)30B/8260						
Acetone	ND ug/	L		10.0	1		04/12/14 06:1	6 67-64-1	
Benzene	ND ug/			1.0	1		04/12/14 06:1	6 71-43-2	
Bromobenzene	ND ug/	L		1.0	1		04/12/14 06:1	6 108-86-1	
Bromochloromethane	ND ug/			1.0	1		04/12/14 06:1	6 74-97-5	
Bromodichloromethane	ND ug/			1.0	1		04/12/14 06:1		
Bromoform	ND ug/			1.0	1		04/12/14 06:1	6 75-25-2	
Bromomethane	ND ug/			5.0	1		04/12/14 06:1		
2-Butanone (MEK)	ND ug/			10.0	1		04/12/14 06:1		
n-Butylbenzene	ND ug/			1.0	1		04/12/14 06:1		
sec-Butylbenzene	ND ug/			1.0	1		04/12/14 06:1		
tert-Butylbenzene	ND ug/			1.0	1		04/12/14 06:1		
Carbon disulfide	ND ug/			5.0	1		04/12/14 06:1		
Carbon tetrachloride	ND ug/			1.0	1		04/12/14 06:1		
Chlorobenzene	ND ug/			1.0	1		04/12/14 06:1		
Chloroethane	ND ug/			1.0	1		04/12/14 06:1		
Chloroform	ND ug/			1.0	1		04/12/14 06:1		
Chloromethane	ND ug/			1.0	1		04/12/14 06:1		
2-Chlorotoluene	ND ug/			1.0	1		04/12/14 06:1		
4-Chlorotoluene	ND ug/			1.0	1		04/12/14 06:1		
1,2-Dibromo-3-chloropropane	ND ug/			2.5	1		04/12/14 06:1		
Dibromochloromethane	ND ug/			2.5 1.0	1		04/12/14 06:1		
1,2-Dibromoethane (EDB)	-			1.0	1		04/12/14 06:1		
Dibromomethane	ND ug/			1.0	1		04/12/14 06:1		
	ND ug/								
	-								
-	0								
	-								
Dichlorodifluoromethane 1,1-Dichloroethane	0			1.0 1.0			04/12/14 06:1 04/12/14 06:1		
1,2-Dichlorobenzene 1,3-Dichlorobenzene 1,4-Dichlorobenzene Dichlorodifluoromethane 1,1-Dichloroethane	ND ug/ ND ug/ ND ug/ ND ug/ ND ug/	L L L		1.0 1.0 1.0 1.0 1.0	1 1 1 1 1		04/12/14 06:1 04/12/14 06:1 04/12/14 06:1 04/12/14 06:1 04/12/14 06:1	6 541-73-1 6 106-46-7 6 75-71-8	



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Sample: GW-085692-040214-CM- MW-2	Lab ID: 60166332001	Collected: 04/02/1	4 13:45	Received: 04/04/14 08:3	5 Matrix: Water	
Parameters	Results Units	Report Limit	DF	Prepared Analyz	ed CAS No.	Qual
8260 MSV	Analytical Method: EPA	5030B/8260				
1,2-Dichloroethane	ND ug/L	1.0	1	04/12/14 (6:16 107-06-2	
1,2-Dichloroethene (Total)	ND ug/L	1.0	1	04/12/14 (6:16 540-59-0	
1,1-Dichloroethene	ND ug/L	1.0	1	04/12/14 (6:16 75-35-4	
cis-1,2-Dichloroethene	ND ug/L	1.0	1	04/12/14 (6:16 156-59-2	
trans-1,2-Dichloroethene	ND ug/L	1.0	1	04/12/14 (6:16 156-60-5	
1,2-Dichloropropane	ND ug/L	1.0	1	04/12/14 (6:16 78-87-5	
1,3-Dichloropropane	ND ug/L	1.0	1	04/12/14 (6:16 142-28-9	
2,2-Dichloropropane	ND ug/L	1.0	1	04/12/14 (6:16 594-20-7	
1,1-Dichloropropene	ND ug/L	1.0	1	04/12/14 (6:16 563-58-6	
cis-1,3-Dichloropropene	ND ug/L	1.0	1	04/12/14 (6:16 10061-01-5	
trans-1,3-Dichloropropene	ND ug/L	1.0	1		6:16 10061-02-6	
Ethylbenzene	ND ug/L	1.0	1		6:16 100-41-4	
Hexachloro-1,3-butadiene	ND ug/L	1.0	1		6:16 87-68-3	
2-Hexanone	ND ug/L	10.0	1		6:16 591-78-6	
Isopropylbenzene (Cumene)	ND ug/L	1.0	1		6:16 98-82-8	
p-Isopropyltoluene	ND ug/L	1.0	1		6:16 99-87-6	
Methylene chloride	ND ug/L	1.0	1		6:16 75-09-2	
4-Methyl-2-pentanone (MIBK)	ND ug/L	10.0	1		6:16 108-10-1	
Methyl-tert-butyl ether	ND ug/L	1.0	1		6:16 1634-04-4	
Naphthalene	ND ug/L	10.0	1		6:16 91-20-3	
n-Propylbenzene	ND ug/L	1.0	1		6:16 103-65-1	
Styrene	ND ug/L	1.0	1		6:16 100-42-5	
1,1,1,2-Tetrachloroethane	ND ug/L	1.0	1		6:16 630-20-6	
1,1,2,2-Tetrachloroethane	ND ug/L	1.0	1		6:16 79-34-5	
Tetrachloroethene	ND ug/L	1.0	1		0:10 79-34-3 06:16 127-18-4	
Toluene	ND ug/L	1.0	1		6:16 108-88-3	
	-	1.0	1		6:16 87-61-6	
1,2,3-Trichlorobenzene	ND ug/L ND ug/L	1.0	1		6:16 120-82-1	
1,2,4-Trichlorobenzene	-	1.0	1			
1,1,1-Trichloroethane 1,1,2-Trichloroethane	ND ug/L	1.0	1		06:16 71-55-6	
	ND ug/L		1		06:16 79-00-5	
Trichloroethene Trichlorofluoromethane	ND ug/L	1.0			06:16 79-01-6	
	ND ug/L	1.0	1		06:16 75-69-4	
1,2,3-Trichloropropane	ND ug/L	2.5	1		06:16 96-18-4	
1,2,4-Trimethylbenzene	ND ug/L	1.0	1		16:16 95-63-6	
1,3,5-Trimethylbenzene	ND ug/L	1.0	1		6:16 108-67-8	
Vinyl chloride	ND ug/L	1.0	1		06:16 75-01-4	
Xylene (Total)	ND ug/L	3.0	1	04/12/14 (6:16 1330-20-7	
<i>Surrogates</i> 4-Bromofluorobenzene (S)	99 %	80-120	1	01/10/11	6:16 460-00-4	
1,2-Dichloroethane-d4 (S)	99 % 104 %	80-120	1		6:16 460-00-4 6:16 17060-07-0	
Toluene-d8 (S)	104 %	80-120 80-120			06:16 2037-26-5	
Preservation pH	109 % 1.0	80-120 0.10	1 1	04/12/14 (
·			I	04/12/14 (0.10	
120.1 Specific Conductance	Analytical Method: EPA					
Specific Conductance	3470 umhos/cm	1.0	1	04/08/14 1	0:11	

REPORT OF LABORATORY ANALYSIS

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Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Sample: GW-085692-040214-CM- MW-2	Lab ID: 6016633200	1 Collected: 04/02/	14 13:45	Received: 04/04/14 08:35	Matrix: Water	
Parameters	Results Units	Report Limit	DF	Prepared Analyzed	CAS No.	Qual
2320B Alkalinity	Analytical Method: SM 2	2320B				
Alkalinity, Total as CaCO3	290 mg/L	20.0	1	04/14/14 12:	58	
2540C Total Dissolved Solids	Analytical Method: SM 2	2540C				
Total Dissolved Solids	3030 mg/L	5.0	1	04/08/14 12:	34	
4500H+ pH, Electrometric	Analytical Method: SM	4500-H+B				
pH at 25 Degrees C	7.3 Std. Units	0.10	1	04/08/14 12:	15	H6
300.0 IC Anions 28 Days	Analytical Method: EPA	300.0				
Bromide	ND mg/L	1.0	1	04/17/14 20:	12 24959-67-9	
Chloride	41.3 mg/L	5.0	5	04/17/14 20:	27 16887-00-6	
Fluoride	0.68 mg/L	0.20	1	04/17/14 20:	12 16984-48-8	
Sulfate	2360 mg/L	200	200	04/17/14 20	43 14808-79-8	
353.2 Nitrogen, NO2/NO3 pres.	Analytical Method: EPA	353.2				
Nitrogen, NO2 plus NO3	ND mg/L	0.10	1	04/08/14 16	24	
365.1 Orthophosphate as P	Analytical Method: EPA	365.1				
Orthophosphate as P	0.10 mg/L	0.10	1	04/04/14 12:	46	



Project: 085692 MARC	OTTE NO 1						
Pace Project No.: 60166332							
QC Batch: GCV/4730		Analysis	Method:	EPA 5030B/80	15B		
QC Batch Method: EPA 5030B/8	015B	Analysis	Description:	Gasoline Rang	e Organics		
Associated Lab Samples: 60166	332001						
METHOD BLANK: 1357152		Mat	trix: Water				
Associated Lab Samples: 60166	332001						
		Blank	Reportin	0			
Parameter	Units	Result	Limit	Analyze	d Quali	fiers	
TPH-GRO	mg/L	1	ND (0.50 04/09/14 1	4:51		
4-Bromofluorobenzene (S)	%	1	00 56-	137 04/09/14 1	4:51		
LABORATORY CONTROL SAMPL	E: 1357153						
		Spike	LCS	LCS	% Rec		
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers	
TPH-GRO	mg/L		0.88	88	62-132		
4-Bromofluorobenzene (S)							



Project:	085692 MARCO	TTE NO '	1										
Pace Project No.:	60166332												
QC Batch:	MERP/8283			Analys	is Method:	E	PA 7470						
QC Batch Method:	EPA 7470			Analys	is Descript	ion: 74	470 Mercury	,Dissolved					
Associated Lab Sar	mples: 6016633	2001											
METHOD BLANK:	1356343			Ν	Aatrix: Wa	er							
Associated Lab Sar	mples: 6016633	2001											
_				Blank		eporting			o				
Parar	neter		Jnits	Resul	t	Limit	Analyz	ed	Qualifiers				
Mercury, Dissolved		ug/L			ND	0.20	04/07/14	14:59					
LABORATORY CO	NTROL SAMPLE:	135634	14										
					1.00		LCS	% Rec					
				Spike	LCS		LUS	70 NEC					
Parar	neter	ι	Jnits	Spike Conc.	Resu		% Rec	Limits		ualifiers			
Parar Mercury, Dissolved		ug/L	Jnits		Resu			Limits		ualifiers	-		
		ug/L		Conc. 5	Resu	lt	% Rec	Limits	Q	ualifiers	-		
Mercury, Dissolved		ug/L PLICATE	: 13563	Conc. 5	Resu	lt 4.1 1356346	% Rec 82	Limits 80	-120 Q		-		
Mercury, Dissolved	ATRIX SPIKE DU	ug/L PLICATE 6016	: 13563 66332001	Conc. 5 45 MS Spike	MSD Spike	lt 4.1 1356346 MS	% Rec 82 MSD	Limits 80-	-120 Q MSD	% Rec	-	Max	
Mercury, Dissolved	ATRIX SPIKE DU	ug/L PLICATE	: 13563	Conc. 5 45 MS	MSD	lt 4.1 1356346	% Rec 82	Limits 80	-120 Q		RPD	Max RPD	Qual



Project:	085692 MARCOT	TE NO 1						
Pace Project No.:	60166332							
QC Batch:	MPRP/26840		Analysis	Method:	EPA 60	10		
QC Batch Method:	EPA 3010		Analysis	Description:	6010 M	ET Dissolved		
Associated Lab Sam	ples: 60166332	2001						
METHOD BLANK:	1360801		Ma	atrix: Water				
Associated Lab Sam	ples: 60166332	2001						
			Blank	Repor	0			
Param	eter	Units	Result	Lim	it /	Analyzed	Qualifiers	
Total Hardness by 23 Dissolved	340B,	ug/L		ND	500 04/	15/14 15:17		
LABORATORY CON	TROL SAMPLE:	1360802						
			Spike	LCS	LCS	% Re	C	
Param	eter	Units	Conc.	Result	% Re	c Limit	s Qualifier	S
Total Hardness by 23 Dissolved	340B,	ug/L		6380	0			



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

QC Batch: MPRP/26	827	Analysis Meth	nod: EF	PA 6020	
QC Batch Method: EPA 3010		Analysis Desc	cription: 60	20 MET Dissolved	
Associated Lab Samples: 601	66332001				
METHOD BLANK: 1360508		Matrix:	Water		
Associated Lab Samples: 601	66332001				
		Blank	Reporting		
Parameter	Units	Result	Limit	Analyzed	Qualifiers
Arsenic, Dissolved	ug/L	ND	1.0	04/16/14 10:57	
Barium, Dissolved	ug/L	ND	1.0	04/16/14 10:57	
Cadmium, Dissolved	ug/L	ND	0.50	04/16/14 10:57	
Chromium, Dissolved	ug/L	ND	1.0	04/16/14 10:57	
Cobalt, Dissolved	ug/L	ND	1.0	04/16/14 10:57	
Copper, Dissolved	ug/L	ND	1.0	04/16/14 10:57	
Lead, Dissolved	ug/L	ND	1.0	04/16/14 10:57	
Manganese, Dissolved	ug/L	1.2	1.0	04/16/14 10:57	
Molybdenum, Dissolved	ug/L	ND	1.0	04/16/14 10:57	

ND

ND

ND

ND

1.0 04/16/14 10:57

1.0 04/16/14 10:57

0.50 04/16/14 10:57

10.0 04/16/14 10:57

LABORATORY CONTROL SAMPLE: 1360509

ug/L

ug/L

ug/L

ug/L

Nickel, Dissolved

Silver, Dissolved

Zinc, Dissolved

Selenium, Dissolved

		Spike	LCS	LCS	% Rec	
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers
rsenic, Dissolved	ug/L	40	40.0	100	80-120	
arium, Dissolved	ug/L	40	41.9	105	80-120	
admium, Dissolved	ug/L	40	40.0	100	80-120	
nromium, Dissolved	ug/L	40	40.5	101	80-120	
obalt, Dissolved	ug/L	40	40.4	101	80-120	
opper, Dissolved	ug/L	40	40.8	102	80-120	
ad, Dissolved	ug/L	40	40.5	101	80-120	
anganese, Dissolved	ug/L	40	40.9	102	80-120	
lybdenum, Dissolved	ug/L	40	41.7	104	80-120	
ckel, Dissolved	ug/L	40	40.2	100	80-120	
elenium, Dissolved	ug/L	40	39.5	99	80-120	
lver, Dissolved	ug/L	20	19.8	99	80-120	
nc, Dissolved	ug/L	100	103	103	80-120	

MATRIX SPIKE SAMPLE:	1360510						
		60166332001	Spike	MS	MS	% Rec	
Parameter	Units	Result	Conc.	Result	% Rec	Limits	Qualifiers
Arsenic, Dissolved	ug/L	1.1	40	43.4	106	75-125	
Barium, Dissolved	ug/L	12.8	40	56.8	110	75-125	
Cadmium, Dissolved	ug/L	ND	40	36.2	90	75-125	
Chromium, Dissolved	ug/L	ND	40	40.1	99	75-125	
Cobalt, Dissolved	ug/L	ND	40	38.3	95	75-125	
Copper, Dissolved	ug/L	ND	40	34.6	85	75-125	

REPORT OF LABORATORY ANALYSIS

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Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

MATRIX SPIKE SAMPLE:	1360510						
Parameter	Units	60166332001 Result	Spike Conc.	MS Result	MS % Rec	% Rec Limits	Qualifiers
Lead, Dissolved	ug/L	ND	40	45.2	113	75-125	
Manganese, Dissolved	ug/L	853	40	951	246	75-125	M1
Molybdenum, Dissolved	ug/L	3.9	40	54.7	127	75-125	M1
Nickel, Dissolved	ug/L	ND	40	35.6	89	75-125	
Selenium, Dissolved	ug/L	ND	40	44.9	112	75-125	
Silver, Dissolved	ug/L	ND	20	17.8	89	75-125	
Zinc, Dissolved	ug/L	ND	100	83.8	82	75-125	



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

QC Batch Method: EPA 50308/8260 Analysis Description: 8260 MSV Water 10 mL Purge Associated Lab Samples: 60166332001 Matrix: Water Matrix: Water Associated Lab Samples: 60166332001 Blank Reporting Analyzed Qualifiers 1,1,1,2-Tetrachloroethane ug/L ND 1.0 04/12/14 03.25 1.1 1,1,2-Tetrachloroethane ug/L ND 1.0 04/12/14 03.25 1.1 1,1-Dichloroethane ug/L ND 1.0 04/12/14 03.25 1.1 1,1-Dichloroethane ug/L ND 1.0 04/12/14 03.25 1.2 1,2-Trithichloroptinane ug/L ND 1.0 04/12/1	QC Batch: MSV/60740	40 Analysis Me		thod: EPA 5030B/8260			
Associated Lab Samples: 60166332001 METHOD BLANK: 1356975 Matrix: Water Associated Lab Samples: 60166332001 Ellank Reporting Limit Analyzed Qualifiers 1,1,1-Trethrachloroethane ugL ND 1.0 04/12/14 03:25 1,1,1-Trichioroethane ugL ND 1.0 04/12/14 03:25 1,1,2-Titchioroethane ugL ND 1.0 04/12/14 03:25 1,1,2-Titchioroethane ugL ND 1.0 04/12/14 03:25 1,1-Dichioroethane ugL ND 1.0 04/12/14 03:25 1,2-Titchioroethane ugL ND 1.0 04/12/14 03:25 1,2-Titchioropropane ugL ND 1.0 04/12/14 03:25 1,2-D	QC Batch Method: EPA 5030B/8	3260	•		8260 MSV Water 10 mL Purge		
METHOD BLANK: 1358975 Matrix: Water Associated Lab Samples: 60166332001 Elink Reporting Limit Analyzed Qualifiers 1.1,1.2-Tetrachiorcethane ug/L ND 1.0 04/12/14 03.25 1.1.1:Trichiorcethane ug/L ND 1.0 04/12/14 03.25 1.1.2:Trichiorcethane ug/L ND 1.0 04/12/14 03.25 1.1.2:Trichiorcethane ug/L ND 1.0 04/12/14 03.25 1.1.2:Trichiorcethane ug/L ND 1.0 04/12/14 03.25 1.1.Dichiorcethene ug/L ND 1.0 04/12/14 03.25 1.2.3.Trichiorcopane ug/L ND 1.0 04/12/14 03.25 1.2.3.Trichiorcopane ug/L ND 1.0 04/12/14 03.25 1.2.4.Trimehytibenzene ug/L ND 1.0 04/12/14 03.25 1.2.4.Trimehytibenzene ug/L ND 1.0 04/12/14 03.25 1.2.2.Trichiorcopane ug/L ND 1.0 04/12/14 03.25 1.2.2.Trichiorcopane ug/L ND 1.0 04/12/14 03.25 1.2.2.Dichiorcentane Ug/L ND <				onp			
Associated Lab Samples: 60166332001 Blank Result Reporting Limit Analyzed Analyzed Qualifiers. 1.1.1.2-Tetrachloroethane ug/L ND 1.0 04/12/14 03:25 0.0		332001					
Parameter Units Result Limit Analyzed Qualifiers 1.1.1.2-Tetrachloroethane ugL ND 1.0 04/12/14 03:25 1.1.1.2-Tetrachloroethane ugL ND 1.0 04/12/14 03:25 1.1.2.2-Tetrachloroethane ugL ND 1.0 04/12/14 03:25 1.1.2-Tetrachloroethane ugL ND 1.0 04/12/14 03:25 1.1.Dichloroethane ugL ND 1.0 04/12/14 03:25 1.2.3-Trichloroptopane ugL ND 1.0 04/12/14 03:25 1.2.3-Trichloroptopane ugL ND 1.0 04/12/14 03:25 1.2.4-Trimethylbenzene ugL ND 1.0 04/12/14 03:25 1.2.Dibromoethane (EDB) ugL ND 1.0 04/12/14 03:25 1.2-Dibromoethane ug/L ND 1.0 04/12/14 03:25 1.2-Dibromoethane ug/L ND 1.0 04/12/14 03:25	METHOD BLANK: 1358975		Matrix:	Water			
Parameter Units Result Limit Analyzed Qualifiers 1,1,1,2-Tetrachloroethane ug/L ND 1.0 04/12/14 03:25 1,1,1-Trichloroethane ug/L ND 1.0 04/12/14 03:25 1,1,2-Zitrachloroethane ug/L ND 1.0 04/12/14 03:25 1,1-Dickhoroethane ug/L ND 1.0 04/12/14 03:25 1,1-Dickhoroethane ug/L ND 1.0 04/12/14 03:25 1,2.3-Trichloroppene ug/L ND 1.0 04/12/14 03:25 1,2.3-Trichloroppone ug/L ND 1.0 04/12/14 03:25 1,2.3-Trichloroppone ug/L ND 1.0 04/12/14 03:25 1,2-Dibromo-3-chloropropane ug/L ND 1.0 04/12/14 03:25 1,2-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroptopane ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroptopane ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroptopane<	Associated Lab Samples: 60166	332001					
Parameter Units Result Limit Analyzed Qualifiers 1,1,1,-Tichtochane ug/L ND 1.0 04/12/14 03:25 1,1,1-Tichtoroethane ug/L ND 1.0 04/12/14 03:25 1,1,2,2-Tetrachloroethane ug/L ND 1.0 04/12/14 03:25 1,1-Dichtoroethane ug/L ND 1.0 04/12/14 03:25 1,1-Dichtoroethane ug/L ND 1.0 04/12/14 03:25 1,2.3-Tichtoropropene ug/L ND 1.0 04/12/14 03:25 1,2.3-Tichtoropropane ug/L ND 1.0 04/12/14 03:25 1,2.3-Tichtoropropane ug/L ND 1.0 04/12/14 03:25 1,2-Dichtorobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dichtoroethane			Blank	Reporting			
1,1.1-Trichloroethane ug/L ND 1.0 04/12/14 03:25 1,1.2.2.7:brichoroethane ug/L ND 1.0 04/12/14 03:25 1,1.2.Trichioroethane ug/L ND 1.0 04/12/14 03:25 1,1.Dichioroethane ug/L ND 1.0 04/12/14 03:25 1,1.Dichioroethane ug/L ND 1.0 04/12/14 03:25 1,2.3.Trichioropropane ug/L ND 1.0 04/12/14 03:25 1,2.3.Trichioropropane ug/L ND 1.0 04/12/14 03:25 1,2.4.Trinchryberzene ug/L ND 1.0 04/12/14 03:25 1,2.4.Trinchryberzene ug/L ND 1.0 04/12/14 03:25 1,2.Dichoros-acheropropane ug/L ND 1.0 04/12/14 03:25 1,2.Dichoros-acheroe ug/L ND 1.0 04/12/14 03:25 1,2.Dichoropropane ug/L ND 1.0 04/12/14 03:25 1,2.Dichoropropane ug/L ND 1.0 04/12/14 03:25 1,3.Dichoropropane ug/L ND 1.0 04/12/14 03:25 1,3.Dichoropropane<	Parameter	Units	Result	Limit	Analyzed	Qualifiers	
1,1,2-Tirchloroethane ug/L ND 1.0 04/12/14 03:25 1,1-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,1-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,1-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,2-3-Tirchlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2.3-Tirchlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2.3-Tirchlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2.4-Tirnethylbenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dirbromo-3-chloropropane ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroethane (EDB) ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,3-Dichlorobenzene<	1,1,1,2-Tetrachloroethane	ug/L	ND	1.0	04/12/14 03:25		
1,1-2:Trichloroethane ug/L ND 1.0 04/12/14 03:25 1,1-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,1-Dichloropropane ug/L ND 1.0 04/12/14 03:25 1,1-Dichloropropane ug/L ND 1.0 04/12/14 03:25 1,2,3-Trichloropropane ug/L ND 1.0 04/12/14 03:25 1,2,3-Trichloropropane ug/L ND 1.0 04/12/14 03:25 1,2-A-Trinchtylbenzene ug/L ND 1.0 04/12/14 03:25 1,2-Ditorno-s-chloropropane ug/L ND 1.0 04/12/14 03:25 1,2-Dichoroethane ug/L ND 1.0 04/12/14 03:25 1,2-Dichoropthane ug/L ND 1.0 04/12/14 03:25 1,3-Dichoroppane ug/L ND 1.0 04/12/14 03:25 1,3-Dichoroppane ug/L	1,1,1-Trichloroethane	ug/L	ND	1.0	04/12/14 03:25		
1,1-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,1-Dichloropropene ug/L ND 1.0 04/12/14 03:25 1,2,3-Trichloroperpene ug/L ND 1.0 04/12/14 03:25 1,2,3-Trichloroperpene ug/L ND 1.0 04/12/14 03:25 1,2,3-Trichloroperpene ug/L ND 1.0 04/12/14 03:25 1,2,4-Trintothybenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dichrobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dichrobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dichrobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dichrobethane (EDB) ug/L ND 1.0 04/12/14 03:25 1,2-Dichrobethane ug/L ND 1.0 04/12/14 03:25 1,2-Dichrobethane ug/L ND 1.0 04/12/14 03:25 1,2-Dichropropane ug/L ND 1.0 04/12/14 03:25 1,3-Dichropropane ug/L ND 1.0 04/12/14 03:25 1,3-Dichropropane ug/L <td>1,1,2,2-Tetrachloroethane</td> <td>ug/L</td> <td>ND</td> <td>1.0</td> <td>04/12/14 03:25</td> <td></td>	1,1,2,2-Tetrachloroethane	ug/L	ND	1.0	04/12/14 03:25		
1,1-Dichloropenene ug/L ND 1.0 04/12/14 03:25 1,1-Dichloropropene ug/L ND 1.0 04/12/14 03:25 1,2.3-Trichloropenzene ug/L ND 1.0 04/12/14 03:25 1,2.4-Trinchlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2.4-Trinchlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 1,3-Dichloropropane ug/L ND 1.0 04/12/14 03:25 1,3-Dichloropropane ug/L ND 1.0 04/12/14 03:25 1,3-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane	1,1,2-Trichloroethane	ug/L	ND	1.0	04/12/14 03:25		
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1,2,3-Trichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2,3-Trichloropenzene ug/L ND 2.5 04/12/14 03:25 1,2,4-Trichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Trichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dibromo-3-chioropropane ug/L ND 1.0 04/12/14 03:25 1,2-Dibromo-s-chioropropane ug/L ND 1.0 04/12/14 03:25 1,2-Dibromo-s-chioropropane ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroethane (EDB) ug/L ND 1.0 04/12/14 03:25 1,2-Dichloropethene (Total) ug/L ND 1.0 04/12/14 03:25 1,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 1,3-Dichloroporpane ug/L ND 1.0 04/12/14 03:25 1,3-Dichloroporpane ug/L ND 1.0 04/12/14 03:25 1,3-Dichloroporpane ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25	1,1-Dichloroethene	ug/L	ND	1.0	04/12/14 03:25		
1,2,3-Trichloropropane ug/L ND 2.5 04/12/14 03:25 1,2,4-Trinchlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2,4-Trinchlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dibromo-3-chloropropane ug/L ND 1.0 04/12/14 03:25 1,2-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroptopane ug/L ND 1.0 04/12/14 03:25 1,3-Dichloroptopane ug/L ND 1.0 04/12/14 03:25 2-Dichloroptopane ug/L ND 1.0 04/12/14 03:25 2-Dichloroptopane ug/L ND 1.0 04/12/14 03:25 2-Dichloroptopane	1,1-Dichloropropene	ug/L	ND	1.0	04/12/14 03:25		
1,2,3-Trichloropropane ug/L ND 2.5 04/12/14 03:25 1,2,4-Trinehlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2,4-Trinehlyblenzene ug/L ND 2.5 04/12/14 03:25 1,2-Dibromo-3-chloropropane ug/L ND 1.0 04/12/14 03:25 1,2-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroptopane ug/L ND 1.0 04/12/14 03:25 1,3-Dichloroptopane ug/L ND 1.0 04/12/14 03:25 1,3-Dichloroptopane ug/L ND 1.0 04/12/14 03:25 1,3-Dichloroptopane ug/L ND 1.0 04/12/14 03:25 2,2-Dichloroptopane ug/L ND 1.0 04/12/14 03:25 2,2-Dichloroptopane ug/L ND 1.0 04/12/14 03:25 2,2-Dichloroptopane ug/L ND 1.0 04/12/14 03:25 2,-Dichloroptopane	1,2,3-Trichlorobenzene	ug/L	ND	1.0	04/12/14 03:25		
1,2,4-Trimethylbenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dibromo-3-chloropropane ug/L ND 1.0 04/12/14 03:25 1,2-Dibromo-3-chloropropane ug/L ND 1.0 04/12/14 03:25 1,2-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 1,3-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,3-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,3-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2,-Ehorotoluene ug/L ND 1.0 04/12/14 03:25 2,-Ehorotoluene ug/L ND 1.0 04/12/14 03:25 2,-Hexanone ug/	1,2,3-Trichloropropane		ND	2.5	04/12/14 03:25		
1,2,4-Trimethylbenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dibromo-3-chloropropane ug/L ND 1.0 04/12/14 03:25 1,2-Dibromoethane (EDB) ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroethane (Total) ug/L ND 1.0 04/12/14 03:25 1,3-Dichloropethane (Total) ug/L ND 1.0 04/12/14 03:25 1,3-Dichloropethane ug/L ND 1.0 04/12/14 03:25 1,3-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2,-Ehorotoluene ug/L ND 1.0 04/12/14 03:25 2,-Ehorotoluene ug/L ND 1.0 04/12/14 03:25 2,-Hexanone	1,2,4-Trichlorobenzene	ug/L	ND	1.0	04/12/14 03:25		
1,2-Dibromo-3-chloropropane ug/L ND 2.5 04/12/14 03:25 1,2-Dibromoethane (EDB) ug/L ND 1.0 04/12/14 03:25 1,2-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroethene (Total) ug/L ND 1.0 04/12/14 03:25 1,3-Dichloropropane ug/L ND 1.0 04/12/14 03:25 1,3-Dichloropropane ug/L ND 1.0 04/12/14 03:25 1,3-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2Dichloropropane ug/L ND 1.0 04/12/14 03:25 2Dichloropropane ug/L ND 1.0 04/12/14 03:25 2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2-Butanone (MEK) ug/L ND 1.0 04/12/14 03:25 2-Hexanone ug/L ND 1.0 04/12/14 03:25 2-Hexanone ug/L	1,2,4-Trimethylbenzene		ND	1.0	04/12/14 03:25		
1,2-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,2-Dichloropthene (Total) ug/L ND 1.0 04/12/14 03:25 1,3-Dichloropthene ug/L ND 1.0 04/12/14 03:25 1,3-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,3-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 2,2-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2,2-Dichlorobluene ug/L ND 1.0 04/12/14 03:25 2,-Hexanone ug/L ND 1.0 04/12/14 03:25 4-Methyl-2-pentanone (MIBK) ug/L ND 1.0 04/12/14 03:25 Bromochloromethane ug/L	1,2-Dibromo-3-chloropropane		ND	2.5	04/12/14 03:25		
1/2-Dichloroethane ug/L ND 1.0 04/12/14 03:25 1,2-Dichloroethene (Total) ug/L ND 1.0 04/12/14 03:25 1,3-Dichloropropane ug/L ND 1.0 04/12/14 03:25 1,3-Dichloropropane ug/L ND 1.0 04/12/14 03:25 1,3-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,3-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2-Butanone (MEK) ug/L ND 1.0 04/12/14 03:25 2-Butanone (MEK) ug/L ND 1.0 04/12/14 03:25 2-Hexanone ug/L ND 1.0	1,2-Dibromoethane (EDB)	ug/L	ND	1.0	04/12/14 03:25		
1,2-Dichloroethene (Total) ug/L ND 1.0 04/12/14 03:25 1,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 1,3-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,3-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,3-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2-Chlorotoluene ug/L ND 1.0 04/12/14 03:25 2-Hexanone ug/L ND 1.0 04/12/14 03:25 2-Hexanone ug/L ND 1.0 04/12/14 03:25 4-Methyl-2-pentanone (MIBK) ug/L ND 1.0 04/12/14 03:25 4-Methyl-2-pentanone (MIBK) ug/L ND 1.0 04/12/14 03:25 Bromochoromethane ug/L ND 1.0 04/12/14 03:25 Bromochoromethane ug/L	1,2-Dichlorobenzene	ug/L	ND	1.0	04/12/14 03:25		
1,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 1,3-5-Trimethylbenzene ug/L ND 1.0 04/12/14 03:25 1,3-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,4-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropopane ug/L ND 1.0 04/12/14 03:25 2,2-Dichorobluene ug/L ND 1.0 04/12/14 03:25 2-Hexanone ug/L ND 1.0 04/12/14 03:25 Bromochoune ug/L ND 1.0 04/12	1,2-Dichloroethane	ug/L	ND	1.0	04/12/14 03:25		
1,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 1,3-5-Trimethylbenzene ug/L ND 1.0 04/12/14 03:25 1,3-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 1,4-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2-Butanone (MEK) ug/L ND 1.0 04/12/14 03:25 2-Hexanone ug/L ND 1.0 04/12/14 03:25 Bromochouree ug/L ND 1.0 04/12/14 0	1,2-Dichloroethene (Total)	ug/L	ND	1.0	04/12/14 03:25		
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1,3-Dichloropropaneug/LND1.004/12/14 03:251,4-Dichlorobenzeneug/LND1.004/12/14 03:252,2-Dichloropropaneug/LND1.004/12/14 03:252-Butanone (MEK)ug/LND1.004/12/14 03:252-Chlorotolueneug/LND1.004/12/14 03:252-Chlorotolueneug/LND1.004/12/14 03:252-Hexanoneug/LND1.004/12/14 03:254-Chlorotolueneug/LND1.004/12/14 03:254-Methyl-2-pentanone (MIBK)ug/LND1.004/12/14 03:25Acetoneug/LND1.004/12/14 03:25Benzeneug/LND1.004/12/14 03:25Bromobenzeneug/LND1.004/12/14 03:25Bromochloromethaneug/LND1.004/12/14 03:25Bromochloromethaneug/LND1.004/12/14 03:25Bromochloromethaneug/LND1.004/12/14 03:25Bromochloromethaneug/LND1.004/12/14 03:25Bromochloromethaneug/LND1.004/12/14 03:25Carbon disulfideug/LND1.004/12/14 03:25Chlorobenzeneug/LND1.004/12/14 03:25Chlorobenzeneug/LND1.004/12/14 03:25Chlorobenzeneug/LND1.004/12/14 03:25Chlorobenzeneug/LND1.004/12/14 03:25C	1,3,5-Trimethylbenzene	ug/L	ND	1.0	04/12/14 03:25		
1,4-Dichlorobenzene ug/L ND 1.0 04/12/14 03:25 2,2-Dichloropropane ug/L ND 1.0 04/12/14 03:25 2-Butanone (MEK) ug/L ND 1.0 04/12/14 03:25 2-Chlorotoluene ug/L ND 1.0 04/12/14 03:25 2-Chlorotoluene ug/L ND 1.0 04/12/14 03:25 2-Hexanone ug/L ND 1.0 04/12/14 03:25 2-Hexanone ug/L ND 1.0 04/12/14 03:25 2-Hexanone ug/L ND 1.0 04/12/14 03:25 4-Chlorotoluene ug/L ND 1.0 04/12/14 03:25 Benzene ug/L ND 1.0 04/12/14 03:25 Benzene ug/L ND 1.0 04/12/14 03:25 Bromochloromethane ug/L ND 1.0 04/12/14 03:25	1,3-Dichlorobenzene	ug/L	ND	1.0	04/12/14 03:25		
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2-Butanone (MEK) ug/L ND 10.0 04/12/14 03:25 2-Chlorotoluene ug/L ND 1.0 04/12/14 03:25 2-Hexanone ug/L ND 1.0 04/12/14 03:25 2-Hexanone ug/L ND 1.0 04/12/14 03:25 4-Chlorotoluene ug/L ND 1.0 04/12/14 03:25 4-Methyl-2-pentanone (MIBK) ug/L ND 1.0 04/12/14 03:25 Acetone ug/L ND 1.0 04/12/14 03:25 Benzene ug/L ND 1.0 04/12/14 03:25 Bromobenzene ug/L ND 1.0 04/12/14 03:25 Bromochloromethane ug/L ND 1.0 04/12/14 03:25 Bromochloromethane ug/L ND 1.0 04/12/14 03:25 Bromotofinormethane ug/L ND 1.0 04/12/14 03:25 Carbon disulfide ug/L ND 5.0 04/12/14 03:25 Carbon disulfide ug/L ND 1.0 04/12/14 03:25	1,4-Dichlorobenzene	ug/L	ND	1.0	04/12/14 03:25		
2-Chlorotoluene ug/L ND 1.0 04/12/14 03:25 2-Hexanone ug/L ND 10.0 04/12/14 03:25 4-Chlorotoluene ug/L ND 1.0 04/12/14 03:25 4-Methyl-2-pentanone (MIBK) ug/L ND 10.0 04/12/14 03:25 Acetone ug/L ND 10.0 04/12/14 03:25 Benzene ug/L ND 1.0 04/12/14 03:25 Bromobenzene ug/L ND 1.0 04/12/14 03:25 Bromochloromethane ug/L ND 1.0 04/12/14 03:25 Bromochloromethane ug/L ND 1.0 04/12/14 03:25 Bromochloromethane ug/L ND 1.0 04/12/14 03:25 Bromothare ug/L ND 1.0 04/12/14 03:25 Bromothare ug/L ND 5.0 04/12/14 03:25 Carbon disulfide ug/L ND 5.0 04/12/14 03:25 Carbon tetrachloride ug/L ND 1.0 04/12/14 03:25 Chlorobenzene ug/L ND 1.0 04/12/14 03:25<	2,2-Dichloropropane	ug/L	ND	1.0	04/12/14 03:25		
2-Hexanoneug/LND10.004/12/14 03:254-Chlorotolueneug/LND1.004/12/14 03:254-Methyl-2-pentanone (MIBK)ug/LND10.004/12/14 03:25Acetoneug/LND1.004/12/14 03:25Benzeneug/LND1.004/12/14 03:25Bromobenzeneug/LND1.004/12/14 03:25Bromochloromethaneug/LND1.004/12/14 03:25Bromodichloromethaneug/LND1.004/12/14 03:25Bromodichloromethaneug/LND1.004/12/14 03:25Bromodichloromethaneug/LND1.004/12/14 03:25Bromodichloromethaneug/LND1.004/12/14 03:25Carbon disulfideug/LND1.004/12/14 03:25Chlorobenzeneug/LND1.004/12/14 03:25Chloroformug/LND1.004/12/14 03:25Chloroformug/LND1.004/12/14 03:25Chloroformug/LND1.004/12/14 03:25Chloromethaneug/LND1.004/12/14 03:25Chloroformug/LND1.004/12/14 03:25Chloromethaneug/LND1.004/12/14 03:25Chloroformug/LND1.004/12/14 03:25Chloromethaneug/LND1.004/12/14 03:25Chloromethaneug/LND1.004/12/14 03:25Chloromethaneug/L<	2-Butanone (MEK)	ug/L	ND	10.0	04/12/14 03:25		
4-Chlorotolueneug/LND1.004/12/14 03:254-Methyl-2-pentanone (MIBK)ug/LND10.004/12/14 03:25Acetoneug/LND1.004/12/14 03:25Benzeneug/LND1.004/12/14 03:25Bromobenzeneug/LND1.004/12/14 03:25Bromochloromethaneug/LND1.004/12/14 03:25Bromodichloromethaneug/LND1.004/12/14 03:25Bromodichloromethaneug/LND1.004/12/14 03:25Bromodichloromethaneug/LND1.004/12/14 03:25Bromodichloromethaneug/LND1.004/12/14 03:25Bromodethaneug/LND1.004/12/14 03:25Carbon disulfideug/LND5.004/12/14 03:25Chlorobenzeneug/LND1.004/12/14 03:25Chloroformug/LND1.004/12/14 03:25Chloroformug/LND1.004/12/14 03:25Chloroformug/LND1.004/12/14 03:25Chloroformug/LND1.004/12/14 03:25Chloromethaneug/LND1.004/12/14 03:25Chloromethaneug/LND1.004/12/14 03:25Chloroformug/LND1.004/12/14 03:25Chloromethaneug/LND1.004/12/14 03:25Chloromethaneug/LND1.004/12/14 03:25Cis-1,2-Dichloropropene <td>2-Chlorotoluene</td> <td>ug/L</td> <td>ND</td> <td>1.0</td> <td>04/12/14 03:25</td> <td></td>	2-Chlorotoluene	ug/L	ND	1.0	04/12/14 03:25		
4-Methyl-2-pentanone (MIBK)ug/LND10.004/12/14 03:25Acetoneug/LND1.004/12/14 03:25Benzeneug/LND1.004/12/14 03:25Bromobenzeneug/LND1.004/12/14 03:25Bromochloromethaneug/LND1.004/12/14 03:25Bromodichloromethaneug/LND1.004/12/14 03:25Bromodichloromethaneug/LND1.004/12/14 03:25Bromodichloromethaneug/LND1.004/12/14 03:25Bromoformug/LND5.004/12/14 03:25Bromomethaneug/LND5.004/12/14 03:25Carbon disulfideug/LND1.004/12/14 03:25Chlorobenzeneug/LND1.004/12/14 03:25Chloroethaneug/LND1.004/12/14 03:25Chloroformug/LND1.004/12/14 03:25Chloroformug/LND1.004/12/14 03:25Chloromethaneug/LND1.004/12/14 03:25Chloroformug/LND1.004/12/14 03:25Chloromethaneug/LND1.004/12/14 03:25Chloroformug/LND1.004/12/14 03:25Chloroformug/LND1.004/12/14 03:25Chloropetheneug/LND1.004/12/14 03:25Chloroformug/LND1.004/12/14 03:25Chloropetheneug/LND<	2-Hexanone	ug/L	ND	10.0	04/12/14 03:25		
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Chloroform ug/L ND 1.0 04/12/14 03:25 Chloromethane ug/L ND 1.0 04/12/14 03:25 cis-1,2-Dichloroethene ug/L ND 1.0 04/12/14 03:25 cis-1,3-Dichloropropene ug/L ND 1.0 04/12/14 03:25							
Chloromethane ug/L ND 1.0 04/12/14 03:25 cis-1,2-Dichloroethene ug/L ND 1.0 04/12/14 03:25 cis-1,3-Dichloropropene ug/L ND 1.0 04/12/14 03:25							
cis-1,2-Dichloroethene ug/L ND 1.0 04/12/14 03:25 cis-1,3-Dichloropropene ug/L ND 1.0 04/12/14 03:25		-					
cis-1,3-Dichloropropene ug/L ND 1.0 04/12/14 03:25							
	,						
Dibromochloromethane ug/l ND 1.0 04/12/14 03:25							
	Dibromochloromethane	ug/L	ND	1.0	04/12/14 03:25		



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

Associated Lab Samples: 60166332001

Matrix: Water

Parameter	Units	Blank Result	Reporting Limit	Analyzed	Qualifiers
Dibromomethane	ug/L		1.0	04/12/14 03:25	
Dichlorodifluoromethane	ug/L	ND	1.0	04/12/14 03:25	
Ethylbenzene	ug/L	ND	1.0	04/12/14 03:25	
Hexachloro-1,3-butadiene	ug/L	ND	1.0	04/12/14 03:25	
Isopropylbenzene (Cumene)	ug/L	ND	1.0	04/12/14 03:25	
Methyl-tert-butyl ether	ug/L	ND	1.0	04/12/14 03:25	
Methylene chloride	ug/L	ND	1.0	04/12/14 03:25	
n-Butylbenzene	ug/L	ND	1.0	04/12/14 03:25	
n-Propylbenzene	ug/L	ND	1.0	04/12/14 03:25	
Naphthalene	ug/L	ND	10.0	04/12/14 03:25	
p-Isopropyltoluene	ug/L	ND	1.0	04/12/14 03:25	
sec-Butylbenzene	ug/L	ND	1.0	04/12/14 03:25	
Styrene	ug/L	ND	1.0	04/12/14 03:25	
tert-Butylbenzene	ug/L	ND	1.0	04/12/14 03:25	
Tetrachloroethene	ug/L	ND	1.0	04/12/14 03:25	
Toluene	ug/L	ND	1.0	04/12/14 03:25	
trans-1,2-Dichloroethene	ug/L	ND	1.0	04/12/14 03:25	
trans-1,3-Dichloropropene	ug/L	ND	1.0	04/12/14 03:25	
Trichloroethene	ug/L	ND	1.0	04/12/14 03:25	
Trichlorofluoromethane	ug/L	ND	1.0	04/12/14 03:25	
Vinyl chloride	ug/L	ND	1.0	04/12/14 03:25	
Xylene (Total)	ug/L	ND	3.0	04/12/14 03:25	
1,2-Dichloroethane-d4 (S)	%	106	80-120	04/12/14 03:25	
4-Bromofluorobenzene (S)	%	96	80-120	04/12/14 03:25	
Toluene-d8 (S)	%	107	80-120	04/12/14 03:25	

LABORATORY CONTROL SAMPLE: 1358976

LABORATORY CONTROL SAMPLE:	1358976					
		Spike	LCS	LCS	% Rec	
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers
1,1,1,2-Tetrachloroethane	ug/L	20	19.5	98	80-124	
1,1,1-Trichloroethane	ug/L	20	17.3	87	80-121	
1,1,2,2-Tetrachloroethane	ug/L	20	18.9	94	73-124	
1,1,2-Trichloroethane	ug/L	20	17.9	90	80-120	
1,1-Dichloroethane	ug/L	20	16.4	82	77-120	
1,1-Dichloroethene	ug/L	20	17.6	88	78-126	
1,1-Dichloropropene	ug/L	20	17.5	87	80-120	
1,2,3-Trichlorobenzene	ug/L	20	20.1	101	75-130	
1,2,3-Trichloropropane	ug/L	20	20.8	104	76-127	
1,2,4-Trichlorobenzene	ug/L	20	20.9	105	79-124	
1,2,4-Trimethylbenzene	ug/L	20	21.8	109	80-122	
1,2-Dibromo-3-chloropropane	ug/L	20	20.4	102	68-131	
1,2-Dibromoethane (EDB)	ug/L	20	20.7	103	80-127	
1,2-Dichlorobenzene	ug/L	20	21.1	105	80-122	
1,2-Dichloroethane	ug/L	20	17.7	88	77-123	
1,2-Dichloroethene (Total)	ug/L	40	37.1	93	80-120	



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

LABORATORY CONTROL SAMPLE: 1358976

Parameter	Units	Spike Conc.	LCS Result	LCS % Rec	% Rec Limits	Qualifiers	
1,2-Dichloropropane	ug/L		18.6	93	80-121		
1,3,5-Trimethylbenzene	ug/L	20	20.8	104	80-121		
1,3-Dichlorobenzene	ug/L	20	20.8	104	80-120		
1,3-Dichloropropane	ug/L	20	18.8	94	80-120		
1,4-Dichlorobenzene	ug/L	20	21.0	105	80-120		
2,2-Dichloropropane	ug/L	20	13.9	70	50-137		
2-Butanone (MEK)	ug/L	100	99.1	99	52-145		
2-Chlorotoluene	ug/L	20	19.9	99	80-120		
2-Hexanone	ug/L	100	104	104	57-139		
4-Chlorotoluene	ug/L	20	21.9	109	80-121		
1-Methyl-2-pentanone (MIBK)	ug/L	100	94.2	94	71-131		
Acetone	ug/L	100	104	104	32-155		
Benzene	ug/L	20	17.1	85	80-120		
Bromobenzene	ug/L	20	21.1	106	80-120		
Bromochloromethane	ug/L	20	16.0	80	77-123		
Bromodichloromethane	ug/L	20	17.0	85	80-120		
Bromoform	ug/L	20	20.2	101	73-124		
Bromomethane	ug/L	20	13.6	68	31-144		
Carbon disulfide	ug/L	20	17.5	87	65-125		
Carbon tetrachloride	ug/L	20	17.6	88	78-128		
chlorobenzene	ug/L	20	19.6	98	80-120		
hloroethane	ug/L	20	18.9	94	55-137		
hloroform	ug/L	20	17.6	88	79-120		
hloromethane	ug/L	20	21.6	108	22-138		
s-1,2-Dichloroethene	ug/L	20	18.4	92	80-120		
s-1,3-Dichloropropene	ug/L	20	16.2	81	80-120		
ibromochloromethane	ug/L	20	17.6	88	80-120		
ibromomethane	ug/L	20	18.5	93	80-122		
lichlorodifluoromethane	ug/L	20	15.1	75	23-120		
thylbenzene	ug/L	20	19.0	95	80-121		
exachloro-1,3-butadiene	ug/L	20	20.2	101	77-129		
opropylbenzene (Cumene)	ug/L	20	21.5	107	80-136		
lethyl-tert-butyl ether	ug/L	20	15.7	79	74-125		
lethylene chloride	ug/L	20	18.8	94	73-126		
-Butylbenzene	ug/L	20	20.6	103	83-123		
-Propylbenzene	ug/L	20	22.2	111	80-122		
laphthalene	ug/L	20	20.2	101	73-130		
-Isopropyltoluene	ug/L	20	22.0	110	80-124		
ec-Butylbenzene	ug/L	20	20.8	104	80-129		
ityrene	ug/L	20	20.3	101	80-120		
ert-Butylbenzene	ug/L	20	20.8	104	80-126		
etrachloroethene	ug/L	20	19.5	98	80-121		
oluene	ug/L	20	19.3	96	80-122		
rans-1,2-Dichloroethene	ug/L	20	18.7	93	79-121		
rans-1,3-Dichloropropene	ug/L	20	19.2	96	80-127		
Trichloroethene	ug/L	20	18.1	90	80-120		
Trichlorofluoromethane	ug/L	20	17.7	88	67-120		
	0	-			59-120		



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

LABORATORY CONTROL SAMPLE:	1358976					
		Spike	LCS	LCS	% Rec	
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers
Xylene (Total)	ug/L	60	59.9	100	80-121	
1,2-Dichloroethane-d4 (S)	%			102	80-120	
4-Bromofluorobenzene (S)	%			100	80-120	
Toluene-d8 (S)	%			104	80-120	



Project: 085692 MARCC	TTE NO 1							
Pace Project No.: 60166332								
QC Batch: OEXT/43505		Analysis	Method:	EF	PA 8015B			
QC Batch Method: EPA 3510C		Analysis	Description	n: EF	PA 8015B			
Associated Lab Samples: 6016633	32001							
METHOD BLANK: 1356074		Ма	trix: Water					
Associated Lab Samples: 6016633	32001							
		Blank	Rep	orting				
Parameter	Units	Result	Li	mit	Analyzed	l Qualif	iers	
TPH-DRO	mg/L		ND	0.50	04/07/14 15	:55		
n-Tetracosane (S)	%		66	22-121	04/07/14 15	:55		
p-Terphenyl (S)	%		68	28-127	04/07/14 15	:55		
LABORATORY CONTROL SAMPLE:	1356075							
		Spike	LCS		LCS	% Rec		
Parameter	Units	Conc.	Result	0	% Rec	Limits	Qualifiers	
TPH-DRO	mg/L	12.5		8.1	65	39-120		
n-Tetracosane (S)	%				71	22-121		
p-Terphenyl (S)	%				73	28-127		



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332					
QC Batch: OEXT/43560)	Analysis Meth	nod: EF	PA 8270	
QC Batch Method: EPA 3510		Analysis Dese	cription: 82	70 Water MSSV	
Associated Lab Samples: 60166	332001				
METHOD BLANK: 1357256		Matrix:	Water		
Associated Lab Samples: 60166	332001				
		Blank	Reporting		
Parameter	Units	Result	Limit	Analyzed	Qualifiers
1,2,4-Trichlorobenzene	ug/L	 ND	10.0	04/10/14 13:00	
1,2-Dichlorobenzene	ug/L	ND	10.0	04/10/14 13:00	
1,3-Dichlorobenzene	ug/L	ND	10.0	04/10/14 13:00	
1,4-Dichlorobenzene	ug/L	ND	10.0	04/10/14 13:00	
2,4,5-Trichlorophenol	ug/L	ND	50.0	04/10/14 13:00	
2,4,6-Trichlorophenol	ug/L	ND	10.0	04/10/14 13:00	
2,4-Dichlorophenol	ug/L	ND	10.0	04/10/14 13:00	
2,4-Dimethylphenol	ug/L	ND	10.0	04/10/14 13:00	
2,4-Dinitrophenol	ug/L	ND	50.0	04/10/14 13:00	
2,4-Dinitrotoluene	ug/L	ND	10.0	04/10/14 13:00	
2.6-Dinitrotoluene	ua/l		10.0	04/10/14 13:00	

1,4-Dichlorobenzene	ug/L	ND	10.0	04/10/14 13:00	
2,4,5-Trichlorophenol	ug/L	ND	50.0	04/10/14 13:00	
2,4,6-Trichlorophenol	ug/L	ND	10.0	04/10/14 13:00	
2,4-Dichlorophenol	ug/L	ND	10.0	04/10/14 13:00	
2,4-Dimethylphenol	ug/L	ND	10.0	04/10/14 13:00	
2,4-Dinitrophenol	ug/L	ND	50.0	04/10/14 13:00	
2,4-Dinitrotoluene	ug/L	ND	10.0	04/10/14 13:00	
2,6-Dinitrotoluene	ug/L	ND	10.0	04/10/14 13:00	
2-Chloronaphthalene	ug/L	ND	10.0	04/10/14 13:00	
2-Chlorophenol	ug/L	ND	10.0	04/10/14 13:00	
2-Methylnaphthalene	ug/L	ND	10.0	04/10/14 13:00	
2-Methylphenol(o-Cresol)	ug/L	ND	10.0	04/10/14 13:00	
2-Nitroaniline	ug/L	ND	50.0	04/10/14 13:00	
2-Nitrophenol	ug/L	ND	10.0	04/10/14 13:00	
3&4-Methylphenol(m&p Cresol)	ug/L	ND	10.0	04/10/14 13:00	
3,3'-Dichlorobenzidine	ug/L	ND	20.0	04/10/14 13:00	
3-Nitroaniline	ug/L	ND	50.0	04/10/14 13:00	
4,6-Dinitro-2-methylphenol	ug/L	ND	50.0	04/10/14 13:00	
4-Bromophenylphenyl ether	ug/L	ND	10.0	04/10/14 13:00	
4-Chloro-3-methylphenol	ug/L	ND	20.0	04/10/14 13:00	
4-Chloroaniline	ug/L	ND	20.0	04/10/14 13:00	
4-Chlorophenylphenyl ether	ug/L	ND	10.0	04/10/14 13:00	
4-Nitroaniline	ug/L	ND	50.0	04/10/14 13:00	
4-Nitrophenol	ug/L	ND	50.0	04/10/14 13:00	
Acenaphthene	ug/L	ND	10.0	04/10/14 13:00	
Acenaphthylene	ug/L	ND	10.0	04/10/14 13:00	
Anthracene	ug/L	ND	10.0	04/10/14 13:00	
Benzo(a)anthracene	ug/L	ND	10.0	04/10/14 13:00	
Benzo(a)pyrene	ug/L	ND	10.0	04/10/14 13:00	
Benzo(b)fluoranthene	ug/L	ND	10.0	04/10/14 13:00	
Benzo(g,h,i)perylene	ug/L	ND	10.0	04/10/14 13:00	
Benzo(k)fluoranthene	ug/L	ND	10.0	04/10/14 13:00	
Benzoic acid	ug/L	ND	50.0	04/10/14 13:00	
Benzyl alcohol	ug/L	ND	20.0	04/10/14 13:00	
bis(2-Chloroethoxy)methane	ug/L	ND	10.0	04/10/14 13:00	
bis(2-Chloroethyl) ether	ug/L	ND	10.0	04/10/14 13:00	
bis(2-Chloroisopropyl) ether	ug/L	ND	10.0	04/10/14 13:00	
bis(2-Ethylhexyl)phthalate	ug/L	ND	10.0	04/10/14 13:00	
Butylbenzylphthalate	ug/L	ND	10.0	04/10/14 13:00	
Carbazole	ug/L	ND	10.0	04/10/14 13:00	

REPORT OF LABORATORY ANALYSIS

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Matrix: Water

Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

METHOD BLANK: 1357256	
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Associated Lab Samples: 60166332001

Blank Reporting Parameter Units Result Limit Analyzed Qualifiers Chrysene ug/L ND 10.0 04/10/14 13:00 ug/L Di-n-butylphthalate ND 10.0 04/10/14 13:00 Di-n-octylphthalate ND 10.0 04/10/14 13:00 ug/L Dibenz(a,h)anthracene ug/L ND 10.0 04/10/14 13:00 ug/L 10.0 04/10/14 13:00 Dibenzofuran ND Diethylphthalate ug/L ND 10.0 04/10/14 13:00 Dimethylphthalate ug/L ND 10.0 04/10/14 13:00 Fluoranthene ug/L ND 10.0 04/10/14 13:00 Fluorene ug/L ND 10.0 04/10/14 13:00 Hexachloro-1,3-butadiene ND 10.0 04/10/14 13:00 ug/L Hexachlorobenzene ug/L ND 10.0 04/10/14 13:00 Hexachlorocyclopentadiene ug/L ND 10.0 04/10/14 13:00 Hexachloroethane ug/L ND 10.0 04/10/14 13:00 ND Indeno(1,2,3-cd)pyrene ug/L 10.0 04/10/14 13:00 ND Isophorone ug/L 10.0 04/10/14 13:00 N-Nitroso-di-n-propylamine ND ug/L 10.0 04/10/14 13:00 N-Nitrosodiphenylamine ug/L ND 10.0 04/10/14 13:00 Naphthalene ug/L ND 10.0 04/10/14 13:00 Nitrobenzene ug/L ND 10.0 04/10/14 13:00 Pentachlorophenol ug/L ND 50.0 04/10/14 13:00 ND Phenanthrene ug/L 10.0 04/10/14 13:00 ND Phenol ug/L 10.0 04/10/14 13:00 Pyrene ug/L ND 10.0 04/10/14 13:00 Pyridine ug/L ND 10.0 04/10/14 13:00 2,4,6-Tribromophenol (S) % 89 29-121 04/10/14 13:00 2-Fluorobiphenyl (S) % 93 19-124 04/10/14 13:00 % 2-Fluorophenol (S) 46 13-120 04/10/14 13:00 % Nitrobenzene-d5 (S) 90 10-135 04/10/14 13:00 Phenol-d6 (S) % 29 10-120 04/10/14 13:00 Terphenyl-d14 (S) % 95 24-131 04/10/14 13:00

LABORATORY CONTROL SAMPLE: 1357257

Parameter	Units	Spike Conc.	LCS Result	LCS % Rec	% Rec Limits	Qualifiers
1,2,4-Trichlorobenzene	ug/L	50	42.6	85	45-120	
1,2-Dichlorobenzene	ug/L	50	42.5	85	45-120	
1,3-Dichlorobenzene	ug/L	50	40.9	82	44-120	
1,4-Dichlorobenzene	ug/L	50	41.6	83	44-120	
2,4,5-Trichlorophenol	ug/L	50	45.6J	91	50-120	
2,4,6-Trichlorophenol	ug/L	50	44.4	89	49-120	
2,4-Dichlorophenol	ug/L	50	42.2	84	48-120	
2,4-Dimethylphenol	ug/L	50	38.8	78	35-120	
2,4-Dinitrophenol	ug/L	50	31.9J	64	21-120	
2,4-Dinitrotoluene	ug/L	50	46.3	93	52-120	
2,6-Dinitrotoluene	ug/L	50	45.2	90	53-120	

REPORT OF LABORATORY ANALYSIS

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Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

LABORATORY CONTROL SAMPLE: 1357257

Parameter	Units	Spike Conc.	LCS Result	LCS % Rec	% Rec Limits	Qualifiers
2-Chloronaphthalene	ug/L	50	45.1	90	49-120	
2-Chlorophenol	ug/L	50	39.5	79	47-120	
2-Methylnaphthalene	ug/L	50	44.6	89	46-120	
2-Methylphenol(o-Cresol)	ug/L	50	36.3	73	40-120	
2-Nitroaniline	ug/L	50	46.1J	92	51-120	
-Nitrophenol	ug/L	50	42.6	85	47-120	
&4-Methylphenol(m&p Cresol)	ug/L	50	33.1	66	34-120	
,3'-Dichlorobenzidine	ug/L	50	74.0	148	28-160	
-Nitroaniline	ug/L	50	51.6	103	43-157	
,6-Dinitro-2-methylphenol	ug/L	50	48.5J	97	42-120	
-Bromophenylphenyl ether	ug/L	50	47.0	94	52-120	
-Chloro-3-methylphenol	ug/L	50	41.8	84	48-120	
-Chloroaniline	ug/L	50	57.2	114	24-160	
-Chlorophenylphenyl ether	ug/L	50	45.4	91	53-120	
Nitroaniline	ug/L	50	46.7J	93	50-120	
-Nitrophenol	ug/L	50	16.8J	34	10-120	
cenaphthene	ug/L	50	45.3	91	50-120	
cenaphthylene	ug/L	50	43.8	88	49-120	
nthracene	ug/L	50	46.4	93	52-120	
enzo(a)anthracene	ug/L	50	47.9	96	53-120	
enzo(a)pyrene	ug/L	50	45.2	90	51-120	
enzo(b)fluoranthene	ug/L	50	44.7	89	51-120	
enzo(g,h,i)perylene	ug/L	50	46.7	93	52-120	
enzo(g,i,i,)pergiene	ug/L	50 50	49.1	98	51-120	
nzoic acid	ug/L	50	ND	7	10-120 L	0
enzyl alcohol	ug/L	50	34.4	69	39-120	_0
s(2-Chloroethoxy)methane	ug/L	50	44.3	89	50-120	
s(2-Chloroethyl) ether	ug/L	50 50	44.6	89	48-120	
s(2-Chloroisopropyl) ether	ug/L	50 50	43.3	87	49-120	
s(2-Ethylhexyl)phthalate	ug/L	50 50	49.4	99	52-123	
utylbenzylphthalate	ug/L	50	49.4	99	52-120	
arbazole	ug/L	50 50	49.4 46.9	99 94	55-120	
hrysene	ug/L ug/L	50 50	40.9	94 98	53-120	
i-n-butylphthalate	ug/L ug/L	50 50	49.1 48.4	98 97	49-125	
i-n-octylphthalate	ug/L ug/L	50 50	46.4 46.6	97 93	49-125 51-121	
ibenz(a,h)anthracene	ug/L ug/L	50 50	40.0	93	51-121	
ibenzofuran	ug/L ug/L	50 50	47.0	94 92	51-120	
iethylphthalate	ug/L ug/L	50 50	45.8 46.2	92 92	53-120	
imethylphthalate	ug/L	50 50	46.0	92	52-120 52-120	
uoranthene	ug/L	50 50	40.0	92 95	53-120	
uorantnene	ug/L ug/L	50 50	47.7 44.9	95 90	53-120 52-120	
exachloro-1,3-butadiene		50 50	44.9 40.2	90 80	52-120 42-120	
exachloro-1,3-butadiene exachlorobenzene	ug/L	50 50	40.2 44.6	80 89	42-120 52-120	
	ug/L					
exachlorocyclopentadiene exachloroethane	ug/L	100 50	75.4 39.1	75 78	26-120 43-120	
ndeno(1,2,3-cd)pyrene	ug/L	50 50	47.3	78 95	43-120 51-120	
	ug/L					
sophorone	ug/L	50 50	43.4	87	50-120	
N-Nitroso-di-n-propylamine	ug/L	50	45.5	91	50-120	



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

LABORATORY CONTROL SAMPLE: 1357257

Parameter	Units	Spike Conc.	LCS Result	LCS % Rec	% Rec Limits	Qualifiers
N-Nitrosodiphenylamine	ug/L		44.9	90	53-120	
Naphthalene	ug/L	50	43.6	87	48-120	
Nitrobenzene	ug/L	50	44.3	89	47-120	
Pentachlorophenol	ug/L	50	40.3J	81	43-120	
Phenanthrene	ug/L	50	45.1	90	53-120	
Phenol	ug/L	50	18.2	36	12-120	
Pyrene	ug/L	50	48.1	96	54-120	
Pyridine	ug/L	50	17.2	34	10-120	
2,4,6-Tribromophenol (S)	%			91	29-121	
2-Fluorobiphenyl (S)	%			90	19-124	
2-Fluorophenol (S)	%			44	13-120	
Nitrobenzene-d5 (S)	%			89	10-135	
Phenol-d6 (S)	%			30	10-120	
Terphenyl-d14 (S)	%			98	24-131	



Project: 0	85692 MARCO	TE NO 1					
Pace Project No.: 6	0166332						
QC Batch:	WET/47146		Analysis Meth	nod: EF	PA 120.1		
QC Batch Method:	EPA 120.1		Analysis Des	Analysis Description: 120.1 Specific Condu		uctance	
Associated Lab Samp	les: 60166332	2001					
METHOD BLANK: 1	356310		Matrix:	Water			
Associated Lab Samp	les: 60166332	2001					
			Blank	Reporting			
Parame	ter	Units	Result	Limit	Analyzed	Qualifiers	_
Specific Conductance		umhos/cm	ND	1.0	04/08/14 10:11		
SAMPLE DUPLICATE	: 1356311						
			60166384001	Dup		Max	
Parame	ter	Units	Result	Result	RPD	RPD	Qualifiers
Specific Conductance		umhos/cm	ND	ND		20	



Project: 0	85692 MARCO	TTE NO 1						
Pace Project No.: 6	0166332							
QC Batch:	WET/47254		Analysis M	Method:	SM 2320B			
QC Batch Method:	SM 2320B		Analysis [Description:	2320B Alkalinit	у		
Associated Lab Samp	les: 6016633	32001						
METHOD BLANK: 1	358655		Mati	rix: Water				
Associated Lab Samp	les: 6016633	32001						
			Blank	Reporting				
Parame	ter	Units	Result	Limit	Analyze	d Quali	ifiers	_
Alkalinity, Total as Ca	03	mg/L	N	ID 20	0.0 04/14/14 1 ⁻	1:40		
LABORATORY CONT	ROL SAMPLE:	1358656						
			Spike	LCS	LCS	% Rec		
Parame	ter	Units	Conc.	Result	% Rec	Limits	Qua	alifiers
Alkalinity, Total as Ca	03	mg/L	500	537	107	90-110		
SAMPLE DUPLICATE	: 1358661							
			6016667500			Max		
Parame	ter	Units	Result	Result	RPD	RPD		Qualifiers
Alkalinity, Total as Ca0	03	mg/L	88	36 8	85	0	10	
SAMPLE DUPLICATE	: 1358665							
			6016663800		555	Max		0.11
			Result	Result	RPD	RPD		Qualifiers
Parame	ter	Units						



Project: 0856	92 MARCOTTE NO 1						
Pace Project No.: 6016	6332						
QC Batch: WE	T/47185	Analysis	Method:	SM 2540C			
QC Batch Method: SN	atch Method: SM 2540C		Description:	2540C Total Di	ssolved Solids		
Associated Lab Samples:	60166332001						
METHOD BLANK: 1356	5722	Ma	atrix: Water				
Associated Lab Samples:	60166332001						
		Blank	Reporting)			
Parameter	Uni	ts Result	Limit	Analyze	d Qual	fiers	
Total Dissolved Solids	mg/L		ND	5.0 04/08/14 1	2:33		
LABORATORY CONTRO	L SAMPLE: 1356723						
		Spike	LCS	LCS	% Rec		
Parameter	Uni	ts Conc.	Result	% Rec	Limits	Qu	alifiers
Total Dissolved Solids	mg/L	1000	987	99	80-120		
SAMPLE DUPLICATE:	1356724						
		601660970	-		Max		
Parameter	Uni	ts Result	Result	RPD	RPD		Qualifiers
Total Dissolved Solids	mg/L	:	389 3	380	2	10	
SAMPLE DUPLICATE:	1356725						
		601663820	-	000	Max		0 ""
Parameter	Uni	ts Result	Result	RPD	RPD		Qualifiers
Total Dissolved Solids			788 7	791			



Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

QC Batch: W	'ET/47180	Analysis Meth	od:	SM 4500-H+B		
QC Batch Method: SI	M 4500-H+B	Analysis Desc	ription:	4500H+B pH		
Associated Lab Samples	s: 60166332001					
SAMPLE DUPLICATE:	1356696					
SAMPLE DUPLICATE:	1356696	60166120001	Dup		Max	
SAMPLE DUPLICATE: Parameter		60166120001 Result	Dup Result	RPD	Max RPD	Qualifiers



Project: 085692 MARCOTTE NO 1

Pace Project No .:	60166332
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QC Batch: WETA	/29042		Analys	is Method	: E	PA 300.0						
QC Batch Method: EPA 3	00.0		Analys	is Descrip	tion: 30	0.0 IC Anio	ons					
Associated Lab Samples:	60166332001											
METHOD BLANK: 136197	1		Ν	Aatrix: Wa	iter							
Associated Lab Samples:	60166332001											
			Blank	K F	Reporting							
Parameter	l	Jnits	Resul	t	Limit	Analyz	ed	Qualifiers				
Bromide	mg/L			ND	1.0	04/17/14	19:41					
Chloride	mg/L			ND	1.0		-					
Fluoride	mg/L			ND	0.20		-					
Sulfate	mg/L			ND	1.0	04/17/14	19:41					
LABORATORY CONTROL S	SAMPLE: 13619	72										
_			Spike	LC		LCS	% Rec					
Parameter	l	Jnits	Conc.	Res	ult	% Rec	Limits	Qı	ualifiers			
	mg/L		5		4.7	94)-110				
Chloride	mg/L		5		4.7	93	90)-110				
Chloride Fluoride	mg/L mg/L		5 2.5		4.7 2.3	93 92	90 90)-110)-110				
Chloride Fluoride	mg/L		5		4.7	93	90 90)-110				
Chloride Fluoride Sulfate	mg/L mg/L mg/L		5 2.5 5		4.7 2.3 4.9	93 92	90 90)-110)-110				
Chloride Fluoride Sulfate	mg/L mg/L mg/L	: 13619	5 2.5 5 73		4.7 2.3	93 92	90 90)-110)-110				
Chloride Fluoride Sulfate	mg/L mg/L mg/L SPIKE DUPLICATE		5 2.5 5 73 MS	MSD	4.7 2.3 4.9 1361974	93 92 98	90 90 90)-110)-110)-110	% Poc		Max	
Chloride Fluoride Sulfate	mg/L mg/L mg/L SPIKE DUPLICATE	:: 13619 66574001 Result	5 2.5 5 73		4.7 2.3 4.9	93 92	90 90)-110)-110	% Rec Limits	RPD	Max RPD	Qu
Chloride Fluoride Sulfate MATRIX SPIKE & MATRIX S Parameter	mg/L mg/L SPIKE DUPLICATE 6010	66574001 Result	5 2.5 5 73 MS Spike Conc.	MSD Spike Conc.	4.7 2.3 4.9 1361974 MS Result	93 92 98 MSD Result	90 90 90 90 90 90)-110)-110)-110)-110 MSD % Rec	Limits	RPD 3	RPD	Qu
Chloride Fluoride Sulfate MATRIX SPIKE & MATRIX S Parameter Bromide	mg/L mg/L mg/L SPIKE DUPLICATE 6016 Units mg/L	66574001 Result ND	5 2.5 5 73 MS Spike Conc. 2500	MSD Spike Conc. 2500	4.7 2.3 4.9 1361974 MS Result 2490	93 92 98 MSD Result 2560	90 90 90 90 90 90 MS % Rec 99)-110)-110)-110)-110 MSD <u>% Rec</u> 102	Limits 80-120	3	RPD 15	Qu
Bromide Chloride Fluoride Sulfate MATRIX SPIKE & MATRIX S Parameter Bromide Chloride Fluoride	mg/L mg/L SPIKE DUPLICATE 6010	66574001 Result	5 2.5 5 73 MS Spike Conc.	MSD Spike Conc.	4.7 2.3 4.9 1361974 MS Result	93 92 98 MSD Result	90 90 90 90 90 90)-110)-110)-110)-110 MSD % Rec	Limits		RPD	Qu

MATRIX SPIKE SAMPLE:	1361975						
		60166574005	Spike	MS	MS	% Rec	
Parameter	Units	Result	Conc.	Result	% Rec	Limits	Qualifiers
Bromide	mg/L	ND	2500	2500	100	80-120	
Chloride	mg/L	5330	2500	8020	107	80-120	
Fluoride	mg/L	ND	1250	1210	97	80-120	
Sulfate	mg/L	ND	2500	2570	88	80-120	



Project: 085692 M	ARCOTTE NO 1						
Pace Project No.: 60166332							
QC Batch: WETA/2	8919	Analysis M	ethod:	EPA 353.2			
QC Batch Method: EPA 353	3.2	Analysis D	escription:	353.2 Nitrate + N	Nitrite, preserve	d	
Associated Lab Samples: 60	0166332001						
METHOD BLANK: 1356711		Matri	x: Water				
Associated Lab Samples: 60	0166332001						
		Blank	Reporting				
Parameter	Units	Result	Limit	Analyzed	Qualifi	ers	
Nitrogen, NO2 plus NO3	mg/L	NE	0.	10 04/08/14 15	:56		
LABORATORY CONTROL SA	MPLE: 1356712						
_		Spike	LCS	LCS	% Rec		
Parameter	Units	Conc.	Result	% Rec	Limits	Qualifiers	
Nitrogen, NO2 plus NO3	mg/L	2	2.1	104	90-110		
MATRIX SPIKE SAMPLE:	1356713						
MATRIX OF IRE OAMIT EE.	1350715	6016622900	01 Spike	MS	MS	% Rec	
Parameter	Units	Result	Conc.	Result	% Rec	Limits	Qualifiers
Nitrogen, NO2 plus NO3	mg/L		1.7 2	2 3.6	92	90-110	
SAMPLE DUPLICATE: 1356	714						
Gram LE DOI LIGATE. 1990	117	60166293003	Dup		Max		
Parameter	Units	Result	Result	RPD	RPD	Qualifiers	
Nitrogen, NO2 plus NO3	mg/L	N	0.	13		20	-



Project: 08569	2 MARCOTTE NO 1						
Pace Project No.: 60166	332						
QC Batch: WET	A/28883	Analysis Me	ethod:	EPA 365.1			
QC Batch Method: EPA	365.1	Analysis De	scription:	365.1 Orthophos	sphate as P		
Associated Lab Samples:	60166332001						
METHOD BLANK: 13552	56	Matrix	: Water				
Associated Lab Samples:	60166332001						
		Blank	Reporting				
Parameter	Units	Result	Limit	Analyzed	Qualif	iers	
Orthophosphate as P	mg/L	ND	0.7	10 04/04/14 12	:49		
LABORATORY CONTROL	SAMPLE: 1355257						
LABORATORT CONTROL	5AWFLL. 1555257	Spike	LCS	LCS	% Rec		
Parameter	Units	•	Result	% Rec	Limits	Qualifiers	
Orthophosphate as P	mg/L	2	1.9	94	90-110		
MATRIX SPIKE SAMPLE:	1355311						
		6016633200	1 Spike	MS	MS	% Rec	
Parameter	Units	Result	Conc.	Result	% Rec	Limits	Qualifiers
Orthophosphate as P	mg/L	0	.10 2	1.9	9	90-110	
SAMPLE DUPLICATE: 1	355258						
		60166320001	Dup		Max		
Parameter	Units	Result	Result	RPD	RPD	Qualifiers	
i ulumeter							



QUALIFIERS

Project: 085692 MARCOTTE NO 1

Pace Project No.: 60166332

DEFINITIONS

DF - Dilution Factor, if reported, represents the factor applied to the reported data due to changes in sample preparation, dilution of the sample aliquot, or moisture content.

ND - Not Detected at or above adjusted reporting limit.

J - Estimated concentration above the adjusted method detection limit and below the adjusted reporting limit.

MDL - Adjusted Method Detection Limit.

PRL - Pace Reporting Limit.

RL - Reporting Limit.

S - Surrogate

1,2-Diphenylhydrazine (8270 listed analyte) decomposes to Azobenzene.

Consistent with EPA guidelines, unrounded data are displayed and have been used to calculate % recovery and RPD values.

LCS(D) - Laboratory Control Sample (Duplicate)

MS(D) - Matrix Spike (Duplicate)

DUP - Sample Duplicate

RPD - Relative Percent Difference

NC - Not Calculable.

SG - Silica Gel - Clean-Up

U - Indicates the compound was analyzed for, but not detected.

N-Nitrosodiphenylamine decomposes and cannot be separated from Diphenylamine using Method 8270. The result reported for each analyte is a combined concentration.

Pace Analytical is TNI accredited. Contact your Pace PM for the current list of accredited analytes.

TNI - The NELAC Institute.

BATCH QUALIFIERS

Batch: OEXT/43505

[M5] A matrix spike/matrix spike duplicate was not performed for this batch due to insufficient sample volume. Batch: GCV/4730

[M5] A matrix spike/matrix spike duplicate was not performed for this batch due to insufficient sample volume. Batch: OEXT/43560

[M5] A matrix spike/matrix spike duplicate was not performed for this batch due to insufficient sample volume.

Batch: MSV/60740

[M5] A matrix spike/matrix spike duplicate was not performed for this batch due to insufficient sample volume.

ANALYTE QUALIFIERS

- H6 Analysis initiated outside of the 15 minute EPA recommended holding time.
- L0 Analyte recovery in the laboratory control sample (LCS) was outside QC limits.
- L2 Analyte recovery in the laboratory control sample (LCS) was below QC limits. Results for this analyte in associated samples may be biased low.
- M1 Matrix spike recovery exceeded QC limits. Batch accepted based on laboratory control sample (LCS) recovery.



QUALITY CONTROL DATA CROSS REFERENCE TABLE

 Project:
 085692 MARCOTTE NO 1

 Pace Project No.:
 60166332

Lab ID	Sample ID	QC Batch Method	QC Batch	Analytical Method	Analytical Batch
60166332001	GW-085692-040214-CM-MW-2	EPA 3510C	OEXT/43505	EPA 8015B	GCSV/16583
60166332001	GW-085692-040214-CM-MW-2	EPA 5030B/8015B	GCV/4730		
60166332001	GW-085692-040214-CM-MW-2	EPA 3010	MPRP/26840	EPA 6010	ICP/20385
60166332001	GW-085692-040214-CM-MW-2	EPA 3010	MPRP/26827	EPA 6020	ICPM/2896
60166332001	GW-085692-040214-CM-MW-2	EPA 7470	MERP/8283	EPA 7470	MERC/8235
60166332001	GW-085692-040214-CM-MW-2	EPA 3510	OEXT/43560	EPA 8270	MSSV/13901
60166332001	GW-085692-040214-CM-MW-2	EPA 5030B/8260	MSV/60740		
60166332001	GW-085692-040214-CM-MW-2	EPA 120.1	WET/47146		
60166332001	GW-085692-040214-CM-MW-2	SM 2320B	WET/47254		
60166332001	GW-085692-040214-CM-MW-2	SM 2540C	WET/47185		
60166332001	GW-085692-040214-CM-MW-2	SM 4500-H+B	WET/47180		
60166332001	GW-085692-040214-CM-MW-2	EPA 300.0	WETA/29042		
60166332001	GW-085692-040214-CM-MW-2	EPA 353.2	WETA/28919		
60166332001	GW-085692-040214-CM-MW-2	EPA 365.1	WETA/28883		



Sample Condition Upon Receipt ESI Tech Spec Client

WO#:60166332

Client Name: COP CRANM Optional Courier: Fed Ex 🗹 UPS 🗆 USPS 🗆 Client 🗆 Commercial 🗅 Pace 🗆 Other Proj Due Date: Tracking #: 5689 1282 1903 Yes 🗹 No 🗆 Pace Shipping Label Used? Proj Name: Custody Seal on Cooler/Box Present: Yes 🗹 No 🗆 Seals intact: Yes 🗸 No 🗆 Other VZPLC Packing Material: Bubble Wrap Bubble Bags Foam 🗆 None 🗆 (-239 / T-194 Thermometer Used: Type of Ice: (Wet) Blue None
Samples received on ice, cooling process has begun. (circle one) **Cooler Temperature:** 4.0 Date and initials of person examining contents: <u>4 /4 / 14 BF</u> Temperature should be above freezing to 6°C Chain of Custody present: Yes No N/A VYes No NA Chain of Custody filled out: Chain of Custody relinquished: Yes No □n/A Yes No N/A Sampler name & signature on COC: Yes No □n/A Samples arrived within holding time: VYes No N/A Short Hold Time analyses (<72hr): 6. Noz Nos, Ortho-p TYes No DNA Rush Turn Around Time requested: Ves ONO ON/A Sufficient volume: IZIYes □No □N/A Correct containers used Pires DNo DN/A Pace containers used: Containers intact: 10 Yes No 12 N/A Unpreserved 5035A soils frozen w/in 48hrs? 11. Yes No NA Filtered volume received for dissolved tests? 12 Ves ONO ON/A Sample labels match COC: Includes date/time/ID/analyses Matrix: wor 13. All containers needing preservation have been checked. Yes No NA All containers needing preservation are found to be in compliance with EPA recommendation. 14 Exceptions: NOA) coliform, TOC, O&G, WI-DRO (water). Initial when Lot # of added Yes DNo Phenolics completed preservative Trip Blank present: Yes No NA Pace Trip Blank lot # (if purchased): 031714-3860 5 Headspace in VOA vials (>6mm): 16 □Yes □No Project sampled in USDA Regulated Area: 17. List State: Client Notification/ Resolution: Copy COC to Client? IN Y 1 Field Data Required? Y / N Temp Log: Record start and finish times Person Contacted: Date/Time: when unpacking cooler, if >20 min, Comments/ Resolution recheck sample temps. Start: 0900 Start: End: 0905 End: Project Manager Review: Date [emp Temp:

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

Pace Analytical

שלשלה שבשביט ופרשה שקשיע שבשרי Pace Project No./ Lab I.D. DRINKING WATER (N/A) 120166332 Samples Intact SAMPLE CONDITIONS > OTHER ę. Cooler (Y/N) Custody Seale 3044 (18) (N/Y) eol **GROUND WATER** 7 Received on Page: Residual Chlorine (Y/V) 40 SE40C LDS O° ni qmeT REGULATORY AGENCY N RCRA 120.1 Specific Conductan Requested Analysis Filtered (Y/N) Hq 'vinits/IA 1.015 TIME 0835 L 365.1 Ortho, P 1 Site Location STATE: 1 anoinA 0.005 NPDES 4/4/14 DATE 553.2 Nitrate/Nitrite UST V470 Mercury L ** sistem . Retais Aosoa SO15B TPH DRO SO15B TPH GRO DATE Signed ACCEPTED BY / AFFILIATION /PRSI KS 8270 SVOCs 1 8260 VOCs IN /A treat sizylenA ć Other Nethanol Alice Flanagan 20 Preservatives 602S2BN ePayables HOBN Manager. Pace Profile #: 5514 IOH Invoice Information: [©]ONH Company Name Pace Quote Reference: Pace Project ^{*}OS^zH Section C Attention: Address: Unpreserved BMIT a # OF CONTAINERS SAMPLE TEMP AT COLLECTION SAMPLER NAME AND SIGNATURE PRINT Name of SAMPLER SIGNATURE of SAMPLER DATE 345 ñ TIME COMPOSITE 1 214 COLLECTED DATE RELINQUISHED BY I AFFILIATION Jeff Walker, Angela Bown TIME COMPOSITE Report To: Christine Mathews START Marcotte No 1 DATE Required Project Information: *085692 (G=GRAB C=COMP) Purchase Order No.: SAMPLE TYPE roject Number. Project Name: AMATRIX CODE Section B Copy To: 2 -011-110-020 Dissolved metals: As, Ba, Cd, Cr, Co, Cu, Pb, Mn, Mo, Ni, 6121 Indian School Rd NE, Ste 200 Fax: (505)884-4932 0914 cmathews@craworld.com Albequerque, NM 87110 ADDITIONAL COMMENTS (A-Z, 0-9 / ,-) Sample IDs MUST BE UNIQUE standard 20 SAMPLE ID COP CRA NM Section D Required Client Information Required Client Information: (505)884-0672 Requested Due Date/TAT: 000 Ag, Zh, HARDNESS BE Section A 0 Company 0 Address; Email To: hone: Page 48 of 48 # WBLI -2 3 4 5 9 ~ 0 9 11 42 00

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F-AtL-Q-020rev 08, 12-Oct-2007

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(YYDDINI)

1 for any invoices not paid within 30 days.

Important Note" By signing this form you are accepting Pace's NET 30 day payment terms and agreeing to late charges of 1.5% per mon

Appendix B

NMOSE Domestic Water Well List





New Mexico Office of the State Engineer Active & Inactive Points of Diversion

(R=POD has been replaced

	(acre ft p ເຮັນໄວ	er annum)		(R=POD has been replaced and no longer serves this file, C=the file is closed)	, (quarters are 1=NW 2=NE 3= (quarters are smallest to large q q q	
WR File Nbr	basin Use Diver		County POD Number SJ SJ 00585	Code Grant	Source 6416 4 Sec Tws R Shallow 08 31N 1	
SJ 00585	DOM	3 HAROLD AND VEVERALY PEPIN	31 21 00000			
SJ 01078	DOM	3 ROSS C. SILVA	SJ <u>SJ 01078</u>		3 08 31N 1	DW 240709 4088789* 🍑
SJ 01081	DOM	3 HANK STIEHL	SJ SJ 01081		3 3 08 31N 1	DW 240508 4088588* 🎡
SJ 01085	DOM	3 DALE BOYD	SJ <u>SJ 01085</u>		4 1 08 31N 1	DW 240946 4089384* 🍈
SJ 01091	DOM	3 JOHN CHRISTENSEN	SJ SJ 01091		3 08 31N 1	DW 240709 4088789* 🎡
SJ 01288	DOM	3 WILLIAM B. HUDSON	SJ <u>SJ 01288</u>		1 1 08 31N 1	DW 240551 4089802* 🛶
SJ 01584	DOM	0 F. R. DICKENS	SJ <u>SJ 01584</u>		3 1 08 31N 1	OW 240537 4089397* 🍥
SJ 01940	DOM	3 HAROLD B. JARRED	SJ SJ 01940		2 1 08 31N 1	OW 240959 4089789* 🧼
SJ 02153	DOM	3 CHARLES GAINES	SJ SJ 02153		2 1 08 31N 1	DW 240959 4089789* 🎡
SJ 02304	DOM	3 EDWARD K. ZINK	SJ SJ 02304		Shallow 2 1 08 31N 1	DW 240959 4089789* 🎡
SJ 02432	DOM	3 BERT HARRIS	SJ <u>SJ 02432</u>		4 1 08 31N 1	DW 240946 4089384* 🍥
SJ 02566	DOM	3 CLARENCE D. KEENOM	SJ SJ 02566		1 2 3 08 31N 1	OW 240832 4089079* 🎡
SJ 03057	DOM	3 ROD KESSELHUTH	SJ SJ 03057		Shallow 4 3 1 08 31N 1	DW 240636 4089296* 🎡
<u>SJ 03714</u>	DOM	3 J.A. WILSON	SJ SJ 03714 POD1		Shallow 1 1 3 08 31N 1	DW 240421 4089091* 🍥
SJ 03923	DOM	1 KAREN BRASFIELD	SJ SJ 03923 POD1		Shallow 4 1 2 08 31N 1	0W 241455 4089675 🍚
SJ 03930	DOM	1 VIRGINIA A. LENBERG	SJ SJ 03930 POD1		Shallow 3 1 3 08 31N 1	0W 240624 4088974 췢
SJ 03988	DOM	1 JOSEPH WILSON	SJ SJ 03988 POD1		2 1 3 08 31N 1	DW 240561 4089123 🍈
SJ 03989	DOM	1 JOSEPH WILSON	SJ SJ 03989 POD1		2 1 3 08 31N 1	DW 240554 4089078 🧼
*UTM location w	as derived from PLS	S - see Help				
						IVE POINTS OF DIVERSION

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ACTIVE & INACTIVE POINTS OF DIVERSION

				(R=POD has been replaced and no longer serves this file,	(quarters are 1-	NIM 2-NE 3-SM	-95)	
(acre ft per annum)			C=the file is closed)	, (quarters are 1=NW 2=NE 3=SW 4=SE) (quarters are smallest to largest) (NAD83 UTM in meters)				
Sub			×	,		,		
WR File Nbr	basin Use Diver	sion Owner	County POD Number	Code Grant	Source 6416 4	Sec Tws Rng	Х	Y
SJ 03990	DOM	1 JOSEPH WILSON	SJ SJ 03990 POD1		213	08 31N 10W	240556	4089022 🧼
SJ 03991	DOM	1 JOE WILSON	SJ SJ 03991 POD1		313	08 31N 10W	240525	4088961 🌑
SJ 03992	DOM ·	1 J A WILSON	SJ <u>SJ 03992 POD1</u>		313	08 31N 10W	240485	4088974 🎡
SJ 03993	DOM	1 J A WILSON	SJ 03993 POD1		113	08 31N 10W	240484	4089021 🍥
SJ 04005	DOL	3 CLARENCE D. KEENOM	SJ 04005 POD1		123	08 31N 10W	240879	4089011 🍈
SJ 04044	DOL	3 ADRIAN GRIEGO	SJ SJ 04044 POD1		421	08 31N 10W	241163	4089709 🍥
Record Count:	24							

The data is furnished by the NMOSE/ISC and is accepted by the recipient with the expressed understanding that the OSE/ISC make no warranties, expressed or implied, concerning the accuracy, completeness, reliability, or suitability for any particular purpose of the data.
7/21/14 10:11 AM Page 2 of 2 ACTIVE & INACTIVE POINTS OF DIVERSION

PLSS Search: Section(s): 8

Sorted by: File Number

Township: 31N

Range: 10W

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