ADEQ
ARKANSAS
Department of Environmental Quality

WATER DIVISION COMPLAINT REPORT

AFIN: **23-00000**	PERMIT #: **N/A**
COUNTY: **23 Faulkner**	PDS #: **015781**
GPS LOCATION: LAT: **34.90373** LONG: **-92.449438** ☒ Discharge Site / ☐ General Area / ☐ Entrance	

COMPLAINANT NAME: **Congressman Tim Griffin**	COMPLAINT AGAINST: **Unknown**
RESPONSE REQUESTED BY COMPLAINANT: **Yes**	SITE ADDRESS:
MAILING ADDRESS: **1105 Deer St., Suite 12**	MAILING ADDRESS:
CITY, STATE, ZIP: **Conway AR 72032**	CITY, STATE, ZIP:
PHONE & EXT: **501.358.3481** FAX:	PHONE & EXT: FAX:
EMAIL:	EMAIL:
HOW WAS COMPLAINT RECEIVED BY ADEQ: **EML**	ADEQ = ADEQ Personnel | EML = E-Mail | FAX = Fax | LTR = Letter PERS = Personal Contact | PHON = Telephone | WEB = ADEQ Website

PERSON RECEIVING REPORT: **Dean VanDerhoff**	DATE: **6/24/13**	SUPERVISOR REFERRAL: **Kerri McCabe**	DATE: **6/24/13**
MEDIA SUPERVISOR REFERRAL:	DATE:	RECEIVING INSPECTOR: (ID & Name) **26036 Risa Parker**	DATE: **6/24/13**

DETAILED DESCRIPTION OF COMPLAINT

Lance Hines, Director of Sales & Business Development,lance@priority1inc.com, 501-374-5960(fax);501-626-1793(cell) notified Congressman Griffin on 6/24/13 that one of Hines' employees observed oil running into the Arkansas River. Subsequently, the EPA (Nicolas Bresica) was notified by Congressman Griffin's office. EPA notified Dean VanDerhoff, ADEQ, Hazardous Waste Division.

DETAILED LOCATION

Confluence of Palarm Creek and the Arkansas River off Hwy 365 (see attached map).

PREVIOUS COMPLAINTS: **No** DATE(S):

FOLLOW-UP ON COMPLAINT

NUMBER OF SITE VISITS: **1** DATES: FIRST **6/24/13** SECOND THIRD

PHOTOS TAKEN: **Yes**	DISCHARGE TO WATER OF THE STATE: **No**	SAMPLES COLLECTED: **Yes**

NAME OF WATERBODY: **Palarm Creek/Arkansas River.**

FAYETTEVILLE SHALE RELATED: **No**	FAYETTEVILLE SHALE VIOLATIONS: **No**

INVESTIGATION & ACTION TAKEN

6/24/13-(Confluence of Palarm Creek/Arkansas River)-Dean VanDerhoff, Erica McAdoo, and I responded to a complaint alleging oil was running into the Arkansas River from Palarm Creek at the confluence. Upon arrival, the following conditions were observed: a distinct mixing zone at the confluence of Palarm Creek with the Arkansas River was observed. Palarm Creek was dark (tea-colored); this is attributed to the decomposition of organic material; specifically, plant matter and the presence of tannins and lignins. Samples were obtained and exceedences of APC&EC Regulation 2: Water Quality Standards for Surface Waters of the State of Arkansas were not detected in the samples obtained at the time of investigation. See NPDES Report Page 6 of 37 for the approximate sampling location and NPDES Report Pages 7-36 of 37 for sample results.

Additional locations investigated were:

(Arkansas River-upstream of confluence with Palarm Creek): water samples were not collected; however, the following readings were obtained using field meters: Dissolved Oxygen (DO): 6.61 mg/L; pH 7.62.

(Boat Launch Area-Palarm Creek-Hwy 365): the following conditions were observed at the boat launch area off Hwy 365: Palarm Creek was dark (tea-colored); the coloration was attributed to the decomposition of organic material; a film was observed around the boat launch structure, this was attributed to bacterial activity in relatively stagnant water causing a biofilm. Water samples were not collected; however, the following readings were obtained using field meters: DO: 5.86 mg/L @ 34.2°C; pH 7.21.

(Palarm Creek located downstream of Lake Conway dam and boat launch area): the following conditions were observed downstream of the dam: Palarm Creek was dark (tea-colored); the coloration was attributed to the decomposition of organic material. Water samples were not collected; however, the following readings were obtained using field meters: DO: 5.04 mg/L @ 31.4°C; pH 7.14.

REFERRED: **No**	TO WHOM: **N/A**	DATE REFERRED:

RESPONSE TRACKING		
RESPONSE REQUESTED: **Yes**	RESPONSE PROVIDED: **Yes**	RESPONSE PROVIDED DATE: **See comment below.**
BY WHOM?(ID and Name):	ASSIGNED TO? (ID and Name):	
RESPONSE COMMENT: **Per verbal notification to me by Dean VanDerhoff on 6/24/13, the response to Congressman Griffin is to be coordinated through the ADEQ Hazardous Waste Division.**		
INSPECTOR: **Risa Parker**		DATE: **6/27/13**
SUPERVISOR: **Jason Bolenbaugh**		DATE: **6/27/2013**

Revised: 5/13/2013

Water Division Photographic Evidence Sheet					
Location:	**Unknown Palarm Creek upstream of its confluence with the Arkansas River.**				
Photographer:	**Risa Parker, ADEQ Water Division**	Date:	**6/24/13**	Time:	**1:30 pm**
Witness:	**Erica McAdoo (Water Division) and Dean VanDerhoff (Hazardous Waste Division).**			Photo #:	**1**
Description:	**The color and clarity of Palarm Creek prior to its confluence with the Arkansas River.**				

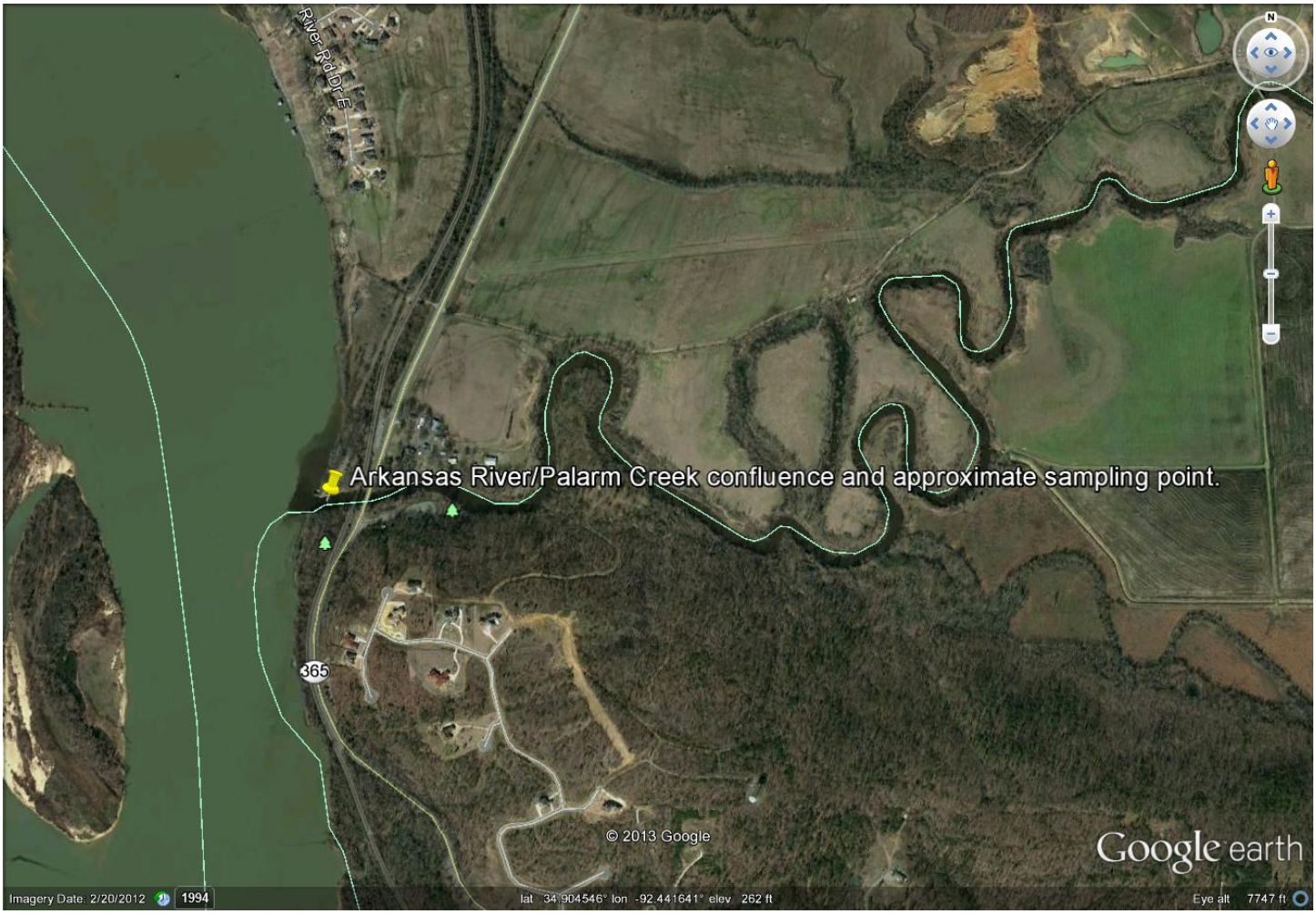
Photographer:	**Risa Parker, ADEQ Water Division**	Date:	**6/24/13**	Time:	**1:33 pm**
Witness:	**Erica McAdoo (Water Division) and Dean VanDerhoff (Hazardous Waste Division).**			Photo #:	**2**
Description:	**Mixing zone of Palarm Creek and the Arkansas River-note difference in coloration of both water bodies. The color of Palarm Creek is attributed to the decomposition of plant matter and the presence of tannins and lignins.**				

Water Division Photographic Evidence Sheet					
Location:	**Unknown**				
Photographer:	**Risa Parker, ADEQ Water Division**	Date:	**6/24/13**	Time:	**1:56 pm**
Witness:	**Erica McAdoo (Water Division) and Dean VanDerhoff (Hazardous Waste Division).**			Photo #:	**3**
Description:	**Film in Palarm Creek at this location attributed to bacteria.**				

Photographer:	**Risa Parker, ADEQ Water Division**	Date:	**6/24/13**	Time:	**1:58 pm**
Witness:	**Erica McAdoo (Water Division) and Dean VanDerhoff (Hazardous Waste Division).**			Photo #:	**4**
Description:	**Film in Palarm Creek at this location attributed to bacteria.**				

Water Division Photographic Evidence Sheet					
Location:	**Palarm Creek located downstream of Lake Conway dam and boat launch area.**				
Photographer:	**Risa Parker, ADEQ Water Division**	Date:	**6/24/13**	Time:	**2:35 pm**
Witness:	**Erica McAdoo (Water Division) and Dean VanDerhoff (Hazardous Waste Division).**			Photo #:	**5**
Description:	**The color of Palarm Creek is attributed to the decomposition of plant matter and the presence of tannins and lignins.**				

Photographer:	**Risa Parker, ADEQ Water Division**	Date:	**6/24/13**	Time:	**2:34 pm**
Witness:	**Erica McAdoo (Water Division) and Dean VanDerhoff (Hazardous Waste Division).**			Photo #:	**6**
Description:	**The color of Palarm Creek is attributed to the decomposition of plant matter and the presence of tannins and lignins.**				


River Rd Dr E
Arkansas River/Palarm Creek confluence and approximate sampling point.
365
© 2013 Google
Google earth
Imagery Date: 2/20/2012
1994
lat 34.904546° lon -92.441641° elev 262 ft
Eye alt 7747 ft

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

Client Report For:	Palarm Creek Complaint 2013 2215-2216
Attention:	
Client Address:	
	,
Report Date:	June 26, 2013
LAB ID:	AR13JUN24-02
Comment:	

Approved By:_____ Date:June 26, 2013

Arkansas Department of Environmental Quality
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North Liitle Rock, AR 72118

Laboratory Contact: Jeff Ruehr
Ruehr@adeq.state.ar.us
501-682-0955

Client:	Special Samples	**Client Sample ID:**	Palarm Creek Complaint
Lab ID:	2013-2215	**Collection Date:**	6/24/2013 1:41:00 PM
		Matrix:	Water

Analyses

Oil and Grease ***EPA1664*** ***Batch: 13062503*** ***Run: 1***

	Result	Reporting Limit	MDL	Qual	Unit
Oil and Grease	<2.5	2.5	2.5		mg/L
Dilution Factor	1				
Analyzed By	Robert Graddy				
Analysis Date/Time	06/25/2013 1000				

Semi-Volatiles by GC/MS ***EPA 3510C/EPA 8270D*** ***Batch: 13062504*** ***Run: 1***

	Result	Reporting Limit	MDL	Qual	Unit
2-Fluorophenol (% Recovery)	37.3	40-110			%
Nitrobenzene-d5 (% Recovery)	65.5	50-110			%
2-Fluorobiphenyl (% Recovery)	66.0	50-110			%
2,4,6-Tribromophenol (% Recovery)	65.6	40-110			%
Terphenyl-d14 (% Recovery)	64.0	50-110			%
Methyl Methanesulfonate	<0.2	0.2	100		ug/L
Ethyl methanesulfonate	<0.2	0.2	100		ug/L
Phenol	<0.2	0.2	100		ug/L
Aniline	<0.2	0.2	100		ug/L
Bis(2-chloroethyl)ether	<0.2	0.2	100		ug/L
2-Chlorophenol	<0.2	0.2	100		ug/L
1,3-Dichlorobenzene	<0.12	0.12	60		ug/L
1,4-Dichlorobenzene	<0.12	0.12	60		ug/L
Benzyl alcohol	0.207	0.16	80		ug/L
1,2-Dichlorobenzene	<0.12	0.12	60		ug/L
2-Methylphenol	<0.1	0.1	50		ug/L
Acetophenone	<0.1	0.1	50		ug/L
4-Methylphenol	<0.1	0.1	50		ug/L
N-Nitrosodi-n-propylamine	<0.2	0.2	100		ug/L
Hexachloroethane	<0.2	0.2	100		ug/L
Nitrobenzene	<0.2	0.2	100		ug/L
N-Nitrosopiperidine	<0.2	0.2	100		ug/L

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Isophorone	<0.1	0.1	50	ug/L
2-Nitrophenol	<0.3	0.3	150	ug/L
2,4-Dimethylphenol	<0.1	0.1	50	ug/L
Bis(2-chloroethoxy)methane	<0.2	0.2	100	ug/L
2,4-Dichlorophenol	<0.2	0.2	100	ug/L
1,2,4-Trichlorobenzene	<0.12	0.12	60	ug/L
Naphthalene	<0.08	0.08	40	ug/L
4-Chloroaniline	<0.1	0.1	50	ug/L
2,6-Dichlorophenol	<0.2	0.2	100	ug/L
Hexachlorobutadiene	<0.2	0.2	100	ug/L
N-Nitrosodibutylamine	<0.2	0.2	100	ug/L
4-Chloro-3-methylphenol	<0.16	0.16	80	ug/L
2-Methylnaphthalene	<0.1	0.1	50	ug/L
1,2,4,5-Tetrachlorobenzene	<0.1	0.1	50	ug/L
Hexachlorocyclopentadiene	<0.16	0.16	80	ug/L
2,4,6-Trichlorophenol	<0.2	0.2	100	ug/L
2,4,5-Trichlorophenol	<0.2	0.2	100	ug/L
2-Chloronaphthalene	<0.1	0.1	50	ug/L
1-Chloronaphthalene	<0.1	0.1	50	ug/L
2-Nitroaniline	<0.2	0.2	100	ug/L
Dimethyl phthalate	<0.2	0.2	100	ug/L
2,6-Dinitrotoluene	<0.2	0.2	100	ug/L
Acenaphthylene	<0.08	0.08	40	ug/L
3-Nitroaniline	<0.2	0.2	100	ug/L
Acenaphthene	<0.1	0.1	50	ug/L
2,4-Dinitrophenol	<4	4	2000	ug/L
Pentachlorobenzene	<0.12	0.12	60	ug/L
4-Nitrophenol	<2	2	1000	ug/L
Dibenzofuran	<0.1	0.1	50	ug/L
2,4-Dinitrotoluene	<0.2	0.2	100	ug/L
2,3,4,6-Tetrachlorophenol	<0.6	0.6	300	ug/L
Diethyl phthalate	<0.2	0.2	100	ug/L
Fluorene	<0.1	0.1	50	ug/L
4-Chlorophenyl phenyl ether	<0.1	0.1	50	ug/L
4-Nitroaniline	<0.2	0.2	100	ug/L
4,6-Dinitro-2-methylphenol	<6	6	3000	ug/L
Diphenylamine	<0.1	0.1	50	ug/L
Azobenzene	<0.08	0.08	40	ug/L
4-Bromophenyl phenyl ether	<0.2	0.2	100	ug/L
Hexachlorobenzene	<0.16	0.16	80	ug/L

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Pentachlorophenol	<1	1	500	ug/L
Pentachloronitrobenzene	<0.2	0.2	100	ug/L
Pronamide	<0.2	0.2	100	ug/L
Phenanthrene	<0.08	0.08	40	ug/L
Anthracene	<0.08	0.08	40	ug/L
Carbazole	<0.1	0.1	50	ug/L
Di-n-butyl phthalate	<0.2	0.2	100	ug/L
Fluoranthene	<0.08	0.08	40	ug/L
Pyrene	<0.08	0.08	40	ug/L
Dimethylaminoazobenzene	<0.2	0.2	100	ug/L
Butyl benzyl phthalate	<0.3	0.3	150	ug/L
Benzo (a) anthracene	<0.1	0.1	50	ug/L
Chrysene	<0.1	0.1	50	ug/L
Bis(2-ethylhexyl)phthalate	<0.3	0.3	150	ug/L
Di-n-octyl phthalate	<0.3	0.3	150	ug/L
Benzo (b) fluoranthene	<0.16	0.16	80	ug/L
7,12-Dimethylbenz (a) anthracene	<0.2	0.2	100	ug/L
Benzo (k) fluoranthene	<0.16	0.16	80	ug/L
Benzo (a) pyrene	<0.16	0.16	80	ug/L
3-Methylcholanthrene	<0.2	0.2	100	ug/L
Indeno (1,2,3-cd) pyrene	<0.2	0.2	100	ug/L
Dibenzo (a,h) anthracene	<0.16	0.16	80	ug/L
Benzo (g,h,i) perylene	<0.16	0.16	80	ug/L
Initial Volume	500			mL
Final Volume	1			mL
Dilution Factor	1			
Analyzed By	Ed Harris			
Analysis Date/Time	6/25/2013 11:52 AM			
Prep By	Ed Harris			
Prep Date/Time	6/25/2013 08:00			

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Client: Special Samples
Lab ID: 2013-2215
Client Sample ID: Palarm Creek Complaint
Collection Date: 6/24/2013 1:41:00 PM
Matrix: Water

Analyses

Volatile Organics by GCMS *EPA 8260C* *Batch: 13062602* *Run: 1*

	Result	Reporting Limit	MDL	Qual	Unit
Dibromofluoromethane (% Recovery)	108	70-130			%
1,2-Dichloroethane-d4 (% Recovery)	107	70-130			%
Toluene-d8 (% Recovery)	93.3	70-130			%
4-Bromofluorobenzene (% Recovery)	104	70-130			%
Dichlorodifluoromethane	<1.12	1.12	1.12		ug/L
Chloromethane	<0.58	0.58	0.58		ug/L
Vinyl chloride	<0.82	0.82	0.82		ug/L
Bromomethane	<3.9	3.9	3.90		ug/L
Chloroethane	<2.68	2.68	2.68		ug/L
Trichlorofluoromethane	<0.51	0.51	0.51		ug/L
1,1-Dichloroethene	<0.43	0.43	0.43		ug/L
Acetone	<10.5	10.5	10.5		ug/L
Methylene chloride	<2.5	2.5	2.5		ug/L
Methyl tert-butyl ether	<0.83	0.83	0.83		ug/L
trans-1,2-Dichloroethene	<0.59	0.59	0.59		ug/L
1,1-Dichloroethane	<0.42	0.42	0.42		ug/L
Methyl ethyl ketone	<12.8	12.8	12.8		ug/L
cis-1,2-Dichloroethene	<1.15	1.15	1.15		ug/L
2,2-Dichloropropane	<0.81	0.81	0.81		ug/L
Bromochloromethane	<0.66	0.66	0.66		ug/L
Chloroform	<0.27	0.27	0.27		ug/L
1,1,1-Trichloroethane	<0.46	0.46	0.46		ug/L
1,1-Dichloropropene	<0.59	0.59	0.59		ug/L
Carbon tetrachloride	<0.6	0.6	0.6		ug/L
Benzene	<0.66	0.66	0.66		ug/L
1,2-Dichloroethane	<1.15	1.15	1.15		ug/L
Trichloroethene	<0.6	0.6	0.60		ug/L
1,2-Dichloropropane	<0.98	0.98	0.98		ug/L
Dibromomethane	<1.78	1.78	1.78		ug/L
Bromodichloromethane	<0.65	0.65	0.65		ug/L
cis-1,3-Dichloropropene	<0.86	0.86	0.86		ug/L
Methyl isobutyl ketone	<8.1	8.1	8.10		ug/L

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Toluene	0.595	0.57	0.57	ug/L
trans-1,3-Dichloropropene	<0.84	0.84	0.84	ug/L
1,1,2-Trichloroethane	<0.78	0.78	0.78	ug/L
2-Hexanone	<9.5	9.5	9.5	ug/L
Tetrachloroethene	<0.96	0.96	0.96	ug/L
1,3-Dichloropropane	<0.94	0.94	0.94	ug/L
Dibromochloromethane	<1.25	1.25	1.25	ug/L
1,2-Dibromoethane (EDB)	<0.68	0.68	0.68	ug/L
Chlorobenzene	<0.62	0.62	0.62	ug/L
Ethylbenzene	<0.51	0.51	0.51	ug/L
1,1,1,2-Tetrachloroethane	<0.57	0.57	0.57	ug/L
m,p-Xylene	<1.2	1.2	1.2	ug/L
o-Xylene	0.527	0.5	0.5	ug/L
Styrene	1.03	0.53	0.53	ug/L
Bromoform	<1.56	1.56	1.56	ug/L
Isopropylbenzene	<0.59	0.59	0.59	ug/L
1,1,2,2-Tetrachloroethane	<0.39	0.39	0.39	ug/L
1,2,3-Trichloropropane	<1.83	1.83	1.83	ug/L
n-Propylbenzene	<0.49	0.49	0.49	ug/L
Bromobenzene	<0.5	0.5	0.5	ug/L
1,3,5-Trimethylbenzene	<0.3	0.3	0.30	ug/L
2-Chlorotoluene	<0.66	0.66	0.66	ug/L
4-Chlorotoluene	<0.8	0.8	0.80	ug/L
tert-Butylbenzene	<0.85	0.85	0.85	ug/L
1,2,4-Trimethylbenzene	<0.46	0.46	0.46	ug/L
sec-Butylbenzene	<0.63	0.63	0.63	ug/L
p-Isopropyltoluene	<0.59	0.59	0.59	ug/L
1,3-Dichlorobenzene	<0.7	0.7	0.70	ug/L
1,4-Dichlorobenzene	<0.53	0.53	0.53	ug/L
n-Butylbenzene	<0.72	0.72	0.72	ug/L
1,2-Dichlorobenzene	<0.7	0.7	0.70	ug/L
1,2-Dibromo-3-chloropropane	<0.86	0.86	0.86	ug/L
1,2,4-Trichlorobenzene	<1.14	1.14	1.14	ug/L
Naphthalene	<1.53	1.53	1.53	ug/L
1,2,3-Trichlorobenzene	<1.3	1.3	1.3	ug/L
Dilution Factor	1			
Analyzed By	Jeff Ruehr			
Analysis Date/Time	6/25/2013 1:36 PM			

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Laboratory Contact: Jeff Ruehr
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501-682-0955

Client: Special Samples
Lab ID: 2013-2216
Client Sample ID: Volatiles Trip Blank
Collection Date: 6/24/2013 1:41:00 PM
Matrix: Water

Analyses

Volatile Organics by GCMS ***EPA 8260C*** ***Batch: 13062602*** ***Run: 1***

	Result	Reporting Limit	MDL	Qual	Unit
Dibromofluoromethane (% Recovery)	106	70-130			%
1,2-Dichloroethane-d4 (% Recovery)	104	70-130			%
Toluene-d8 (% Recovery)	95.9	70-130			%
4-Bromofluorobenzene (% Recovery)	101	70-130			%
Dichlorodifluoromethane	<1.12	1.12	1.12		ug/L
Chloromethane	<0.58	0.58	0.58		ug/L
Vinyl chloride	<0.82	0.82	0.82		ug/L
Bromomethane	<3.9	3.9	3.90		ug/L
Chloroethane	<2.68	2.68	2.68		ug/L
Trichlorofluoromethane	<0.51	0.51	0.51		ug/L
1,1-Dichloroethene	<0.43	0.43	0.43		ug/L
Acetone	<10.5	10.5	10.5		ug/L
Methylene chloride	<2.5	2.5	2.5		ug/L
Methyl tert-butyl ether	<0.83	0.83	0.83		ug/L
trans-1,2-Dichloroethene	<0.59	0.59	0.59		ug/L
1,1-Dichloroethane	<0.42	0.42	0.42		ug/L
Methyl ethyl ketone	<12.8	12.8	12.8		ug/L
cis-1,2-Dichloroethene	<1.15	1.15	1.15		ug/L
2,2-Dichloropropane	<0.81	0.81	0.81		ug/L
Bromochloromethane	<0.66	0.66	0.66		ug/L
Chloroform	<0.27	0.27	0.27		ug/L
1,1,1-Trichloroethane	<0.46	0.46	0.46		ug/L
1,1-Dichloropropene	<0.59	0.59	0.59		ug/L
Carbon tetrachloride	<0.6	0.6	0.6		ug/L
Benzene	<0.66	0.66	0.66		ug/L
1,2-Dichloroethane	<1.15	1.15	1.15		ug/L
Trichloroethene	<0.6	0.6	0.60		ug/L
1,2-Dichloropropane	<0.98	0.98	0.98		ug/L
Dibromomethane	<1.78	1.78	1.78		ug/L
Bromodichloromethane	<0.65	0.65	0.65		ug/L
cis-1,3-Dichloropropene	<0.86	0.86	0.86		ug/L
Methyl isobutyl ketone	<8.1	8.1	8.10		ug/L

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Toluene	<0.57	0.57	0.57	ug/L
trans-1,3-Dichloropropene	<0.84	0.84	0.84	ug/L
1,1,2-Trichloroethane	<0.78	0.78	0.78	ug/L
2-Hexanone	<9.5	9.5	9.5	ug/L
Tetrachloroethene	<0.96	0.96	0.96	ug/L
1,3-Dichloropropane	<0.94	0.94	0.94	ug/L
Dibromochloromethane	<1.25	1.25	1.25	ug/L
1,2-Dibromoethane (EDB)	<0.68	0.68	0.68	ug/L
Chlorobenzene	<0.62	0.62	0.62	ug/L
Ethylbenzene	<0.51	0.51	0.51	ug/L
1,1,1,2-Tetrachloroethane	<0.57	0.57	0.57	ug/L
m,p-Xylene	<1.2	1.2	1.2	ug/L
o-Xylene	<0.5	0.5	0.5	ug/L
Styrene	1.04	0.53	0.53	ug/L
Bromoform	<1.56	1.56	1.56	ug/L
Isopropylbenzene	<0.59	0.59	0.59	ug/L
1,1,2,2-Tetrachloroethane	<0.39	0.39	0.39	ug/L
1,2,3-Trichloropropane	<1.83	1.83	1.83	ug/L
n-Propylbenzene	<0.49	0.49	0.49	ug/L
Bromobenzene	<0.5	0.5	0.5	ug/L
1,3,5-Trimethylbenzene	<0.3	0.3	0.30	ug/L
2-Chlorotoluene	<0.66	0.66	0.66	ug/L
4-Chlorotoluene	<0.8	0.8	0.80	ug/L
tert-Butylbenzene	<0.85	0.85	0.85	ug/L
1,2,4-Trimethylbenzene	<0.46	0.46	0.46	ug/L
sec-Butylbenzene	<0.63	0.63	0.63	ug/L
p-Isopropyltoluene	<0.59	0.59	0.59	ug/L
1,3-Dichlorobenzene	<0.7	0.7	0.70	ug/L
1,4-Dichlorobenzene	<0.53	0.53	0.53	ug/L
n-Butylbenzene	<0.72	0.72	0.72	ug/L
1,2-Dichlorobenzene	<0.7	0.7	0.70	ug/L
1,2-Dibromo-3-chloropropane	<0.86	0.86	0.86	ug/L
1,2,4-Trichlorobenzene	<1.14	1.14	1.14	ug/L
Naphthalene	<1.53	1.53	1.53	ug/L
1,2,3-Trichlorobenzene	<1.3	1.3	1.3	ug/L
Dilution Factor	1			
Analyzed By	Jeff Ruehr			
Analysis Date/Time	6/25/2013 1:11 PM			

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Client:	Special Samples	**Client Sample ID:**	Palarm Creek Complaint
Lab ID:	2013-2215	**Collection Date:**	6/24/2013 1:41:00 PM
		Matrix:	Water

Analyses

Field Data ***Batch: 13062604*** ***Run: 1***

	Result	Reporting Limit	MDL	Qual	Unit
Dissolved Oxygen	4.43				mg/L
pH	6.79				SU
Temperature	30.7				C
Analyzed By	Erica McAdoo				
Analysis Date/Time	6/24/2013 13:47				

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Analytical Quality Control Results Report

Batch: 13062602					**VOA - water**
Palarm Creek Complaint					***LIMS ID: 2013-2215***
Volatiles - water DUP					***Run: 1***
Parameter	***Result***	***DL***	***RL***	***Accuracy Control***	***Precision Control***
Dibromofluoromethane (% Recovery)	103 %			70 - 130	
1,2-Dichloroethane-d4 (% Recovery)	102 %			70 - 130	
Toluene-d8 (% Recovery)	92.2 %			70 - 130	
4-Bromofluorobenzene (% Recovery)	99.4 %			70 - 130	
Dichlorodifluoromethane	<1.12 ug/L	1.12	1.12		
Dichlorodifluoromethane (RPD)	0 %				0 - 20
Chloromethane (RPD)	0 %				0 - 20
Chloromethane	<0.58 ug/L	0.58	0.58		
Vinyl chloride	<0.82 ug/L	0.82	0.82		
Vinyl chloride (RPD)	0 %				0 - 20
Bromomethane (RPD)	0 %				0 - 20
Bromomethane	<3.9 ug/L	3.9	3.9		
Chloroethane	<2.68 ug/L	2.68	2.68		
Chloroethane (RPD)	0 %				0 - 20
Trichlorofluoromethane (RPD)	0 %				0 - 20
Trichlorofluoromethane	<0.51 ug/L	0.51	0.51		
1,1-Dichloroethene (RPD)	0 %				0 - 20
1,1-Dichloroethene	<0.43 ug/L	0.43	0.43		
Acetone	<10.5 ug/L	10.5	10.5		
Acetone (RPD)	38.6 %				0 - 20
Methylene chloride (RPD)	7.1 %				0 - 20
Methylene chloride	<2.5 ug/L	2.5	2.5		
Methyl tert-butyl ether	<0.83 ug/L	0.83	0.83		
Methyl tert-butyl ether (RPD)	0 %				0 - 20
trans-1,2-Dichloroethene (RPD)	0 %				0 - 20
trans-1,2-Dichloroethene	<0.59 ug/L	0.59	0.59		
1,1-Dichloroethane	<0.42 ug/L	0.42	0.42		
1,1-Dichloroethane (RPD)	0 %				0 - 20
Methyl ethyl ketone (RPD)	0 %				0 - 20
Methyl ethyl ketone	<12.8 ug/L	12.8	12.8		
cis-1,2-Dichloroethene	<1.15 ug/L	1.15	1.15		

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cis-1,2-Dichloroethene (RPD)	0 %				0 - 20
2,2-Dichloropropane (RPD)	0 %				0 - 20
2,2-Dichloropropane	<0.81 ug/L	0.81	0.81		
Bromochloromethane	<0.66 ug/L	0.66	0.66		
Bromochloromethane (RPD)	0 %				0 - 20
Chloroform (RPD)	0 %				0 - 20
Chloroform	<0.27 ug/L	0.27	0.27		
1,1,1-Trichloroethane	<0.46 ug/L	0.46	0.46		
1,1,1-Trichloroethane (RPD)	0 %				0 - 20
1,1-Dichloropropene (RPD)	0 %				0 - 20
1,1-Dichloropropene	<0.59 ug/L	0.59	0.59		
Carbon tetrachloride	<0.6 ug/L	0.6	0.6		
Carbon tetrachloride (RPD)	0 %				0 - 20
Benzene (RPD)	0 %				0 - 20
Benzene	<0.66 ug/L	0.66	0.66		
1,2-Dichloroethane	<1.15 ug/L	1.15	1.15		
1,2-Dichloroethane (RPD)	0 %				0 - 20
Trichloroethene (RPD)	0 %				0 - 20
Trichloroethene	<0.6 ug/L	0.6	0.6		
1,2-Dichloropropane	<0.98 ug/L	0.98	0.98		
1,2-Dichloropropane (RPD)	0 %				0 - 20
Dibromomethane (RPD)	0 %				0 - 20
Dibromomethane	<1.78 ug/L	1.78	1.78		
Bromodichloromethane	<0.65 ug/L	0.65	0.65		
Bromodichloromethane (RPD)	0 %				0 - 20
cis-1,3-Dichloropropene (RPD)	0 %				0 - 20
cis-1,3-Dichloropropene	<0.86 ug/L	0.86	0.86		
Methyl isobutyl ketone	<8.1 ug/L	8.1	8.1		
Methyl isobutyl ketone (RPD)	0 %				0 - 20
Toluene (RPD)	0.5 %				0 - 20
Toluene	0.592 ug/L	0.57	0.57		
trans-1,3-Dichloropropene	<0.84 ug/L	0.84	0.84		
trans-1,3-Dichloropropene (RPD)	0 %				0 - 20
1,1,2-Trichloroethane (RPD)	0 %				0 - 20
1,1,2-Trichloroethane	<0.78 ug/L	0.78	0.78		
2-Hexanone	<9.5 ug/L	9.5	9.5		
2-Hexanone (RPD)	0 %				0 - 20
Tetrachloroethene (RPD)	0 %				0 - 20
Tetrachloroethene	<0.96 ug/L	0.96	0.96		
1,3-Dichloropropane	<0.94 ug/L	0.94	0.94		

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1,3-Dichloropropane (RPD)	0 %				0 - 20
Dibromochloromethane (RPD)	0 %				0 - 20
Dibromochloromethane	<1.25 ug/L	1.25	1.25		
1,2-Dibromoethane (EDB)	<0.68 ug/L	0.68	0.68		
1,2-Dibromoethane (EDB) (RPD)	0 %				0 - 20
Chlorobenzene (RPD)	0 %				0 - 20
Chlorobenzene	<0.62 ug/L	0.62	0.62		
Ethylbenzene	<0.51 ug/L	0.51	0.51		
Ethylbenzene (RPD)	0 %				0 - 20
1,1,1,2-Tetrachloroethane (RPD)	0 %				0 - 20
1,1,1,2-Tetrachloroethane	<0.57 ug/L	0.57	0.57		
m,p-Xylene	<1.2 ug/L	1.2	1.2		
m,p-Xylene (RPD)	26.9 %				0 - 20
o-Xylene (RPD)	0.2 %				0 - 20
o-Xylene	0.526 ug/L	0.5	0.5		
Styrene	1.03 ug/L	0.53	0.53		
Styrene (RPD)	0.5 %				0 - 20
Bromoform (RPD)	0 %				0 - 20
Bromoform	<1.56 ug/L	1.56	1.56		
Isopropylbenzene	<0.59 ug/L	0.59	0.59		
Isopropylbenzene (RPD)	0 %				0 - 20
1,1,2,2-Tetrachloroethane (RPD)	0 %				0 - 20
1,1,2,2-Tetrachloroethane	<0.39 ug/L	0.39	0.39		
1,2,3-Trichloropropane	<1.83 ug/L	1.83	1.83		
1,2,3-Trichloropropane (RPD)	0 %				0 - 20
n-Propylbenzene (RPD)	200 %				0 - 20
n-Propylbenzene	<0.49 ug/L	0.49	0.49		
Bromobenzene	<0.5 ug/L	0.5	0.5		
Bromobenzene (RPD)	0 %				0 - 20
1,3,5-Trimethylbenzene (RPD)	0 %				0 - 20
1,3,5-Trimethylbenzene	<0.3 ug/L	0.3	0.3		
2-Chlorotoluene	<0.66 ug/L	0.66	0.66		
2-Chlorotoluene (RPD)	0 %				0 - 20
4-Chlorotoluene (RPD)	0 %				0 - 20
4-Chlorotoluene	<0.8 ug/L	0.8	0.8		
tert-Butylbenzene	<0.85 ug/L	0.85	0.85		
tert-Butylbenzene (RPD)	0 %				0 - 20
1,2,4-Trimethylbenzene (RPD)	0 %				0 - 20
1,2,4-Trimethylbenzene	<0.46 ug/L	0.46	0.46		
sec-Butylbenzene	<0.63 ug/L	0.63	0.63		

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sec-Butylbenzene (RPD)	0 %				0 - 20
p-Isopropyltoluene (RPD)	0 %				0 - 20
p-Isopropyltoluene	<0.59 ug/L	0.59	0.59		
1,3-Dichlorobenzene	<0.7 ug/L	0.7	0.7		
1,3-Dichlorobenzene (RPD)	0 %				0 - 20
1,4-Dichlorobenzene (RPD)	0 %				0 - 20
1,4-Dichlorobenzene	<0.53 ug/L	0.53	0.53		
n-Butylbenzene	<0.72 ug/L	0.72	0.72		
n-Butylbenzene (RPD)	0 %				0 - 20
1,2-Dichlorobenzene (RPD)	0 %				0 - 20
1,2-Dichlorobenzene	<0.7 ug/L	0.7	0.7		
1,2-Dibromo-3-chloropropane	<0.86 ug/L	0.86	0.86		
1,2-Dibromo-3-chloropropane (RPD)	0 %				0 - 20
1,2,4-Trichlorobenzene (RPD)	0 %				0 - 20
1,2,4-Trichlorobenzene	<1.14 ug/L	1.14	1.14		
Naphthalene	<1.53 ug/L	1.53	1.53		
Naphthalene (RPD)	0 %				0 - 20
1,2,3-Trichlorobenzene (RPD)	0 %				0 - 20
1,2,3-Trichlorobenzene	<1.3 ug/L	1.3	1.3		
Dilution Factor	1				
Analyzed By	Jeff Ruehr				
Analysis Date/Time	6/25/2013 2:01 PM				

Palarm Creek Complaint ***LIMS ID: 2013-2215***

Volatiles - water MS ***Run: 1***

Parameter	***Result***	***DL***	***RL***	***Accuracy Control***	***Precision Control***
Dibromofluoromethane (% Recovery)	104 %			70 - 130	
1,2-Dichloroethane-d4 (% Recovery)	110 %			70 - 130	
Toluene-d8 (% Recovery)	89.1 %			70 - 130	
4-Bromofluorobenzene (% Recovery)	104 %			70 - 130	
1,1-Dichloroethene (% Recovery)	93.2 %			70 - 130	
Benzene (% Recovery)	96.7 %			70 - 130	
Trichloroethene (% Recovery)	93.2 %			70 - 130	
Toluene (% Recovery)	81.2 %			70 - 130	
Chlorobenzene (% Recovery)	87.3 %			70 - 130	
Dilution Factor	1				
Analyzed By	Jeff Ruehr				
Analysis Date/Time	6/25/2013 2:27 PM				

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Palarm Creek Complaint — ***LIMS ID: 2013-2215***

Volatiles - water MSD — ***Run: 1***

Parameter	*Result*	*DL*	*RL*	*Accuracy Control*	*Precision Control*
Dibromofluoromethane (% Recovery)	104 %			70 - 130	
1,2-Dichloroethane-d4 (% Recovery)	108 %			70 - 130	
Toluene-d8 (% Recovery)	92.0 %			70 - 130	
4-Bromofluorobenzene (% Recovery)	105 %			70 - 130	
1,1-Dichloroethene (% Recovery)	95.9 %			70 - 130	
1,1-Dichloroethene (RPD)	2.8 %				0 - 20
Benzene (RPD)	1.0 %				0 - 20
Benzene (% Recovery)	97.7 %			70 - 130	
Trichloroethene (RPD)	3.0 %				0 - 20
Trichloroethene (% Recovery)	96.1 %			70 - 130	
Toluene (% Recovery)	83.0 %			70 - 130	
Toluene (RPD)	2.2 %				0 - 20
Chlorobenzene (RPD)	1.1 %				0 - 20
Chlorobenzene (% Recovery)	88.3 %			70 - 130	
Dilution Factor	1				
Analyzed By	Jeff Ruehr				
Analysis Date/Time	6/25/2013 2:52 PM				

LCS — ***LIMS ID: 13062602-LCS-01***

Volatiles - water LCS — ***Run: 1***

Parameter	*Result*	*DL*	*RL*	*Accuracy Control*	*Precision Control*
Dibromofluoromethane (% Recovery)	101 %			70 - 130	
1,2-Dichloroethane-d4 (% Recovery)	109 %			70 - 130	
Toluene-d8 (% Recovery)	91.9 %			70 - 130	
4-Bromofluorobenzene (% Recovery)	98.2 %			70 - 130	
Dichlorodifluoromethane (% Recovery)	112 %			60 - 130	
Chloromethane (% Recovery)	86.2 %			60 - 130	
Vinyl chloride (% Recovery)	105 %			60 - 130	
Bromomethane (% Recovery)	70.2 %			60 - 130	
Chloroethane (% Recovery)	76.7 %			60 - 130	
Trichlorofluoromethane (% Recovery)	96.3 %			60 - 130	
1,1-Dichloroethene (% Recovery)	97.5 %			60 - 130	
Acetone (% Recovery)	127 %			60 - 130	
Methylene chloride (% Recovery)	111 %			60 - 130	
Methyl tert-butyl ether (% Recovery)	99.9 %			60 - 130	

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trans-1,2-Dichloroethene (% Recovery)	97.8 %	60 - 130
1,1-Dichloroethane (% Recovery)	110 %	60 - 130
Methyl ethyl ketone (% Recovery)	119 %	60 - 130
cis-1,2-Dichloroethene (% Recovery)	99.7 %	60 - 130
2,2-Dichloropropane (% Recovery)	103 %	60 - 130
Bromochloromethane (% Recovery)	109 %	60 - 130
Chloroform (% Recovery)	100 %	60 - 130
1,1,1-Trichloroethane (% Recovery)	99.5 %	60 - 130
1,1-Dichloropropene (% Recovery)	102 %	60 - 130
Carbon tetrachloride (% Recovery)	98.2 %	60 - 130
Benzene (% Recovery)	103 %	60 - 130
1,2-Dichloroethane (% Recovery)	109 %	60 - 130
Trichloroethene (% Recovery)	98.7 %	60 - 130
1,2-Dichloropropane (% Recovery)	102 %	60 - 130
Dibromomethane (% Recovery)	103 %	60 - 130
Bromodichloromethane (% Recovery)	96.8 %	60 - 130
cis-1,3-Dichloropropene (% Recovery)	104 %	60 - 130
Methyl isobutyl ketone (% Recovery)	107 %	60 - 130
Toluene (% Recovery)	94.8 %	60 - 130
trans-1,3-Dichloropropene (% Recovery)	103 %	60 - 130
1,1,2-Trichloroethane (% Recovery)	93.3 %	60 - 130
2-Hexanone (% Recovery)	112 %	60 - 130
Tetrachloroethene (% Recovery)	102 %	60 - 130
1,3-Dichloropropane (% Recovery)	99.4 %	60 - 130
Dibromochloromethane (% Recovery)	95.6 %	60 - 130
1,2-Dibromoethane (EDB) (% Recovery)	95.8 %	60 - 130
Chlorobenzene (% Recovery)	99.6 %	60 - 130
Ethylbenzene (% Recovery)	90.1 %	60 - 130
1,1,1,2-Tetrachloroethane (% Recovery)	93.0 %	60 - 130
m,p-Xylene (% Recovery)	95.6 %	60 - 130
o-Xylene (% Recovery)	97.7 %	60 - 130
Styrene (% Recovery)	92.7 %	60 - 130
Bromoform (% Recovery)	109 %	60 - 130
Isopropylbenzene (% Recovery)	109 %	60 - 130
1,1,2,2-Tetrachloroethane (% Recovery)	98.4 %	60 - 130
1,2,3-Trichloropropane (% Recovery)	109 %	60 - 130
n-Propylbenzene (% Recovery)	90.0 %	60 - 130
Bromobenzene (% Recovery)	106 %	60 - 130
1,3,5-Trimethylbenzene (% Recovery)	94.7 %	60 - 130
2-Chlorotoluene (% Recovery)	94.5 %	60 - 130

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4-Chlorotoluene (% Recovery)	96.1 %	60 - 130
tert-Butylbenzene (% Recovery)	94.2 %	60 - 130
1,2,4-Trimethylbenzene (% Recovery)	95.8 %	60 - 130
sec-Butylbenzene (% Recovery)	95.1 %	60 - 130
p-Isopropyltoluene (% Recovery)	89.9 %	60 - 130
1,3-Dichlorobenzene (% Recovery)	97.3 %	60 - 130
1,4-Dichlorobenzene (% Recovery)	92.5 %	60 - 130
n-Butylbenzene (% Recovery)	91.9 %	60 - 130
1,2-Dichlorobenzene (% Recovery)	98.2 %	60 - 130
1,2-Dibromo-3-chloropropane (% Recovery)	106 %	60 - 130
1,2,4-Trichlorobenzene (% Recovery)	99.5 %	60 - 130
Naphthalene (% Recovery)	95.3 %	60 - 130
1,2,3-Trichlorobenzene (% Recovery)	101 %	60 - 130
Dilution Factor	1	
Analyzed By	Jeff Ruehr	
Analysis Date/Time	6/25/2013 11:30	

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Analytical Quality Control Results Report

Batch: 13062503	Oil and Grease - water
Palarm Creek Complaint	LIMS ID: 2013-2215
Oil and Grease - water DUP	Run: 1

Parameter	Result	DL	RL	Accuracy Control	Precision Control
Oil and Grease	<2.5 mg/L	2.5	2.5		
Oil and Grease (RPD)	0 %				0 - 20
Dilution Factor	1				
Analyzed By	Robert Graddy				
Analysis Date/Time	06/25/2013 1000				

Palarm Creek Complaint	LIMS ID: 2013-2215
Oil and Grease - water MS	Run: 1

Parameter	Result	DL	RL	Accuracy Control	Precision Control
Oil and Grease (% Recovery)	102 %			70 - 130	
Dilution Factor	1				
Analyzed By	Robert Graddy				
Analysis Date/Time	06/25/2013 1000				

Palarm Creek Complaint	LIMS ID: 2013-2215
Oil and Grease - water MSD	Run: 1

Parameter	Result	DL	RL	Accuracy Control	Precision Control
Oil and Grease (% Recovery)	101 %			70 - 130	
Oil and Grease (RPD)	1.5 %				0 - 20
Dilution Factor	1				
Analyzed By	Robert Graddy				
Analysis Date/Time	06/25/2013 1000				

MB	LIMS ID: 13062503-MB-01
Oil and Grease - water MB	Run: 1

Parameter	Result	DL	RL	Accuracy Control	Precision Control
Oil and Grease	<2.5 mg/L	2.5	2.5		
Dilution Factor	1				
Analyzed By	Robert Graddy				
Analysis Date/Time	06/25/2013 1000				

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LCS					*LIMS ID: 13062503-LCS-01*
Oil and Grease - water LCS					***Run: 1***
Parameter	***Result***	***DL***	***RL***	***Accuracy Control***	***Precision Control***
Oil and Grease (% Recovery)	101 %			70 - 130	
Dilution Factor	1				
Analyzed By	Robert Graddy				
Analysis Date/Time	06/25/2013 1000				

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Analytical Quality Control Results Report

Batch: 13062504					**Semi-VOA water (Prep)**
Palarm Creek Complaint					***LIMS ID: 2013-2215***
Semi Volatiles - water DUP					***Run: 1***
Parameter	***Result***	***DL***	***RL***	***Accuracy Control***	***Precision Control***
Initial Volume	500 mL				
Final Volume	1 mL				
Prep By	Ed Harris				
Prep Date/Time	6/25/2013 08:00				
2-Fluorophenol (% Recovery)	36.0 %			40 - 110	
Nitrobenzene-d5 (% Recovery)	63.1 %			40 - 125	
2-Fluorobiphenyl (% Recovery)	64.6 %			40 - 110	
2,4,6-Tribromophenol (% Recovery)	58.3 %			40 - 110	
Terphenyl-d14 (% Recovery)	62.1 %			40 - 125	
Methyl Methanesulfonate	<0.2 ug/L	0.2	0.2		
Methyl Methanesulfonate (RPD)	0 %				0 - 40
Ethyl methanesulfonate (RPD)	0 %				0 - 40
Ethyl methanesulfonate	<0.2 ug/L	0.2	0.2		
Phenol	<0.2 ug/L	0.2	0.2		
Phenol (RPD)	38.2 %				0 - 40
Aniline (RPD)	3.0 %				0 - 40
Aniline	<0.2 ug/L	0.2	0.2		
Bis(2-chloroethyl)ether	<0.24 ug/L	0.2	0.24		
Bis(2-chloroethyl)ether (RPD)	0 %				0 - 40
2-Chlorophenol (RPD)	0 %				0 - 40
2-Chlorophenol	<0.2 ug/L	0.2	0.2		
1,3-Dichlorobenzene	<0.12 ug/L	0.12	0.12		
1,3-Dichlorobenzene (RPD)	0 %				0 - 40
1,4-Dichlorobenzene (RPD)	0 %				0 - 40
1,4-Dichlorobenzene	<0.12 ug/L	0.12	0.12		
Benzyl alcohol	0.202 ug/L	0.16	0.16		
Benzyl alcohol (RPD)	2.1 %				0 - 40
1,2-Dichlorobenzene (RPD)	0 %				0 - 40
1,2-Dichlorobenzene	<0.12 ug/L	0.12	0.12		
2-Methylphenol	<0.1 ug/L	0.1	0.1		
2-Methylphenol (RPD)	0 %				0 - 40
Acetophenone (RPD)	17.3 %				0 - 40
Acetophenone	<0.1 ug/L	0.1	0.1		

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4-Methylphenol	<0.1 ug/L	0.1	0.1		
4-Methylphenol (RPD)	0 %				0 - 40
N-Nitrosodi-n-propylamine (RPD)	0 %				0 - 40
N-Nitrosodi-n-propylamine	<0.2 ug/L	0.2	0.2		
Hexachloroethane	<0.2 ug/L	0.2	0.2		
Hexachloroethane (RPD)	0 %				0 - 40
Nitrobenzene (RPD)	0 %				0 - 40
Nitrobenzene	<0.2 ug/L	0.2	0.2		
N-Nitrosopiperidine	<0.2 ug/L	0.2	0.2		
N-Nitrosopiperidine (RPD)	0 %				0 - 40
Isophorone (RPD)	0 %				0 - 40
Isophorone	<0.1 ug/L	0.1	0.1		
2-Nitrophenol	<0.3 ug/L	0.3	0.3		
2-Nitrophenol (RPD)	0 %				0 - 40
2,4-Dimethylphenol (RPD)	200 %				0 - 40
2,4-Dimethylphenol	<0.1 ug/L	0.1	0.1		
Bis(2-chloroethoxy)methane	<0.2 ug/L	0.2	0.2		
Bis(2-chloroethoxy)methane (RPD)	0 %				0 - 40
2,4-Dichlorophenol (RPD)	0 %				0 - 40
2,4-Dichlorophenol	<0.2 ug/L	0.2	0.2		
1,2,4-Trichlorobenzene	<0.12 ug/L	0.12	0.12		
1,2,4-Trichlorobenzene (RPD)	0 %				0 - 40
Naphthalene (RPD)	0 %				0 - 40
Naphthalene	<0.08 ug/L	0.08	0.08		
4-Chloroaniline	<0.1 ug/L	0.1	0.1		
4-Chloroaniline (RPD)	0 %				0 - 40
2,6-Dichlorophenol (RPD)	0 %				0 - 40
2,6-Dichlorophenol	<0.2 ug/L	0.2	0.2		
Hexachlorobutadiene	<0.2 ug/L	0.2	0.2		
Hexachlorobutadiene (RPD)	0 %				0 - 40
N-Nitrosodibutylamine	<0.24 ug/L	0.2	0.24		
N-Nitrosodibutylamine (RPD)	0 %				0 - 40
4-Chloro-3-methylphenol (RPD)	0 %				0 - 40
4-Chloro-3-methylphenol	<0.16 ug/L	0.16	0.16		
2-Methylnaphthalene	<0.1 ug/L	0.1	0.1		
2-Methylnaphthalene (RPD)	0 %				0 - 40
1,2,4,5-Tetrachlorobenzene (RPD)	0 %				0 - 40
1,2,4,5-Tetrachlorobenzene	<0.1 ug/L	0.1	0.1		
Hexachlorocyclopentadiene	<0.16 ug/L	0.16	0.16		
Hexachlorocyclopentadiene (RPD)	0 %				0 - 40

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2,4,6-Trichlorophenol (RPD)	0 %				0 - 40
2,4,6-Trichlorophenol	<0.4 ug/L	0.2	0.4		
2,4,5-Trichlorophenol	<0.2 ug/L	0.2	0.2		
2,4,5-Trichlorophenol (RPD)	0 %				0 - 40
2-Chloronaphthalene (RPD)	0 %				0 - 40
2-Chloronaphthalene	<0.1 ug/L	0.1	0.1		
1-Chloronaphthalene	<0.1 ug/L	0.1	0.1		
1-Chloronaphthalene (RPD)	0 %				0 - 40
2-Nitroaniline (RPD)	0 %				0 - 40
2-Nitroaniline	<0.2 ug/L	0.2	0.2		
Dimethyl phthalate	<0.2 ug/L	0.2	0.2		
Dimethyl phthalate (RPD)	0 %				0 - 40
2,6-Dinitrotoluene (RPD)	0 %				0 - 40
2,6-Dinitrotoluene	<0.2 ug/L	0.2	0.2		
Acenaphthylene	<0.08 ug/L	0.08	0.08		
Acenaphthylene (RPD)	0 %				0 - 40
3-Nitroaniline (RPD)	0 %				0 - 40
3-Nitroaniline	<0.2 ug/L	0.2	0.2		
Acenaphthene	<0.1 ug/L	0.1	0.1		
Acenaphthene (RPD)	0 %				0 - 40
2,4-Dinitrophenol (RPD)	0 %				0 - 40
2,4-Dinitrophenol	<4 ug/L	4	4		
Pentachlorobenzene	<0.12 ug/L	0.12	0.12		
Pentachlorobenzene (RPD)	0 %				0 - 40
4-Nitrophenol (RPD)	0 %				0 - 40
4-Nitrophenol	<2 ug/L	2	2		
Dibenzofuran	<0.1 ug/L	0.1	0.1		
Dibenzofuran (RPD)	200 %				0 - 40
2,4-Dinitrotoluene (RPD)	0 %				0 - 40
2,4-Dinitrotoluene	<0.2 ug/L	0.2	0.2		
2,3,4,6-Tetrachlorophenol	<0.6 ug/L	0.6	0.6		
2,3,4,6-Tetrachlorophenol (RPD)	0 %				0 - 40
Diethyl phthalate (RPD)	2.5 %				0 - 40
Diethyl phthalate	<0.4 ug/L	0.2	0.4		
Fluorene	<0.1 ug/L	0.1	0.1		
Fluorene (RPD)	0 %				0 - 40
4-Chlorophenyl phenyl ether (RPD)	0 %				0 - 40
4-Chlorophenyl phenyl ether	<0.1 ug/L	0.1	0.1		
4-Nitroaniline	<0.3 ug/L	0.2	0.3		
4-Nitroaniline (RPD)	0 %				0 - 40

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4,6-Dinitro-2-methylphenol (RPD)	0 %				0 - 40
4,6-Dinitro-2-methylphenol	<6 ug/L	6	6		
Diphenylamine	<0.1 ug/L	0.1	0.1		
Diphenylamine (RPD)	0 %				0 - 40
Azobenzene (RPD)	0 %				0 - 40
Azobenzene	<0.08 ug/L	0.08	0.08		
4-Bromophenyl phenyl ether	<0.2 ug/L	0.2	0.2		
4-Bromophenyl phenyl ether (RPD)	0 %				0 - 40
Hexachlorobenzene (RPD)	0 %				0 - 40
Hexachlorobenzene	<0.16 ug/L	0.16	0.16		
Pentachlorophenol	<1 ug/L	1	1		
Pentachlorophenol (RPD)	0 %				0 - 40
Pentachloronitrobenzene (RPD)	0 %				0 - 40
Pentachloronitrobenzene	<0.2 ug/L	0.2	0.2		
Pronamide	<0.2 ug/L	0.2	0.2		
Pronamide (RPD)	0 %				0 - 40
Phenanthrene (RPD)	1.3 %				0 - 40
Phenanthrene	<0.08 ug/L	0.08	0.08		
Anthracene	<0.08 ug/L	0.08	0.08		
Anthracene (RPD)	0 %				0 - 40
Carbazole (RPD)	0 %				0 - 40
Carbazole	<0.1 ug/L	0.1	0.1		
Di-n-butyl phthalate	<0.2 ug/L	0.2	0.2		
Di-n-butyl phthalate (RPD)	7.0 %				0 - 40
Fluoranthene (RPD)	0 %				0 - 40
Fluoranthene	<0.08 ug/L	0.08	0.08		
Pyrene	<0.08 ug/L	0.08	0.08		
Pyrene (RPD)	0 %				0 - 40
Dimethylaminoazobenzene (RPD)	0 %				0 - 40
Dimethylaminoazobenzene	<0.2 ug/L	0.2	0.2		
Butyl benzyl phthalate	<0.4 ug/L	0.4	0.4		
Butyl benzyl phthalate (RPD)	0 %				0 - 40
Benzo (a) anthracene (RPD)	0 %				0 - 40
Benzo (a) anthracene	<0.1 ug/L	0.1	0.1		
Chrysene	<0.1 ug/L	0.1	0.1		
Chrysene (RPD)	0 %				0 - 40
Bis(2-ethylhexyl)phthalate (RPD)	28.4 %				0 - 40
Bis(2-ethylhexyl)phthalate	<0.3 ug/L	0.3	0.3		
Di-n-octyl phthalate	<0.3 ug/L	0.3	0.3		
Di-n-octyl phthalate (RPD)	28.3 %				0 - 40

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Benzo (b) fluoranthene (RPD)	0 %				0 - 40
Benzo (b) fluoranthene	<0.16 ug/L	0.16	0.16		
7,12-Dimethylbenz (a) anthracene	<0.2 ug/L	0.2	0.2		
7,12-Dimethylbenz (a) anthracene (RPD)	0 %				0 - 40
Benzo (k) fluoranthene (RPD)	0 %				0 - 40
Benzo (k) fluoranthene	<0.16 ug/L	0.16	0.16		
Benzo (a) pyrene	<0.16 ug/L	0.16	0.16		
Benzo (a) pyrene (RPD)	0 %				0 - 40
3-Methylcholanthrene (RPD)	0 %				0 - 40
3-Methylcholanthrene	<0.2 ug/L	0.2	0.2		
Indeno (1,2,3-cd) pyrene	<0.2 ug/L	0.2	0.2		
Indeno (1,2,3-cd) pyrene (RPD)	0 %				0 - 40
Dibenzo (a,h) anthracene (RPD)	0 %				0 - 40
Dibenzo (a,h) anthracene	<0.16 ug/L	0.16	0.16		
Benzo (g,h,i) perylene	<0.16 ug/L	0.16	0.16		
Benzo (g,h,i) perylene (RPD)	0 %				0 - 40
Dilution Factor	1				
Analyzed By	Ed Harris				
Analysis Date/Time	6/25/2013 12:21 PM				

Palarm Creek Complaint ***LIMS ID: 2013-2215***

Semi Volatiles - water MS ***Run: 1***

Parameter	***Result***	***DL***	***RL***	***Accuracy Control***	***Precision Control***
Initial Volume	500 mL				
Final Volume	1 mL				
Prep By	Ed Harris				
Prep Date/Time	6/25/2013 08:00				
2-Fluorophenol (% Recovery)	35.3 %			40 - 110	
Nitrobenzene-d5 (% Recovery)	78.2 %			40 - 125	
2-Fluorobiphenyl (% Recovery)	70.0 %			40 - 125	
2,4,6-Tribromophenol (% Recovery)	66.6 %			40 - 125	
Terphenyl-d14 (% Recovery)	62.9 %			40 - 125	
Phenol (% Recovery)	21.9 %			25 - 125	
2-Chlorophenol (% Recovery)	36.0 %			25 - 125	
1,4-Dichlorobenzene (% Recovery)	63.6 %			25 - 125	
N-Nitrosodi-n-propylamine (% Recovery)	68.4 %			25 - 125	
1,2,4-Trichlorobenzene (% Recovery)	62.7 %			25 - 125	
4-Chloro-3-methylphenol (% Recovery)	38.6 %			25 - 125	
Acenaphthene (% Recovery)	67.9 %			25 - 125	

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4-Nitrophenol (% Recovery)	28.8 %			25 - 125	
2,4-Dinitrotoluene (% Recovery)	65.6 %			25 - 125	
Pentachlorophenol (% Recovery)	55.6 %			25 - 125	
Pyrene (% Recovery)	55.7 %			25 - 125	
Dilution Factor	1				
Analyzed By	Ed Harris				
Analysis Date/Time	6/25/2013 12:50 PM				

Palarm Creek Complaint ***LIMS ID: 2013-2215***

Semi Volatiles - water MSD ***Run: 1***

Parameter	***Result***	***DL***	***RL***	***Accuracy Control***	***Precision Control***
Initial Volume	500 mL				
Final Volume	1 mL				
Prep By	Ed Harris				
Prep Date/Time	6/25/2013 08:00				
2-Fluorophenol (% Recovery)	37.2 %			40 - 110	
Nitrobenzene-d5 (% Recovery)	68.1 %			40 - 125	
2-Fluorobiphenyl (% Recovery)	61.0 %			40 - 125	
2,4,6-Tribromophenol (% Recovery)	69.0 %			40 - 125	
Terphenyl-d14 (% Recovery)	60.2 %			40 - 125	
Phenol (% Recovery)	23.8 %			25 - 125	
Phenol (RPD)	8.5 %				0 - 40
2-Chlorophenol (% Recovery)	40.0 %			25 - 125	
2-Chlorophenol (RPD)	10.4 %				0 - 40
1,4-Dichlorobenzene (RPD)	7.4 %				0 - 40
1,4-Dichlorobenzene (% Recovery)	59.0 %			25 - 125	
N-Nitrosodi-n-propylamine (% Recovery)	68.3 %			25 - 125	
N-Nitrosodi-n-propylamine (RPD)	0.2 %				0 - 40
1,2,4-Trichlorobenzene (RPD)	0.5 %				0 - 40
1,2,4-Trichlorobenzene (% Recovery)	62.3 %			25 - 125	
4-Chloro-3-methylphenol (% Recovery)	45.9 %			25 - 125	
4-Chloro-3-methylphenol (RPD)	17.3 %				0 - 40
Acenaphthene (% Recovery)	64.4 %			25 - 125	
Acenaphthene (RPD)	5.2 %				0 - 40
4-Nitrophenol (% Recovery)	31.0 %			25 - 125	
4-Nitrophenol (RPD)	7.3 %				0 - 40
2,4-Dinitrotoluene (% Recovery)	63.2 %			25 - 125	
2,4-Dinitrotoluene (RPD)	3.7 %				0 - 40
Pentachlorophenol (% Recovery)	55.4 %			25 - 125	

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Pentachlorophenol (RPD)	0.3 %		0 - 40
Pyrene (% Recovery)	58.8 %	25 - 125	
Pyrene (RPD)	5.5 %		0 - 40
Dilution Factor	1		
Analyzed By	Ed Harris		
Analysis Date/Time	6/25/2013 1:19 PM		

MB ***LIMS ID: 13062504-MB-01***

Semi Volatiles - water MB ***Run: 1***

Parameter	*Result*	*DL*	*RL*	*Accuracy Control*	*Precision Control*
Initial Volume	500 mL				
Final Volume	1 mL				
Prep By	Ed Harris				
Prep Date/Time	6/25/2013 08:00				
2-Fluorophenol (% Recovery)	36.4 %			40 - 110	
Nitrobenzene-d5 (% Recovery)	65.4 %			40 - 125	
2-Fluorobiphenyl (% Recovery)	60.7 %			40 - 125	
2,4,6-Tribromophenol (% Recovery)	60.4 %			40 - 125	
Terphenyl-d14 (% Recovery)	79.7 %			40 - 125	
Methyl Methanesulfonate	<0.2 ug/L	0.2	0.2		
Ethyl methanesulfonate	<0.2 ug/L	0.2	0.2		
Phenol	<0.2 ug/L	0.2	0.2		
Aniline	<0.2 ug/L	0.2	0.2		
Bis(2-chloroethyl)ether	<0.2 ug/L	0.2	0.2		
2-Chlorophenol	<0.2 ug/L	0.2	0.2		
1,3-Dichlorobenzene	<0.12 ug/L	0.12	0.12		
1,4-Dichlorobenzene	<0.12 ug/L	0.12	0.12		
Benzyl alcohol	<0.16 ug/L	0.16	0.16		
1,2-Dichlorobenzene	<0.12 ug/L	0.12	0.12		
2-Methylphenol	<0.1 ug/L	0.1	0.1		
Acetophenone	<0.1 ug/L	0.1	0.1		
4-Methylphenol	<0.1 ug/L	0.1	0.1		
N-Nitrosodi-n-propylamine	<0.2 ug/L	0.2	0.2		
Hexachloroethane	<0.2 ug/L	0.2	0.2		
Nitrobenzene	<0.2 ug/L	0.2	0.2		
N-Nitrosopiperidine	<0.2 ug/L	0.2	0.2		
Isophorone	<0.1 ug/L	0.1	0.1		
2-Nitrophenol	<0.3 ug/L	0.3	0.3		
2,4-Dimethylphenol	<0.1 ug/L	0.1	0.1		

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Bis(2-chloroethoxy)methane	<0.2 ug/L	0.2	0.2
2,4-Dichlorophenol	<0.2 ug/L	0.2	0.2
1,2,4-Trichlorobenzene	<0.12 ug/L	0.12	0.12
Naphthalene	<0.08 ug/L	0.08	0.08
4-Chloroaniline	<0.1 ug/L	0.1	0.1
2,6-Dichlorophenol	<0.2 ug/L	0.2	0.2
Hexachlorobutadiene	<0.2 ug/L	0.2	0.2
N-Nitrosodibutylamine	<0.2 ug/L	0.2	0.2
4-Chloro-3-methylphenol	<0.16 ug/L	0.16	0.16
2-Methylnaphthalene	<0.12 ug/L	0.12	0.12
1,2,4,5-Tetrachlorobenzene	<0.1 ug/L	0.1	0.1
Hexachlorocyclopentadiene	<0.16 ug/L	0.16	0.16
2,4,6-Trichlorophenol	<0.2 ug/L	0.2	0.2
2,4,5-Trichlorophenol	<0.2 ug/L	0.2	0.2
2-Chloronaphthalene	<0.1 ug/L	0.1	0.1
1-Chloronaphthalene	<0.1 ug/L	0.1	0.1
2-Nitroaniline	<0.2 ug/L	0.2	0.2
Dimethyl phthalate	<0.2 ug/L	0.2	0.2
2,6-Dinitrotoluene	<0.2 ug/L	0.2	0.2
Acenaphthylene	<0.08 ug/L	0.08	0.08
3-Nitroaniline	<0.2 ug/L	0.2	0.2
Acenaphthene	<0.1 ug/L	0.1	0.1
2,4-Dinitrophenol	<4 ug/L	4	4
Pentachlorobenzene	<0.12 ug/L	0.12	0.12
4-Nitrophenol	<2 ug/L	2	2
Dibenzofuran	<0.1 ug/L	0.1	0.1
2,4-Dinitrotoluene	<0.2 ug/L	0.2	0.2
2,3,4,6-Tetrachlorophenol	<0.6 ug/L	0.6	0.6
Diethyl phthalate	<0.2 ug/L	0.2	0.2
Fluorene	<0.1 ug/L	0.1	0.1
4-Chlorophenyl phenyl ether	<0.1 ug/L	0.1	0.1
4-Nitroaniline	<0.2 ug/L	0.2	0.2
4,6-Dinitro-2-methylphenol	<6 ug/L	6	6
Diphenylamine	<0.1 ug/L	0.1	0.1
Azobenzene	<0.08 ug/L	0.08	0.08
4-Bromophenyl phenyl ether	<0.2 ug/L	0.2	0.2
Hexachlorobenzene	<0.16 ug/L	0.16	0.16
Pentachlorophenol	<1 ug/L	1	1
Pentachloronitrobenzene	<0.2 ug/L	0.2	0.2
Pronamide	<0.2 ug/L	0.2	0.2

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Phenanthrene	<0.08 ug/L	0.08	0.08	
Anthracene	<0.08 ug/L	0.08	0.08	
Carbazole	<0.1 ug/L	0.1	0.1	
Di-n-butyl phthalate	0.637 ug/L	0.2	0.2	
Fluoranthene	<0.08 ug/L	0.08	0.08	
Pyrene	<0.08 ug/L	0.08	0.08	
Dimethylaminoazobenzene	<0.2 ug/L	0.2	0.2	
Butyl benzyl phthalate	<0.3 ug/L	0.3	0.3	
Benzo (a) anthracene	<0.2 ug/L	0.2	0.2	
Chrysene	<0.1 ug/L	0.1	0.1	
Bis(2-ethylhexyl)phthalate	<0.3 ug/L	0.3	0.3	
Di-n-octyl phthalate	<0.3 ug/L	0.3	0.3	
Benzo (b) fluoranthene	<0.16 ug/L	0.16	0.16	
7,12-Dimethylbenz (a) anthracene	<0.2 ug/L	0.2	0.2	
Benzo (k) fluoranthene	<0.16 ug/L	0.16	0.16	
Benzo (a) pyrene	<0.16 ug/L	0.16	0.16	
3-Methylcholanthrene	<0.2 ug/L	0.2	0.2	
Indeno (1,2,3-cd) pyrene	<0.2 ug/L	0.2	0.2	
Dibenzo (a,h) anthracene	<0.16 ug/L	0.16	0.16	
Benzo (g,h,i) perylene	<0.16 ug/L	0.16	0.16	
Dilution Factor	1			
Analyzed By	Ed Harris			
Analysis Date/Time	6/25/2013 10:53			

LCS ***LIMS ID: 13062504-LCS-01***

Semi Volatiles - water LCS ***Run: 1***

Parameter	***Result***	***DL***	***RL***	***Accuracy Control***	***Precision Control***
Initial Volume	500 mL				
Final Volume	1 mL				
Prep By	Ed Harris				
Prep Date/Time	6/25/2013 08:00				
2-Fluorophenol (% Recovery)	42.2 %			40 - 110	
Nitrobenzene-d5 (% Recovery)	71.7 %			50 - 125	
2-Fluorobiphenyl (% Recovery)	66.1 %			50 - 125	
2,4,6-Tribromophenol (% Recovery)	71.0 %			40 - 125	
Terphenyl-d14 (% Recovery)	73.4 %			50 - 125	
Methyl Methanesulfonate (% Recovery)	70.4 %			50 - 150	
Ethyl methanesulfonate (% Recovery)	80.7 %			50 - 150	
Phenol (% Recovery)	48.1 %			50 - 150	

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Aniline (% Recovery)	64.8 %	50 - 150
Bis(2-chloroethyl)ether (% Recovery)	79.1 %	50 - 150
2-Chlorophenol (% Recovery)	79.4 %	50 - 150
1,3-Dichlorobenzene (% Recovery)	60.6 %	50 - 150
1,4-Dichlorobenzene (% Recovery)	62.6 %	50 - 150
Benzyl alcohol (% Recovery)	96.0 %	50 - 150
1,2-Dichlorobenzene (% Recovery)	65.7 %	50 - 150
2-Methylphenol (% Recovery)	78.1 %	50 - 150
Acetophenone (% Recovery)	85.4 %	50 - 150
4-Methylphenol (% Recovery)	76.8 %	50 - 150
N-Nitrosodi-n-propylamine (% Recovery)	84.8 %	50 - 150
Hexachloroethane (% Recovery)	54.7 %	50 - 150
Nitrobenzene (% Recovery)	83.9 %	50 - 150
N-Nitrosopiperidine (% Recovery)	96.0 %	50 - 150
Isophorone (% Recovery)	90.3 %	50 - 150
2-Nitrophenol (% Recovery)	85.6 %	50 - 150
2,4-Dimethylphenol (% Recovery)	8.5 %	50 - 150
Bis(2-chloroethoxy)methane (% Recovery)	80.0 %	50 - 150
2,4-Dichlorophenol (% Recovery)	85.1 %	50 - 150
1,2,4-Trichlorobenzene (% Recovery)	65.1 %	50 - 150
Naphthalene (% Recovery)	74.1 %	50 - 150
4-Chloroaniline (% Recovery)	75.8 %	50 - 150
2,6-Dichlorophenol (% Recovery)	86.5 %	50 - 150
Hexachlorobutadiene (% Recovery)	57.6 %	50 - 150
N-Nitrosodibutylamine (% Recovery)	101 %	50 - 150
4-Chloro-3-methylphenol (% Recovery)	87.1 %	50 - 150
2-Methylnaphthalene (% Recovery)	76.7 %	50 - 150
1,2,4,5-Tetrachlorobenzene (% Recovery)	72.4 %	50 - 150
Hexachlorocyclopentadiene (% Recovery)	66.7 %	50 - 150
2,4,6-Trichlorophenol (% Recovery)	81.0 %	50 - 150
2,4,5-Trichlorophenol (% Recovery)	83.6 %	50 - 150
2-Chloronaphthalene (% Recovery)	78.7 %	50 - 150
1-Chloronaphthalene (% Recovery)	83.5 %	50 - 150
2-Nitroaniline (% Recovery)	90.5 %	50 - 150
Dimethyl phthalate (% Recovery)	86.4 %	50 - 150
2,6-Dinitrotoluene (% Recovery)	86.6 %	50 - 150
Acenaphthylene (% Recovery)	80.5 %	50 - 150
3-Nitroaniline (% Recovery)	93.0 %	50 - 150
Acenaphthene (% Recovery)	81.6 %	50 - 150
2,4-Dinitrophenol (% Recovery)	108 %	50 - 150

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501-682-0955

Pentachlorobenzene (% Recovery)	77.6 %	50 - 150
4-Nitrophenol (% Recovery)	44.0 %	50 - 150
Dibenzofuran (% Recovery)	82.4 %	50 - 150
2,4-Dinitrotoluene (% Recovery)	90.6 %	50 - 150
2,3,4,6-Tetrachlorophenol (% Recovery)	90.9 %	50 - 150
Diethyl phthalate (% Recovery)	94.7 %	50 - 150
Fluorene (% Recovery)	78.3 %	50 - 150
4-Chlorophenyl phenyl ether (% Recovery)	79.9 %	50 - 150
4-Nitroaniline (% Recovery)	92.3 %	50 - 150
4,6-Dinitro-2-methylphenol (% Recovery)	100 %	50 - 150
Diphenylamine (% Recovery)	79.9 %	50 - 150
Azobenzene (% Recovery)	75.6 %	50 - 150
4-Bromophenyl phenyl ether (% Recovery)	74.3 %	50 - 150
Hexachlorobenzene (% Recovery)	75.1 %	50 - 150
Pentachlorophenol (% Recovery)	80.6 %	50 - 150
Pentachloronitrobenzene (% Recovery)	81.6 %	50 - 150
Pronamide (% Recovery)	97.4 %	50 - 150
Phenanthrene (% Recovery)	81.2 %	50 - 150
Anthracene (% Recovery)	82.8 %	50 - 150
Carbazole (% Recovery)	100 %	50 - 150
Di-n-butyl phthalate (% Recovery)	130 %	50 - 150
Fluoranthene (% Recovery)	92.7 %	50 - 150
Pyrene (% Recovery)	79.7 %	50 - 150
Dimethylaminoazobenzene (% Recovery)	91.2 %	50 - 150
Butyl benzyl phthalate (% Recovery)	96.8 %	50 - 150
Benzo (a) anthracene (% Recovery)	94.8 %	50 - 150
Chrysene (% Recovery)	90.0 %	50 - 150
Bis(2-ethylhexyl)phthalate (% Recovery)	96.8 %	50 - 150
Di-n-octyl phthalate (% Recovery)	102 %	50 - 150
Benzo (b) fluoranthene (% Recovery)	102 %	50 - 150
7,12-Dimethylbenz (a) anthracene (% Recovery)	83.4 %	50 - 150
Benzo (k) fluoranthene (% Recovery)	105 %	50 - 150
Benzo (a) pyrene (% Recovery)	98.3 %	50 - 150
3-Methylcholanthrene (% Recovery)	95.9 %	50 - 150
Indeno (1,2,3-cd) pyrene (% Recovery)	85.5 %	50 - 150
Dibenzo (a,h) anthracene (% Recovery)	91.7 %	50 - 150
Benzo (g,h,i) perylene (% Recovery)	87.7 %	50 - 150
Dilution Factor	1	
Analyzed By	Ed Harris	
Analysis Date/Time	6/25/2013 11:22	

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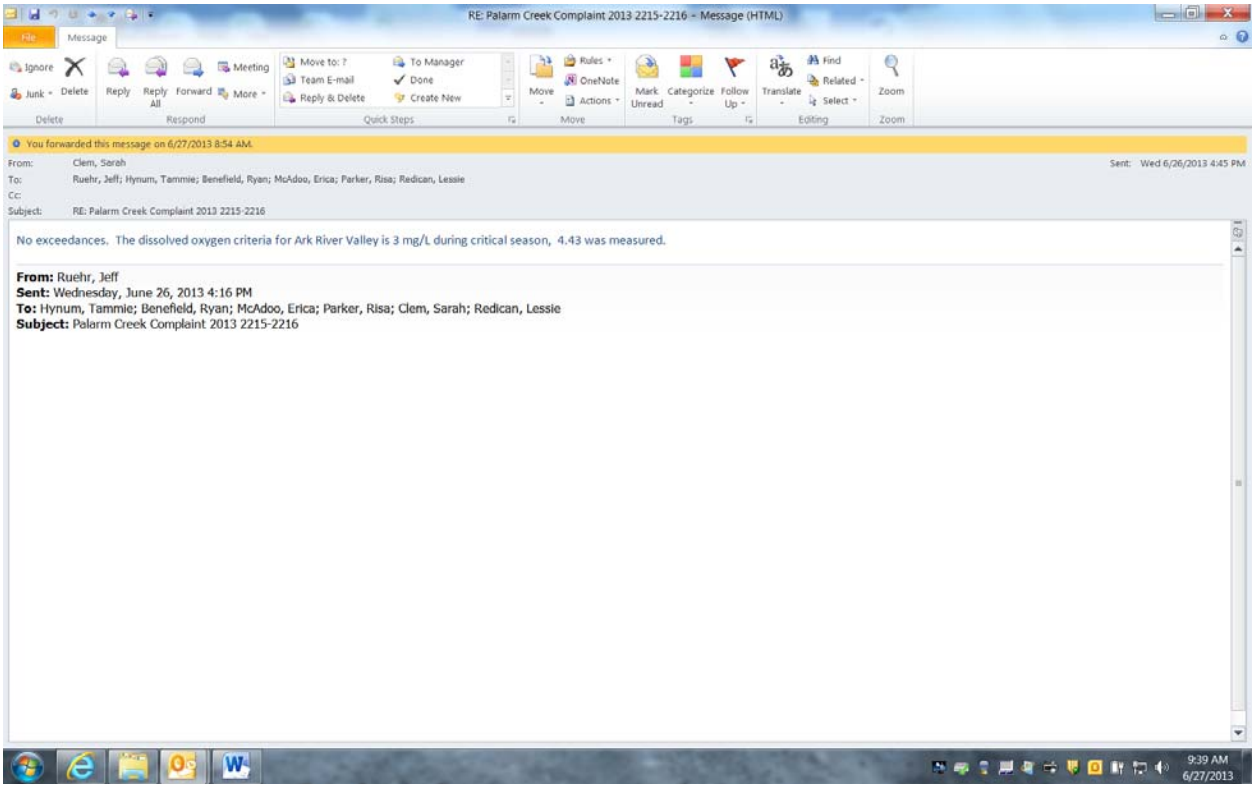
RE: Palarm Creek Complaint 2013 2215-2216 - Message (HTML)
You forwarded this message on 6/27/2013 8:54 AM.
From: Clem, Sarah
Sent: Wed 6/26/2013 4:45 PM
To: Ruehr, Jeff; Hynum, Tammie; Benefield, Ryan; McAdoo, Erica; Parker, Risa; Redican, Lessie
Cc:
Subject: RE: Palarm Creek Complaint 2013 2215-2216
No exceedances. The dissolved oxygen criteria for Ark River Valley is 3 mg/L during critical season, 4.43 was measured.
From: Ruehr, Jeff
Sent: Wednesday, June 26, 2013 4:16 PM
To: Hynum, Tammie; Benefield, Ryan; McAdoo, Erica; Parker, Risa; Clem, Sarah; Redican, Lessie
Subject: Palarm Creek Complaint 2013 2215-2216
9:39 AM
6/27/2013