



**CONESTOGA-ROVERS
& ASSOCIATES**

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December 15, 2011

Reference No. 622237

Mr. Bruce Wanstall
Alaska Department of Environmental Conservation
410 Willoughby Avenue, Suite 302
Juneau, Alaska 99801

Re: Annual 2011 Groundwater Monitoring and
Ozone System Maintenance Report
Chevron Site 8-2307
9203 Cessna Drive
Juneau, Alaska
ADEC File ID 1513.26.046

Dear Mr. Bruce Wanstall:

Conestoga-Rovers & Associates (CRA) is submitting this *Annual 2011 Groundwater Monitoring and Ozone System Maintenance Report* for the site referenced above (Figure 1) on behalf of Chevron Environmental Management Company. CRA performed groundwater monitoring and sampling in accordance with the Alaska Department of Environmental Conservation's May 2010 *Draft Field Guidance*. The *Groundwater Elevation Map* is presented on Figure 2. Groundwater monitoring and sampling data are presented in Table 1. CRA's *Annual 2011* monitoring data package is included as Attachment A. Lancaster Laboratories' June 28, 2011 *Analytical Results* are included as Attachment B. Site photos are included as Attachment C. Standard operation procedures for groundwater monitoring and sampling are included as Attachment D and the ADEC laboratory data review checklist and memorandum are included as Attachment E.

RESULTS OF ANNUAL 2011 EVENT

On May 26, 2011, CRA monitored and sampled the site well per the established schedule.

Results of the current monitoring event indicate the following:

- Depth to Water 4.46 feet below grade

Equal
Employment Opportunity
Employer



December 15, 2011

Reference No. 622237

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Results of the current sampling event are presented below in Table A:

TABLE A: GROUNDWATER ANALYTICAL DATA SUMMARY						
<i>Well ID</i>	<i>DRO (mg/l)</i>	<i>GRO (mg/l)</i>	<i>Benzene (mg/l)</i>	<i>Toluene (mg/l)</i>	<i>Ethylbenzene (mg/l)</i>	<i>Total Xylenes (mg/l)</i>
<i>Table C Cleanup Levels</i>	1.5	2.2	0.005	1.0	0.7	10
MW-4	19/24	0.54/0.56	NA	NA	NA	NA
NA Not Analyzed						

OZONE SYSTEM MAINTENANCE AND GEOCHEMICAL ANALYSES

System Maintenance

CRA installed an ozone injection system in monitoring well MW-4 on August 28, 2007 to reduce dissolved-phase DRO concentrations in groundwater. CRA upgraded the ozone emitter on May 10, 2010. CRA replaced one broken solar panel, after which, each individual solar panel and battery was tested to confirm proper operation. Individual battery output was approximately 14.5 volts per battery. Solar panel output was approximately 0.3 volts per panel. The system was inspected on May 26, 2011 and all components were found to be safe and operational. All implemented safety equipment (i.e. orange safety cones, snow poles, and caution tape) were intact and visible. The westernmost panel was cracked and will be replaced in 2012.. Site photos are presented in Attachment C.

Groundwater Geochemical Results

CRA collected field geochemical parameters and groundwater samples on May 26, 2011 to evaluate natural attenuation and the ozone injection system's effectiveness. Temperature, dissolved oxygen (DO), oxidation reduction potential (ORP), conductivity and pH were measured in the field (Attachment A). Groundwater samples were analyzed for carbon dioxide, sulfate, nitrite, nitrate, and alkalinity. The geochemical analytical results are presented in Table 2.



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CONCLUSIONS

Low concentrations of nitrate (0.04 mg/L) and sulfate (2.4 mg/L) indicate hydrocarbon biodegradation is active. High alkalinity and a reducing environment also indicate a high level of hydrocarbon consumption. DRO and GRO concentrations have remained stable.

ANTICIPATED FUTURE ACTIVITIES

Groundwater Monitoring

CRA will continue annual monitoring and sampling and submit a groundwater monitoring and sampling report presenting the results.

Alaska Qualified Personnel in accordance with *Title 18 Alaska Administrative Code (AAC) 78, Articles 2, 6, and 9* conducted all project work.

Please contact John Riggi at (720) 975-9121 if you have any questions or require additional information.

Sincerely,

CONESTOGA-ROVERS & ASSOCIATES

Diane Escobedo
Staff Geologist

John Riggi, P.G.
Senior Project Geologist

Jeffrey Cloud
Chemist

JR/aa/4
Encl.



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Figure 1	Vicinity Map
Figure 2	Groundwater Elevation Map
Table 1	Groundwater Analytical Results
Table 2	Geochemical Analytical Results
Attachment A	Monitoring Data Package
Attachment B	Laboratory Analytical Report
Attachment C	Site Photos
Attachment D	Standard Operation Procedures for Groundwater Monitoring and Sampling
Attachment E	ADEC Laboratory Data Review Checklist and Memorandum

cc: Dan Carrier, Chevron EMC (*electronic copy*)
Oge Nkenke (Chevron Filing)
Bev Niemann
Allen Hesse

FIGURES

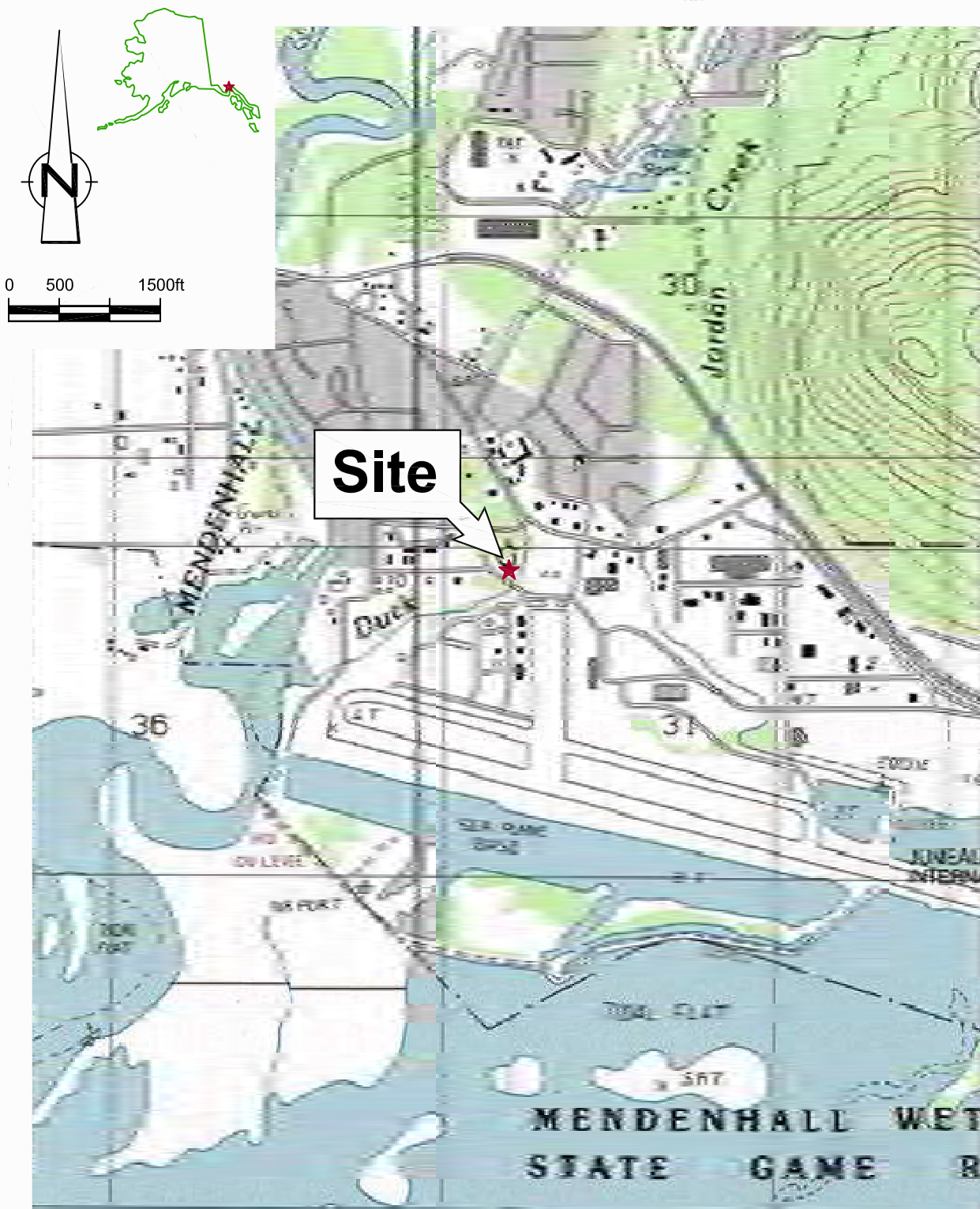


Figure 1
VICINITY MAP
FORMER DELTA WESTERN BULK TERMINAL 8-2307
9203 Cessna Drive
Juneau, Alaska



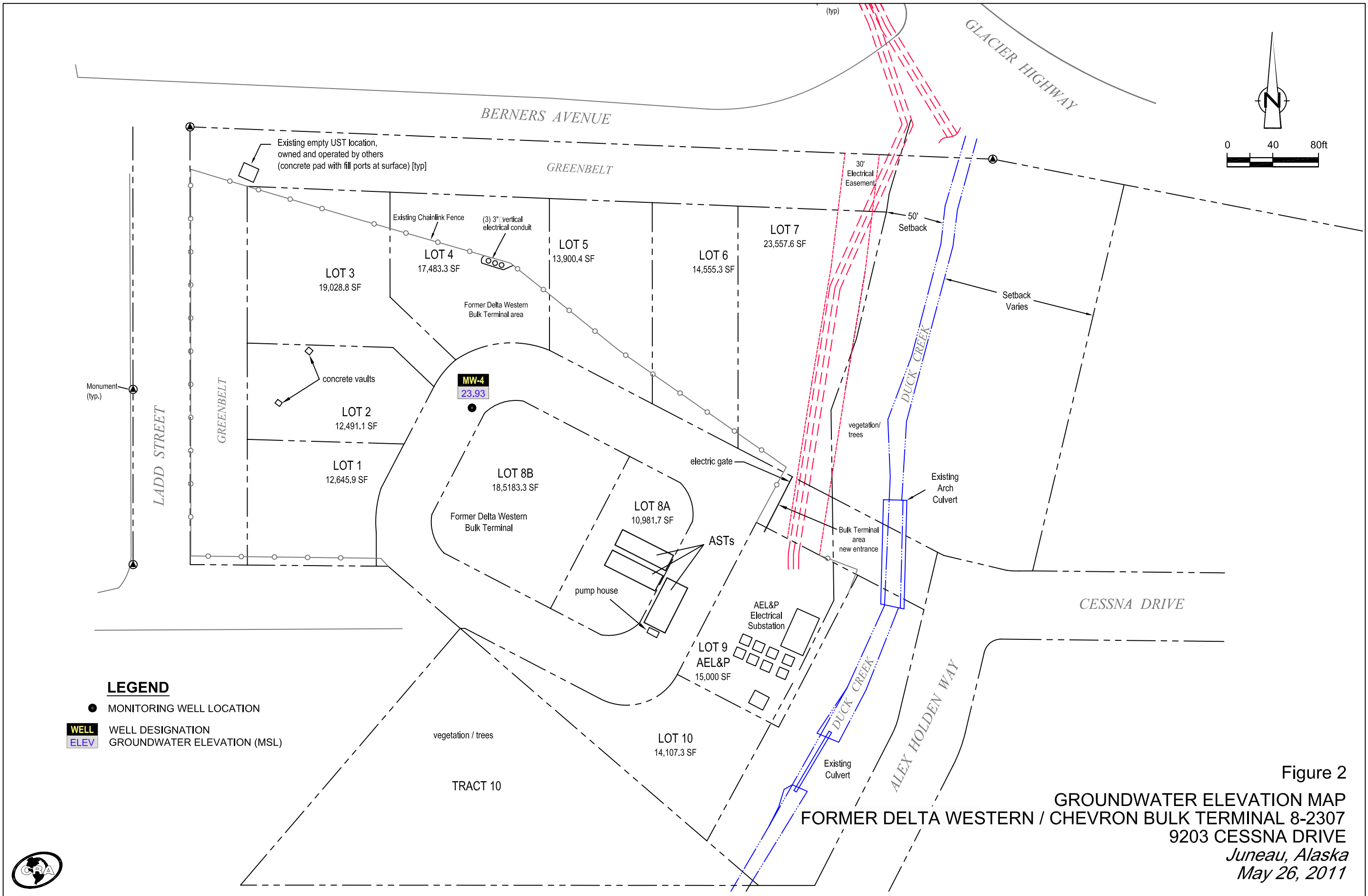


Figure 2

GROUNDWATER ELEVATION MAP
 FORMER DELTA WESTERN / CHEVRON BULK TERMINAL 8-2307
 9203 CESSNA DRIVE
 Juneau, Alaska
 May 26, 2011



TABLES

TABLE 1

GROUNDWATER ANALYTICAL RESULTS
DELTA WESTERN/FORMER CHEVRON BULK TERMINAL 8-2307
9203 CESSNA DRIVE
JUNEAU, ALASKA

Location	Date	TOC	DTW	GWE	HYDROCARBONS			PRIMARY VOCs				ADDITIONAL VOC'S
					RRO	DRO	GRO	Benzene	Toluene	Ethyl-benzene	Total Xylenes	VOCs
					mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L
ADEC Groundwater Cleanup Levels ^a					1.1	1.5	2.2	0.005	1.0	0.7	10.0	
MW-1	06/06/2000	25.19	8.55	16.64	<0.75	0.269	<0.05	<0.0005	<0.0005	<0.0005	<0.001	ND
MW-1	12/09/2000	25.19	7.72	17.47	<0.75	0.201	<0.05	0.00025	0.00082	<0.0005	<0.001	-
MW-1	03/24/2001	25.19	8.48	16.71	-	0.298	<0.05	0.00214	<0.0005	<0.0005	<0.001	-
MW-1	06/19/2001	25.19	8.93	16.26	-	0.121	<0.05	0.0002	<0.0005	<0.0005	<0.001	-
MW-1	06/17/2002	25.19	8.60	16.59	-	0.181	<0.05	0.00115	<0.0005	<0.0005	<0.001	-
MW-1	12/11/2002	25.19	7.53	17.66	-	0.1	<0.05	<0.0002	<0.0005	<0.0005	<0.001	-
MW-1	06/25/2003	25.19	4.72	20.47	-	0.1	<0.05	0.00457	<0.0005	<0.0005	<0.001	-
MW-1	12/09/2003	25.19	8.43	16.76	-	0.139	<0.05	<0.0002	<0.0005	<0.0005	<0.001	-
MW-1	05/18/2004	25.19	9.38	15.81	-	0.391	<0.05	<0.0002	<0.0005	<0.0005	<0.001	-
MW-1	05/27/2005	Well Destroyed										
MW-2	06/06/2000	28.73	13.20	15.53	<0.75	0.22	<0.05	<0.0005	<0.0005	<0.0005	<0.001	ND
MW-2	12/09/2000	28.73	12.12	16.61	<0.75 / <0.75	0.22	<0.05 / <0.05	0.000395 / 0.000353	0.000951 / 0.001	<0.0005 / <0.0005	<0.001 / 0.001	-
MW-2	03/24/2001	28.73	13.28	15.45	-	0.176	<0.05	<0.0002	<0.0005	<0.0005	<0.001	-
MW-2	06/19/2001	28.73	13.72	15.01	-	0.274	0.058	0.000213	<0.0005	<0.0005	0.0011	-
MW-2	06/17/2002	28.73	13.13	15.60	-	0.393	<0.05	<0.0002	<0.0005	<0.0005	<0.001	-
MW-2	12/11/2002	28.73	9.00	19.73	-	0.159	<0.05	<0.0002	<0.0005	<0.0005	0.001	-
MW-2	06/25/2003	28.73	14.34	14.39	-	0.209	0.08	<0.0002	<0.0005	<0.0005	<0.001	-
MW-2	12/09/2003	28.73	13.15	15.58	-	0.132	<0.05	<0.0002	<0.0005	<0.0005	0.001	-
MW-2	05/18/2004	28.73	9.40	19.33	-	0.391	<0.05	<0.0002	0.00062	<0.0005	0.00101	-
MW-2	05/27/2005	Well Destroyed										
MW-3	06/06/2000	28.21	12.09	16.12	<0.75	0.144	<0.05	<0.0005	<0.0005	<0.0005	<0.001	ND
MW-3	12/10/2000	28.21	11.29	16.92	<0.75	0.439	<0.05	0.000223	<0.0005	<0.0005	<0.001	-
MW-3	03/24/2001	28.21	12.11	16.10	-	0.188	<0.05 / <0.05	<0.0002 / <0.0002	<0.0005 / <0.0005	<0.0005 / <0.0005	<0.001 / <0.001	-
MW-3	06/19/2001	28.21	12.53	15.68	-	0.163	<0.05 / <0.05	<0.0002 / <0.0002	<0.0005 / <0.0005	<0.0005 / <0.0005	<0.001 / <0.001	-
MW-3	06/17/2002	28.21	12.18	16.03	-	0.105	<0.05	<0.0002	<0.0005	<0.0005	<0.001	-
MW-3	12/11/2002	28.21	11.00	17.21	-	0.122	<0.05	<0.0002	<0.0005	<0.0005	<0.001	-
MW-3	06/25/2003	28.21	13.26	14.95	-	0.1	<0.05	0.00067	<0.0005	<0.0005	<0.001	-
MW-3	12/09/2003	28.21	11.98	16.23	-	0.186	<0.05	<0.0002	<0.0005	<0.0005	<0.001	-
MW-3	05/18/2004	28.21	12.97	15.24	-	0.391	<0.05	<0.0002	<0.0005	0.000528	0.00162	-
MW-3	05/27/2005	Well Destroyed										
MW-4	06/19/2001	28.39	4.08	24.31	-	22.2	0.948	0.00148	<0.00125	0.00398	0.0821	-
MW-4	06/17/2002	28.39	4.17	24.22	-	33.2	1.05	0.001	<0.0005	0.0517	0.0979	-
MW-4	12/11/2002	28.39	2.25	26.14	-	29.3	0.921	0.0091	0.00125	0.0448	0.088	-
MW-4	06/25/2003	28.39	4.14	24.25	-	22.2	0.833	0.00128	0.00107	0.0485	0.0795	-

TABLE 1

GROUNDWATER ANALYTICAL RESULTS
DELTA WESTERN/FORMER CHEVRON BULK TERMINAL 8-2307
9203 CESSNA DRIVE
JUNEAU, ALASKA

Location	Date	TOC	DTW	GWE	HYDROCARBONS			PRIMARY VOCs			ADDITIONAL VOC'S	
					RRO	DRO	GRO	Benzene	Toluene	Ethyl-benzene	Total Xylenes	VOCs
					Units	ft msl	fbg	ft msl	mg/L	mg/L	mg/L	mg/L
ADEC Groundwater Cleanup Levels ^a					1.1	1.5	2.2	0.005	1.0	0.7	10.0	
MW-4	12/09/2003	28.39	3.66	24.73	-	27.5	0.537	0.00149	<0.0025	0.0517	0.0547	-
MW-4	05/18/2004	28.39	4.74	23.65	-	12.8	2.2	0.00139	<0.001	0.0492	0.0786	-
MW-4	05/27/2005	28.39	5.50	22.89	-	24	0.51 / 0.52	0.0008 / 0.0008	<0.0005 / <0.0005	0.037 / 0.038	0.04 / 0.041	-
MW-4	06/01/2006	28.39	5.80	22.59	-	26	0.55	<0.0005	<0.0005	0.035	0.045	-
MW-4	08/28/2007	28.39	4.18	24.21	-	17	0.5 / 0.5	<0.001 / <0.001	<0.001 / <0.001	0.02 / 0.02	0.05 / 0.04	-
MW-4	06/24/2008	28.39	5.20	23.19	-	37	-	<0.001 / <0.001	0.001 / 0.001	0.02 / 0.02	0.03 / 0.03	-
MW-4	08/25/2008	28.39	2.23	26.16	-	9.12	-	<0.0005 / <0.0005	<0.0005 / <0.0005	0.0161 / 0.0142	0.0264 / 0.0211	-
MW-4	06/23/2009	28.39	6.41	21.98	-	12.2 / 13.1	0.578 / 0.476	-	-	-	-	-
MW-4	05/11/2010	28.39	4.44	23.95	-	17 / 13	0.57 J / 0.81 J	-	-	-	-	-
MW-4	05/26/2011	28.39	4.46	23.93	-	19 / 24	0.54 / 0.56	-	-	-	-	-
Trip Blank	12/09/2000	-	-	-	-	-	<0.05	<0.0002	<0.0005	<0.0005	<0.001	-
Trip Blank	03/24/2001	-	-	-	-	-	<0.05	<0.0002	<0.0005	<0.0005	<0.001	-
Trip Blank	06/19/2001	-	-	-	-	-	<0.05	<0.0002	<0.0005	<0.0005	<0.001	-
Trip Blank-1	12/09/2003	-	-	-	-	0.165	<0.05 / <0.05	<0.0002 / <0.0002	<0.0005 / <0.0005	<0.0005 / <0.0005	<0.001 / <0.001	-
Trip Blank-2	12/09/2003	-	-	-	-	-	<0.05	<0.0002	<0.0005	<0.0005	<0.001	-
Trip Blank	05/18/2004	-	-	-	-	-	<0.01	<0.0005	<0.0005	<0.0005	<0.0015	-
Trip Blank	06/01/2006	-	-	-	-	-	<0.01	<0.0005	<0.0005	<0.0005	<0.0015	-
Trip Blank	08/28/2007	-	-	-	-	-	<0.01	<0.001	<0.001	<0.001	<0.002	<0.0005
Trip Blank	06/24/2008	-	-	-	-	-	-	<0.001	<0.001	<0.001	<0.002	-
Trip Blank	08/25/2008	-	-	-	-	-	-	<0.0005	<0.0005	<0.0005	<0.001	-
Trip Blank	06/23/2009	-	-	-	-	-	<0.0100	-	-	-	-	-
Trip Blank	05/11/2010	-	-	-	-	-	<0.010	-	-	-	-	-

Notes and Abbreviations

VOCs = Volatile Organic Compounds

TOC = Top of Casing

DTW = Depth to Water

GWE = Groundwater Elevation

RRO = Residual Range Organics AK103

DRO = Diesel Range Organics by Alaska Series Method AK102

GRO = Gasoline Range Organics by Alaska Series Method AK101

Benzene, Toluene, Ethylbenzene, and Total Xylenes by Environmental Protection Agency (EPA) Method 8021B or 8260B

Total Xylenes = Sum of m-, o-, and p-xylenes

ft msl = Feet Above Mean Sea Level

fbg = Feet Below Grade

TABLE 1

GROUNDWATER ANALYTICAL RESULTS
 DELTA WESTERN/FORMER CHEVRON BULK TERMINAL 8-2307
 9203 CESSNA DRIVE
 JUNEAU, ALASKA

Location	Date	TOC	DTW	GWE	HYDROCARBONS			Benzene	PRIMARY VOCs			ADDITIONAL VOC'S	
					RRO	DRO	GRO		Toluene	Ethyl-benzene	Total Xylenes	VOCs	
	Units	ft msl	fbg	ft msl	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L
ADEC Groundwater Cleanup Levels ^a					1.1	1.5	2.2	0.005	1.0	0.7	10.0		

mg/L = Milligrams per Liter

ADEC = Alaska Department of Environmental Conservation

^a = Levels established in ADEC Table C Groundwater Cleanup Levels (18 AAC 75.345)

<x = Constituent not detected above x milligrams per liter

- = Not Measured/Not Analyzed

ND = Not detected above laboratory method detection limits

x / y = Sample Results / Blind Duplicate Results

BOLD = Indicates concentration above the ADEC Table C Groundwater Cleanup Level

TABLE 2

GEOCHEMICAL ANALYTICAL RESULTS
DELTA WESTERN/FORMER CHEVRON BULK TERMINAL 8-2307
9203 CESSNA DRIVE
JUNEAU, ALASKA

Location	Date	MNA PARAMETERS											
		Ferrous Iron	Nitrite/Nitrate	Nitrate	Nitrite	Carbon Dioxide	Alkalinity,	Alkalinity,	Alkalinity,	Alkalinity,	Sulfate	pH*	Conductivity*
							Total (as CaCO3)	Bicarbonate	Carbonate	Hydroxide			
							mg/L	mg/L	mg/L	mg/L			
ADEC Groundwater Cleanup Levels ^a													
MW-4	06/24/2008	--	--	--	--	--	--	--	--	--	--	6.3	393.5
MW-4	08/26/2008	3.9	0.531	--	--	81.8	149	149	5	5	11.7	6.65	383.8
MW-4	10/12/2009	--	0.0910 J	--	--	--	137	137	<0.640	<0.640	0.920 J	6.33	--
MW-4	05/11/2010	--	--	--	--	--	--	--	--	--	3.5 J / 3.6 J	6.91	-74.8
MW-4	05/26/2011	--	--	0.040 UJ	<0.015	89	127	--	--	--	2.4 J	6.40	0.279

Notes and Abbreviations

MNA = Monitored Natural Attenuation

CaCO₃ = Calcium Carbonate

* = Average of readings during purging

mg/L = Milligrams per Liter

ADEC = Alaska Department of Environmental Conservation

^a = Levels established in ADEC Table C Groundwater Cleanup Levels (18 AAC 75.345)

-- = Not Measured/Not Analyzed

J = Estimated value

ATTACHMENT A

MONITORING DATA PACKAGE

Groundwater Monitoring Field Sheet

Well ID	Time	DTP	DTW	Depth to Bottom	Product Thickness	Amount of Product Removed	Casing Diam.	Comments
MW-4	13:25	—	4.46	9.45	—	—	2"	

Project Name: 8-2307Project Number/Task: 622237/GWSTechnician: D. ESCOBEDO/N. GRECODate: 5/26/11



CONESTOGA-ROVERS
& ASSOCIATES

WELL SAMPLING FORM

DISPOSABLE BAILER SAMPLING

Site ID: 8-2307	CRA Mgr: John Riggi	Well ID: MW-4
CRA Project No.: 622237	Date: 5/26/11	Field Staff: DE/NG
Street Address: 9203 Cessna Dr.	City, State: Juneau, AK	Purging Device: <input type="radio"/> Teflon Disp. Bailer <input checked="" type="radio"/> poly bailer Sampling Method: <input type="radio"/> Teflon Disp. Bailer <input checked="" type="radio"/> poly bailer
Depth to Water: 4.46	Depth to Bottom: 9.45	Water Column Height: 4.99
Volume/ft: 0.16	1 Casing Volume: 0.79	3 Casing Volumes: 2.40
Well Diameter: 2"	Did Well Dewater?: no	Total Gallons Purged: 2.5
Start Purge Time: 13:32	Stop Purge Time: 13:38	Total Time: 6 minutes

1 Casing Volume = Water column height x Volume/ft.

Well Diam.	Volume/ft (gallons)
2"	0.16
4"	0.65
6"	1.47

NO PURGE APPROVED BY ADEC? ☐ YES ☒ NO (If NO, please enter parameters below.)

Time	Volume Purged (gallons)	Temp. (°C) ± 10%	DO ± 10%	pH ± 0.1	Cond. (mS) ± 3%	ORP (mv) ± 10	Comments
13:33	2 1/4 gal	9.95	85.5	6.20	0.691	-208.3	
13:34	3 1/4 gal	9.00	34.3	6.40	0.309	-196.3	
13:35	2	8.70	27.7	6.42	0.279	-186.1	
13:37	2.5	8.80	31.5	6.38	0.278	-176.4	

*** A minimum of three parameters must be monitored and recorded. ***

NOTE: If well is purged dry, DO NOT collect sample until it has recharged to approximately 80% of its pre-purge volume.

FIELD KIT RESULTS: Ferrous Iron _____ mg/L Nitrate _____ mg/L

Sample ID	Date	Time	Analytes / Analytical Method
MW-4 - 052611	5/26/11	13:20	<input checked="" type="radio"/> DRO by AK102 <input type="radio"/> SVOCs by TCL8270 <input type="radio"/> PAHs by 8270 <input type="radio"/> RRO by AK103 <input type="radio"/> Lead by 6010 <input type="radio"/> PAHs by 8270SIM <input checked="" type="radio"/> Alkalinity by 310.1 <input type="radio"/> Methane by RSK175 <input checked="" type="radio"/> Nitrate/Nitrite by 353.2 <input checked="" type="radio"/> Sulfate by 300 <input checked="" type="radio"/> CO ₂ <input type="radio"/> <input checked="" type="radio"/> GRO by AK101 <input type="radio"/> BTEX by 8260B <input type="radio"/> EDB by 8011 <input type="radio"/> BTEX by 8021B <input type="radio"/> MtBE by 8260B <input type="radio"/> 1,2-DCA by 8260B <input type="radio"/> HVOCs by 8260B
Additional Comments:			

ATTACHMENT B

LABORATORY ANALYTICAL REPORT



2425 New Holland Pike, PO Box 12425, Lancaster, PA 17605-2425 • 717-656-2300 Fax: 717-656-2681 • www.lancasterlabs.com

Type III Data Package

Prepared for:

ChevronTexaco
6001 Bollinger Canyon Rd L4310
San Ramon CA 94583

Project: 82307
Water Samples
Collected on 05/26/11

SDG# AKE89

GROUP	SAMPLE NUMBERS
1249125	6301430-6301432

PA Cert. # 36-00037
NY Cert. # 10670
NJ Cert. # PA011
NC Cert. # 521
TX Cert. # T104704194-08A-TX

Through our technical processes and second person review of data, we have established that our data/deliverables are in compliance with the methods and project requirements unless otherwise noted or previously resolved with the client.

Authorized by:

Date: 06/28/2011

Dana M. Kauffman
Manager

Any questions or concerns you might have regarding this data package should be directed to your client representative, Natalie Luciano at Ext. 1881.

Total Number of Pages

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SDG# AKE89

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Sample Reference List for SDG Number AKE89
with a Data Package Type of III
10880 - ChevronTexaco
Project: 82307

Lab Sample Number	Lab Sample Code	Client Sample Description
6301430	CJMW5	MW-4-052611 Grab Water Sample Facility# 82307
6301431	CJDU1	DUP-1-052611 Grab Water Sample Facility# 82307
6301432	CJTB1	Trip_Blank-1-052611 Water Sample Facility# 82307

Chevron Generic Analysis Request/Chain of Custody



For Lancaster Laboratories use only
 Acct. # 10880 Sample # 6301430-32

SCR#

DL 7649

LAB

Analyses Requested

Grp # 1248125Facility #: Chevron 8-2307Site Address: 9203 Lessem Dr.Chevron PM: Dan Carrier Lead Consultant: CRAConsultant/Office: CRA DenverConsultant Pjt. Mgr.: John RiggConsultant Phone #: 720 975 9121 Fax #: 720 975 9150Sampler: Diane EscobedoService Order #: ☐ Non SAR:

Sample Identification

MA-04-052011

PAP-1-052011

Trip Blank-1-052011

Matrix

☐ Potable
☐ NPDESSoil ☐ Air ☐

Water

Oil ☐

Total Number of Containers

BTX + MTBE 8021 ☐ 8260 ☐ Naphth ☐

8260 full scan

Sulfate

Oxygenates

TPH & Alkalinity to pH 4.5

TPH ☐ Extended Rtg.
☐ Silica Gel CleanupLead Total ☐ Diss. ☐ MethodVPH/EPH ☐ CO₂ by headspaceNWTPH H₂CO₃ ☐ quantification

DRO AK 102

GRO AK 101

Nitrate Nitrogen

Nitrite Nitrogen

Preservation Codes

H = HCl T = Thioisulfate
N = HNO₃ B = NaOH
S = H₂SO₄ O = Other☐ J value reporting needed☐ Must meet lowest detection limits possible for 8260 compounds

8021 MTBE Confirmation

☐ Confirm MTBE + Naphthalene☐ Confirm highest hit by 8260☐ Confirm all hits by 8260☐ Run oxy's on highest hit☐ Run oxy's on all hits

Comments / Remarks

Report MDL

mg/L

Relinquished by: Date: 5/27/11 Time: 0800Relinquished by:



Environmental Sample Administration Receipt Documentation Log

Client/Project: CRA

Shipping Container Sealed: YES NO

Date of Receipt: 5/28/11

Custody Seal Present *: YES NO

Time of Receipt: 0900

* Custody seal was intact unless otherwise noted in the discrepancy section

Source Code: 50-1

Package: Chilled Not Chilled

Temperature of Shipping Containers							
Cooler #	Thermometer ID	Temperature (°C)	Temp Bottle (TB) or Surface Temp (ST)	Wet Ice (WI) or Dry Ice (DI) or Ice Packs (IP)	Ice Present? Y/N	Loose (L) Bagged Ice (B) or NA	Comments
1	0419261	4.9°C	TB	WI	Y	B	
2							
3							
4							
5							
6							

Number of Trip Blanks received NOT listed on chain of custody 0

Paperwork Discrepancy/Unpacking Problems:

REL. 2 EXTRA MIN-04 VIALS & MISSING 2 DUP-01 VIALS
① kmz 5-28-11

Unpacker Signature/Emp#:

1454

Date/Time: 5/28/11 / 0935

AKES9: 8883

Issued by Dept. 6042 Management

Batchlog Summary 11153A53A

QC	ID	Sample Code	Amt	SS/IS S	Amt (mL)	MS Sol.	Amt (mL)	FV (mL)	SW	DF	PH	BC	Comments
6297939MS	AA	FTA01MS	1.00	SS1112525A	0.0002	MS1112625A	0.002001	1.00	1.00	≤ 2		026A	
6297939MS (AB)	AB	FTA01MS	1.00	SS1112525A	0.0002	MS1115125A	0.000220	1.00	1.00	≤ 2		026B	
6297940MSD	AA	FTA01MSI	1.00	SS1112525A	0.0002	MS1112625A	0.002001	1.00	1.00	≤ 2		026A	
6297940MSD (AB)	AB	FTA01MSI	1.00	SS1112525A	0.0002	MS1115125A	0.000220	1.00	1.00	≤ 2		026B	
BLANKA	AA	BLKFI	1.00	SS1112525A	0.0002			1.00	1.00				
LCSA	AA	LCSE4	1.00	SS1112525A	0.0002	MS1112625A	0.002001	1.00	1.00				
LCSB	AA	LCSE5	1.00	SS1112525A	0.0002	MS1115125A	0.000220	1.00	1.00				

Sample#	ID	Sample Code	Amt	SS/IS Std.	Amt (mL)	FV (mL)	SW	DF	PH	BC	HS	Due Date	Hold Date	P	Analyses	Comments
6297938	AA	FTA01	1.00	SS1112525A	0.0002	1.00		1.00	≤ 2	026A	□	6/7	6/7	N	02102	
6297942	AA	FTA02	1.00	SS1112525A	0.0002	1.00		1.00	≤ 2	026A	□	6/7	6/7	N	02102	
6297943	AA	FTAEB	1.00	SS1112525A	0.0002	1.00		1.00	≤ 2	026A	□	6/7	6/7	N	02102	
6297944	AA	FTATB	1.00	SS1112525A	0.0002	1.00		1.00	≤ 2	026A	□	6/7	6/7	N	02102	
6297949	AA	FTEBA	1.00	SS1112525A	0.0002	1.00		1.00	≤ 2	026A	□	6/7	6/7	N	02102	
6297950	AA	FTTBA	1.00	SS1112525A	0.0002	1.00		1.00	≤ 2	026A	□	6/7	6/7	N	02102	
6298649	AA	14573	1.00	SS1112525A	0.0010	1.00		5.00	≤ 2	041B	□	8/8		P	02102	
6299141	AA	GAAM8	1.00	SS1112525A	0.0002	1.00		1.00	≤ 2	104A	□	6/6	6/7	P	01440 02102	
6299142	AA	GAAD1	1.00	SS1112525A	0.0002	1.00		1.00	≤ 2	104A	□	6/6	6/7	P	01440 02102	
6299143	AA	GAAM9	1.00	SS1112525A	0.0002	1.00		1.00	≤ 2	104A	□	6/6	6/7	P	01440 02102	
6299144	AA	GAATB	1.00	SS1112525A	0.0002	1.00		1.00	≤ 2	104A	□	6/6	6/7	P	01440 02102	
6301430	AA	CJMW5	1.00	SS1112525A	0.0002	1.00		1.00	≤ 2	104A	□	6/8	6/9	P	01440	
6301431	AA	CJDU1	1.00	SS1112525A	0.0002	1.00		1.00	≤ 2	104A	□	6/8	6/9	P	01440	
6301432	AA	CJTB1	1.00	SS1112525A	0.0002	1.00		1.00	≤ 2	104A	□	6/8	6/9	P	01440	

Spike Solutions:

MS1112625A Waters MI working Spike
 MS1115125A Waters GRO Spike #2
 SS1112525A Waters 2 Component Sur. Sol.

Analyst:

cem1991

Date:

6/3/11

Verifier:

MDJ2001

Date:

6/3/11

Comments

ANALYSIS 06/03/11

01146 GC VOA Water Prep

An undiluted aliquot of the water sample or a dilution of the sample is purged with an inert gas and the volatiles are collected on an adsorbent trap that is subsequently desorbed onto a gas chromatographic column.

Reference: Test Methods for Evaluating Solid Wastes, SW-846 Method 5030B, December 1996.

01438 TPH-GRO AK water C6-C10

The volatile compounds are extracted by bubbling an inert gas through the sample and collecting them on a sorbent trap. The trap is thermally desorbed onto a capillary column and analysis is performed using gas chromatography with a flame ionization detector (FID) and, optionally, a photoionization detector (PID) in series. Quantitation for Gasoline Range Organics (GRO) is performed using the total peak area detected within the hydrocarbon range defined in the method.

Reference: Method AK101 for the Determination of Gasoline Range Organics, April 8, 2002

00219 Nitrite Nitrogen

Nitrite ions react with sulfanilamide to yield a diazo compound which couples with N-1-naphthylethylene diamine dihydrochloride to form a soluble, highly-colored dye. The result is determined colorimetrically.

Reference: Method 353.2, Methods for Chemical Analysis of Water and Wastes USEPA 600, Revision 2.0, 1993

00220 Nitrate Nitrogen

Nitrate ions are reduced to nitrite by passing through a cadmium coil. The nitrite ions then react with sulfanilamide to yield a diazo compound which couples with N-1-naphthylethylene diamine dihydrochloride to form a soluble, highly-colored dye. The result is determined colorimetrically.

Reference: Method 353.2, Methods for Chemical Analysis of Water and Wastes USEPA 600, Revision 2.0, 1993

00228 Sulfate

A small volume of sample is introduced into an ion chromatograph. The anions are then separated and measured by a system consisting of a guard column, separator column, suppressor, and conductivity detector.

Reference: Method 300.0, Methods for Chemical Analysis of Water and Wastes USEPA 600, Revision 2.1, 1993

00201 Alkalinity to pH 8.3

Alkalinity is determined by titrating the sample with standardized sulfuric acid to pH of 8.3 for the phenolphthalein alkalinity.

Reference: Standard Methods for the Examination of Water and Wastewater, 20th Edition, 1998, Method 2320 B

00202 Alkalinity to pH 4.5

Alkalinity is determined by titrating the sample with standardized sulfuric acid to a pH of 4.5 for the total alkalinity.

Reference: Standard Methods for the Examination of Water and Wastewater, 20th Edition, 1998, Method 2320 B

01741 TPH-DRO AK water C10-C25

Sample extracts in methylene chloride are analyzed by capillary chromatography using flame ionization detection. Quantitation is performed using the total peak area detected within the hydrocarbon ranges defined in the method.

Reference: Alaska Method 102/103 for Determination of Diesel Range Organics, April 8, 2002.

11184 AK DRO Waters Extraction

An aliquot of sample is extracted with methylene chloride using either separatory funnel extraction or micro extraction technique.

Reference: Alaska Method 102/103 for Determination of Diesel Range Organics, April 8, 2002.

08097 CO2 by Headspace

An aliquot of sample is placed in a headspace vial and warmed to 35C. A portion of the headspace is analyzed on a gas chromatograph using a capillary column and thermal conductivity detection.

Reference: Test Methods for Evaluating Solid Wastes SW-846, Method 8015B Modified, December 1996.



ANALYTICAL RESULTS

Prepared by:

Lancaster Laboratories
2425 New Holland Pike
Lancaster, PA 17605-2425

Prepared for:

ChevronTexaco
6001 Bollinger Canyon Rd L4310
San Ramon CA 94583

June 07, 2011

Project: 82307

Submittal Date: 05/28/2011

Group Number: 1249125

SDG: AKE89

PO Number: 0015074818

Release Number: CARRIER

State of Sample Origin: AK

Client Sample Description

MW-4-052611 Grab Water Sample
DUP-1-052611 Grab Water Sample
Trip_Blank-1-052611 Water Sample

Lancaster Labs (LLI) #

6301430
6301431
6301432

The specific methodologies used in obtaining the enclosed analytical results are indicated on the Laboratory Sample Analysis Record.

ELECTRONIC COPY TO	CRA	Attn: Nick Greco
ELECTRONIC COPY TO	Chevron	Attn: CRA EDD
ELECTRONIC COPY TO	CRA	Attn: Jeffrey Cloud
ELECTRONIC COPY TO	CRA	Attn: Sarah Gillette
ELECTRONIC COPY TO	Chevron	Attn: Anna Avina
ELECTRONIC COPY TO	CRA	Attn: Diane Escobedo
ELECTRONIC COPY TO	CRA	Attn: John Riggi
1 COPY TO	Data Package Group	

AKE89 0007



Questions? Contact your Client Services Representative
Natalie R Luciano at (717) 656-2300 Ext. 1881

Respectfully Submitted,

Robert Heisey

Robert Heisey
Senior Specialist

AKES9 0000



Sample Description: MW-4-052611 Grab Water Sample
 Facility# 82307
 9203 Cessna Dr - Juneau, AK

LLI Sample # WW 6301430
 LLI Group # 1249125
 Account # 10880

Project Name: 82307

Collected: 05/26/2011 13:20 by DE

ChevronTexaco

6001 Bollinger Canyon Rd L4310
 San Ramon CA 94583

Submitted: 05/28/2011 09:00

Reported: 06/07/2011 10:11

CJMW5 SDG#: AKE89-01

CAT No.	Analysis Name	CAS Number	As Received Result	As Received Method Detection Limit*	As Received Limit of Quantitation	Dilution Factor
GC Volatiles	AK 101		mg/l	mg/l	mg/l	
01438	TPH-GRO AK water C6-C10	n.a.	0.54	0.010	0.10	1
GC Extractable TPH	AK 102/AK 103		mg/l	mg/l	mg/l	
	04/08/02					
01741	TPH-DRO AK water C10-C25	n.a.	19	1.2	6.2	25
GC Miscellaneous	SW-846 8015B modified		mg/l	mg/l	mg/l	
08097	CO2 by Headspace	124-38-9	89	4.0	12	1
Wet Chemistry	EPA 300.0		mg/l	mg/l	mg/l	
00228	Sulfate	14808-79-8	2.4 J	1.5	5.0	5
	EPA 353.2		mg/l	mg/l	mg/l	
00220	Nitrate Nitrogen	14797-55-8	N.D.	0.040	0.10	1
00219	Nitrite Nitrogen	14797-65-0	N.D.	0.015	0.050	1
	SM20 2320 B		mg/l as CaCO3	mg/l as CaCO3	mg/l as CaCO3	
00202	Alkalinity to pH 4.5	n.a.	127	0.46	2.0	1
00201	Alkalinity to pH 8.3	n.a.	N.D.	0.46	2.0	1

General Sample Comments

State of Alaska Lab Certification No. UST-061

All QC is compliant unless otherwise noted. Please refer to the Quality Control Summary for overall QC performance data and associated samples.

Laboratory Sample Analysis Record

CAT No.	Analysis Name	Method	Trial#	Batch#	Analysis Date and Time	Analyst	Dilution Factor
01438	TPH-GRO AK water C6-C10	AK 101	1	11153A53A	06/02/2011 20:03	Carrie E Miller	1
01146	GC VOA Water Prep	SW-846 5030B	1	11153A53A	06/02/2011 20:03	Carrie E Miller	1
01741	TPH-DRO AK water C10-C25	AK 102/AK 103	1	111520013A	06/03/2011 22:10	Heather E Williams	25
		04/08/02					
08097	CO2 by Headspace	SW-846 8015B modified	1	111520033A	06/02/2011 17:19	Elizabeth J Marin	1
11184	AK DRO Waters Extraction	AK 102/AK 103	1	111520013A	06/02/2011 03:00	Sherry L Morrow	1
		04/08/02					
00228	Sulfate	EPA 300.0	1	11152196901C	06/02/2011 09:46	Ashley M Adams	5
00220	Nitrate Nitrogen	EPA 353.2	1	11152106102A	06/01/2011 20:41	Venia B McFadden	1
00219	Nitrite Nitrogen	EPA 353.2	1	11148105101A	05/28/2011 12:11	Joseph E McFadden	1
00202	Alkalinity to pH 4.5	SM20 2320 B	1	11153020201B	06/02/2011 06:26	Susan A Engle	1
00201	Alkalinity to pH 8.3	SM20 2320 B	1	11153020201B	06/02/2011 06:26	Susan A Engle	1



Sample Description: DUP-1-052611 Grab Water Sample
 Facility# 82307
 9203 Cessna Dr - Juneau, AK

LLI Sample # WW 6301431
 LLI Group # 1249125
 Account # 10880

Project Name: 82307

Collected: 05/26/2011 by DE

ChevronTexaco

6001 Bollinger Canyon Rd L4310
 San Ramon CA 94583

Submitted: 05/28/2011 09:00

Reported: 06/07/2011 10:11

CJDU1 SDG#: AKE89-02FD

CAT No.	Analysis Name	CAS Number	As Received Result	As Received Method Detection Limit*	As Received Limit of Quantitation	Dilution Factor
GC Volatiles	AK 101		mg/l	mg/l	mg/l	
01438	TPH-GRO AK water C6-C10	n.a.	0.56	0.010	0.10	1
GC Extractable TPH	AK 102/AK 103		mg/l	mg/l	mg/l	
	04/08/02					
01741	TPH-DRO AK water C10-C25	n.a.	24	2.4	12	50

General Sample Comments

State of Alaska Lab Certification No. UST-061

All QC is compliant unless otherwise noted. Please refer to the Quality Control Summary for overall QC performance data and associated samples.

Laboratory Sample Analysis Record

CAT No.	Analysis Name	Method	Trial#	Batch#	Analysis Date and Time	Analyst	Dilution Factor
01438	TPH-GRO AK water C6-C10	AK 101	1	11153A53A	06/03/2011 00:31	Carrie E Miller	1
01146	GC VOA Water Prep	SW-846 5030B	1	11153A53A	06/03/2011 00:31	Carrie E Miller	1
01741	TPH-DRO AK water C10-C25	AK 102/AK 103	1	111520013A	06/03/2011 22:37	Heather E Williams	50
		04/08/02					
11184	AK DRO Waters Extraction	AK 102/AK 103	1	111520013A	06/02/2011 03:00	Sherry L Morrow	1
		04/08/02					

AKE89-02FD



Sample Description: Trip_Blank-1-052611 Water Sample
Facility# 82307
 9203 Cessna Dr - Juneau, AK

LLI Sample # WW 6301432
LLI Group # 1249125
Account # 10880

Project Name: 82307

Collected: 05/26/2011

ChevronTexaco

Submitted: 05/28/2011 09:00

6001 Bollinger Canyon Rd L4310
 San Ramon CA 94583

Reported: 06/07/2011 10:11

CJTBI SDG#: AKE89-03TB*

CAT No.	Analysis Name	CAS Number	As Received Result	As Received Method Detection Limit*	As Received Limit of Quantitation	Dilution Factor
GC Volatiles	AK 101		mg/l	mg/l	mg/l	
01438	TPH-GRO AK water C6-C10	n.a.	N.D.	0.010	0.10	1

General Sample Comments

State of Alaska Lab Certification No. UST-061

All QC is compliant unless otherwise noted. Please refer to the Quality Control Summary for overall QC performance data and associated samples.

Laboratory Sample Analysis Record

CAT No.	Analysis Name	Method	Trial#	Batch#	Analysis Date and Time	Analyst	Dilution Factor
01438	TPH-GRO AK water C6-C10	AK 101	1	11153A53A	06/02/2011 18:42	Carrie E Miller	1
01146	GC VOA Water Prep	SW-846 5030B	1	11153A53A	06/02/2011 18:42	Carrie E Miller	1

AKE89 8011

Volatiles by GC-GRO

Case Narrative Conformance/ Non-Conformance Summary



CLIENT: ChevronTexaco
SDG: AKE89

Volatiles by GC

Fraction: Volatiles by GC-GRO

TPH-GRO AK water C6-C10

<u>Sample #</u>	<u>Client ID</u>	<u>Matrix</u>		<u>Comments</u>
		<u>Liquid</u>	<u>Solid</u>	
6301430	MW-4-052611 Grab Water Sample	X		
6301431	DUP-1-052611 Grab Water Sample	X		Field Duplicate Sample
6301432	Trip_Blank-1-052611 Water Sample	X		Trip Blank

See QC Reference List for Associated Batch QC Samples

SAMPLE PREPARATION:

No problems were encountered with the preparation of the samples.

ANALYSIS:

There were no dilutions performed for analyses associated with samples in this SDG.

No problems were encountered with the analysis of the samples.

QUALITY CONTROL AND NONCONFORMANCE SUMMARY:

Site specific matrix QC samples were not submitted for this SDG. The batch matrix QC was performed on samples from another project. Therefore the matrix effects would not be relevant and matrix QC is not provided in the data package. Laboratory spike data (LCS) are provided.

All QC is within specifications

DATA INTERPRETATION:

No further interpretation is necessary for the data submitted.

Abbreviation Key

UNSPK = Unspiked (for MS/MSD)	LOQ = Limit of Quantitation
MS = Matrix Spike	MDL = Method Detection Limit
MSD = Matrix Spike Duplicate	ND = Not Detected
BKG = Background (for Duplicate)	J = Estimated Value
D = Duplicate (DUP)	E= out of calibration range
LCS = Lab Control Sample	
LCSD = Lab Control Sample Duplicate	* = Out of Specification

AKE89 8814



CLIENT: ChevronTexaco
SDG: AKE89

Volatiles by GC

Fraction: Volatiles by GC-GRO

Nancy E. Saunders
NANCY E. SAUNDERS
Specialist

Narrative Reviewed and Approved 6/22/11 by
(Date)

AKE89 0015



CONFORMANCE/NON-CONFORMANCE SUMMARY

SDG: AKE89

	Indicate Yes, No, N/A
1. Chromatograms labeled / Compounds identified (Field Samples & Method Blanks)	YES
2. Retention times for chromatograms provided	YES
3. Standards summary meet criteria	YES
4. Calibration - Initial calibration performed before sample analysis and continuing calibration performed within 24 hours of sample analysis.	YES
5. Blank contamination If yes, list compounds and concentrations in each blank: N/A	NO
6. Surrogate recoveries meet criteria If not met, list those compounds and the recoveries that fall outside the acceptable range: N/A If not met, were the calculations checked and the results qualified as "estimated"? N/A	YES
7. Matrix Spike / Matrix Spike Duplicate recoveries meet criteria. If not met, list those compounds and the recoveries that fall outside the acceptable range:	N/A
8. Laboratory Control Sample / Laboratory Control Sample Duplicate meet criteria. If not met, list those compounds and the recoveries that fall outside the acceptable range:	YES
9. Retention time summaries for primary and confirmation analyses meet criteria	N/A
10. Were samples run on dissimilar columns?	N/A
11. Extraction holding time met If not met, list number of days exceeded for each sample: N/A	N/A
12. Analysis holding time met If not met, list number of days exceeded for each sample: N/A	YES

Additional Comments:

Summary reviewed and approved by:

Nancy E. Saunders
Dana Kauffman/Manager Data Deliverables

6/22/11
Date

AKE89 8816

Quality Control and QC Summary Forms



Quality Control Reference List
Volatiles by GC

CLIENT: ChevronTexaco
SDG: AKE89

Fraction: Volatiles by GC-GRO

Analysis	Batch Number	Sample Number	Analysis Date
TPH-GRO AK water C6-C10	11153A53A	BLKFI	06/02/2011 15:09:00
		LCSE5	06/02/2011 16:02:00
		6301430	06/02/2011 20:03:00
		6301431	06/03/2011 00:31:00
		6301432	06/02/2011 18:42:00

AKE89 8018

Fraction: Volatiles by GC-GRO

11153A53 / BLKFI					
Analyte	Analysis Date	Blank Results	Units	MDL	LOQ
TPH-GRO AK water C6-C10	06/02/11	N.D.	ug/l	10	100

AKE89 8819

Fraction: Volatiles by GC-GRO

Sample	Trifluorotoluene-F	
	Spike Added	30 ug/l
	% Recovery	Limits
6301430	69	60 - 120
6301431	72	60 - 120
6301432	68	60 - 120
BLKFI	69	60 - 120
LCSE5	88	60 - 120

AKE89 0820

SDG: AKE89
Matrix: LIQUID
Volatiles by GC
Fraction: Volatiles by GC-GRO

LCS: LCSE5		Batch: 11153A53A (Sample number(s): 6301430-6301432)						
Analyte	Spike Added ug/l	LCS Conc ug/l	LCSD Conc ug/l	LCS %Rec	LCSD %Rec	%Rec Limits	%RPD	%RPD Limits
TPH-GRO AK water C6-C10	1100	1200		109		60-120		

AKE89 0021

6D

INITIAL CALIBRATION - RETENTION TIME SUMMARY

Lab Name: Lancaster Laboratories

Contract:

Lab Code:

Case No.:

SAS No.:

SDG No.:

Instrument: 10995FCalibration File: ALK53105GC Column (1): J&W DB-VRX ID: 75 (mm)

Update File:

Date(s) Analyzed: 4/16/2011 4/16/2011

COMPOUND	RT OF STANDARDS							MIDPOINT RT	RT WINDOW	
	LEVEL 1	LEVEL 2	LEVEL 3	LEVEL 4	LEVEL 5	LEVEL 6	LEVEL 7		FROM	TO
Trifluorotoluene-F	3.84	3.84	3.84	3.84	3.84	3.84	3.84	3.84	3.81	3.88
1-Chloro-3-fluorobenzene	5.11	5.11	5.11	5.11	5.11	5.11	5.11	5.11	5.08	5.15

AKES9 8822

6E

INITIAL CALIBRATION - CALIBRATION FACTOR SUMMARY

Lab Name: Lancaster Laboratories

Contract:

Lab Code:

Case No.:

SAS No.:

SDG No.:

Instrument: 10995FCalibration File: ALK53105GC Column (1) : J&W DB-VRX ID: 75 (mm)Date(s) Analyzed: 4/16/2011 4/16/2011

COMPOUND	CALIBRATION FACTORS							MEAN	%RSD
	LEVEL 1	LEVEL 2	LEVEL 3	LEVEL 4	LEVEL 5	LEVEL 6	LEVEL 7		
Trifluorotoluene-F	2.11E+04	2.13E+04	1.99E+04	1.85E+04	1.97E+04			2.01E+04	5.6
1-Chloro-3-fluorobenzene	1.32E+04	1.22E+04	1.28E+04	1.73E+04	2.15E+04	2.72E+04	3.53E+04	1.99E+04	43.7

Average % RSD: 24.7

ANES9 0023

6F

INITIAL CALIBRATION OF MULTICOMPONENT ANALYTES

Lab Name: Lancaster Laboratories

Contract:

Lab Code:

Case No.:

SAS No.:

SDG No.:

Instrument: 10995FCalibration File: ALK53105GC Column (1): J&W DB-VRX ID: 75 (mm)Date(s) Analyzed: 4/16/2011 4/16/2011

COMPOUND	PEAK	RT	RT WINDOW		CALIBRATION FACTOR	AVERAGE CF	LEVEL	AMOUNT	PEAK AREA	%RSD
			FROM	TO						
GRO	1		2.52	6.50	18332	15960	1	21.4	392315	10.3
					18068		2	53	957511	
					16109		3	107	1723709	
					15481		4	536	8297848	
					15031		5	1072	16113350	
					14303		6	2680	38332888	
					14398		7	5380	77173928	

MDJ2001
4-20-11

AKB89 8824

Chrom Perfect Calibration File

File Name: I:\Cal\53\ALK53105.cal
Version: 17

Creator: LCP/895
Description: ALASKA
Reason for change:

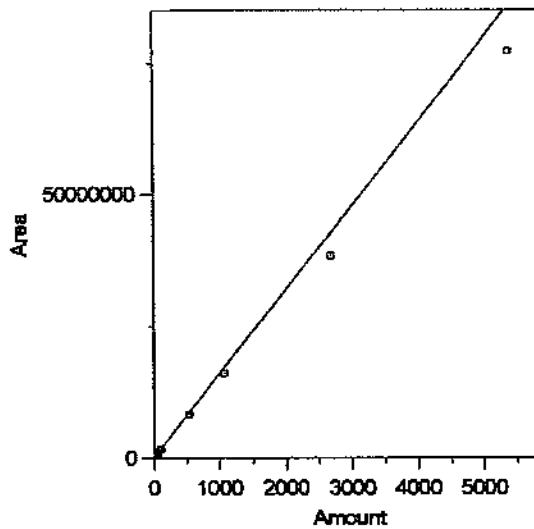
External standard calibration

Standard injection volume: 1
Standard sample weight: 1
Area reject threshold: 0
Reference peak area reject threshold: 0
Amount units: PPB
No default component

Method of calculating data point averages: Current update equal to cal data
Print calibration update report

All levels are normal data points.

I GRO



Expected retention time: 2 minutes
 Search window: 0 minutes
 No retention time reference component
 Group number: 3
 GRO
 High alarm limit: 0
 Low alarm limit: 0
 Component constant: 530

Single peak quantification by area

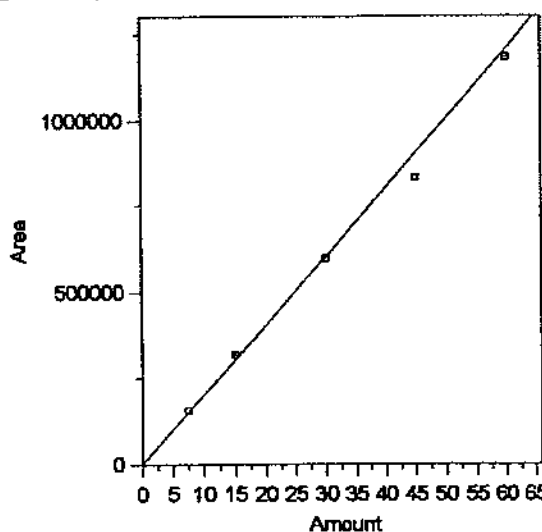
$$Y = 15960.25 X + 0$$

Average CF fit with equal weighting, forced to origin

Coefficient of determination: 0.9812068
 Average error: 8.284%
 Average CF: 15960.25
 RSD: 10.347%

Level	Amount	Response	Cal Factor	Error, %	Source	Date and time
1	21.4	392315	18332.48	14.863	Manual	4/20/2011 4:17:20 PM
2	53	957511	18066.25	13.195	Manual	4/20/2011 4:17:28 PM
3	107	1723709	16109.43	0.935	Manual	4/20/2011 4:17:34 PM
4	536	8297846	15481.06	-3.002	Manual	4/20/2011 4:17:41 PM
5	1072	1.611335E+07	15031.11	-5.822	Manual	4/20/2011 4:17:45 PM
6	2680	3.833289E+07	14303.32	-10.382	Manual	4/20/2011 4:17:52 PM
7	5360	7.717393E+07	14398.12	-9.788	Manual	4/20/2011 4:18:01 PM

2 SURR-TFT-F



Expected retention time: 3.843 minutes
 Search window: 0.035 minutes
 No retention time reference component
 Group number: 1
 SURR-TFT-F
 High alarm limit: 0
 Low alarm limit: 0
 Component constant: 30

Single peak quantification by area

$$Y = 20104.43 X + 0$$

Average CF fit with equal weighting, forced to origin
 Coefficient of determination: 0.990984
 Average error: 4.254%
 Average CF: 20104.43
 RSD: 5.567%

Level	Amount	Response	Cal Factor	Error, %	Source	Date and time
1	7.5	158219.3	21095.91	4.932	Y:\Active\CP53\53105B.0044.BND	4/20/2011 4:14:51 PM
2	15	318767.8	21251.19	5.704	Y:\Active\CP53\53105B.0045.BND	4/20/2011 4:14:54 PM
3	30	597895.8	19929.86	-0.868	Y:\Active\CP53\53105B.0046.BND	4/20/2011 4:14:57 PM
4	45	832875.6	18508.35	-7.939	Y:\Active\CP53\53105B.0047.BND	4/20/2011 4:15:00 PM
5	60	1184212	19736.87	-1.828	Y:\Active\CP53\53105B.0048.BND	4/20/2011 4:15:04 PM
6	(-1)	(0)	-	-	C:\Dep\25\54144B.0012.raw	6/20/2005 4:19:15 PM
7	(-1)	(0)	-	-	C:\Dep\25\54144B.0014.raw	6/20/2005 4:19:19 PM

AKB9 8827

7E

CALIBRATION VERIFICATION SUMMARY

Lab Name: Lancaster Laboratories

Contract:

Lab Code:

Case No.:

SAS No.:

SDG No.:

Instrument: 10995F

Detector: FID

Init. Calib Date(s): 04/16/11

04/16/11

GC Column (1) : J&W DB-VRX ID: 75 (mm)

Date Analyzed: 04/16/11

Lab File ID: 53105B.0054.RAW

Time Analyzed: 16:59

Lab Standard ID: GICVXCW

Initial Calibration: ALK53105

Method: ALASKA

COMPOUND	RT	RT WINDOW FROM TO	CALC AMOUNT	NOM AMOUNT	%D	Limits
GRO		2.52 6.50	953.92	1100.00	-13.3	-25 to +25

Average of %D: 13.3

ARE89 8828

7E

CALIBRATION VERIFICATION SUMMARY

Lab Name: Lancaster Laboratories

Contract:

Lab Code:

Case No.:

SAS No.:

SDG No.:

Instrument: 10995F

Detector: FID

Init. Calib Date(s): 04/16/11

04/16/11

GC Column (1): J&W DB-VRX ID: 75 (mm)

Date Analyzed: 06/02/11

Lab File ID: 53152B.0042.RAW

Time Analyzed: 12:29

Lab Standard ID: WGCCCXXV

Initial Calibration: ALK53105

Method: ALASKA

COMPOUND	RT	RT WINDOW FROM TO		CALC AMOUNT	NOM AMOUNT	%D	Limits
Trifluorotoluene-F	3.85	3.81	3.88	23.49	30.00	-21.7	-43 to +46
GRO		2.52	6.50	577.14	536	7.7	-25 to +25

Average of %D: 14.7

AKES9 8829

7E

CALIBRATION VERIFICATION SUMMARY

Lab Name: Lancaster Laboratories

Contract

Lab Code:

Case No.:

SAS No.:

SDG No.:

Instrument: 10995F

Detector: FID

Init. Calib Date(s): 06/02/11

06/03/11

GC Column (1): J&W DB-VRX ID: 75 (mm)

Date Analyzed: 06/02/11

Lab File ID: 53153B.0017.RAW

Time Analyzed: 21:23

Lab Standard ID: WGCCCXX

Initial Calibration: ALK53105

Method: ALASKA

COMPOUND	RT	RT WINDOW FROM TO		CALC AMOUNT	NOM AMOUNT	%D	Limits
Trifluorotoluene-F	3.84	3.81	3.88	23.15	30.00	-22.8	-43 to +46
GRO		2.52	6.50	556.97	536.00	3.9	-25 to +25

Average of %D: 13.4

AKES9 0030

7E

CALIBRATION VERIFICATION SUMMARY

Lab Name: Lancaster Laboratories

Contract:

Lab Code:

Case No.:

SAS No.:

SDG No.:

Instrument: 10995F

Detector: FID

Init. Calib Date(s): 06/02/11

06/03/11

GC Column (1): J&W DB-VRX ID: 75 (mm)

Date Analyzed: 06/03/11

Lab File ID: 53153B.0025.RAW

Time Analyzed: 0:58

Lab Standard ID: WGCCCXY

Initial Calibration: ALK53105

Method: ALASKA

COMPOUND	RT	RT WINDOW FROM TO		CALC AMOUNT	NOM AMOUNT	%D	Limits
Trifluorotoluene-F	3.84	3.81	3.88	22.09	30.00	-26.4	-43 to +46
GRO		2.52	6.50	520.24	536.00	-2.9	-25 to +25

Average of %D: 14.7

AKZ89.8531

8D ANALYTICAL SEQUENCE

Sequence: 53105B

Lab Name: Lancaster laboratories

Contract:

Lab Code:

Case No.:

SAS No:

SDG No.:

GC Column: JW DB-VRX

ID: 75

Instrument: 10995F

THIS ANALYTICAL SEQUENCE OF BLANKS, SAMPLES AND STANDARDS IS GIVEN BELOW:

	Sample Code No.	Lab Sample ID	Date Analyzed	Time Analyzed	Calibration File	TFTF
001	AA	GRO MAR	04/16/2011	11:38:00	GX53105	3.84
002	AA	IBLK	04/16/2011	12:04:29	TPH53105	3.84
003	WGRO1AA	WGRO11125I	04/16/2011	12:31:31	ALK53105	3.84
004	WGRO2AA	WGRO21125H	04/16/2011	12:58:16	ALK53105	3.84
005	WGRO3AA	WGRO31125H	04/16/2011	13:25:07	ALK53105	3.84
006	WGRO4AA	WGRO41125H	04/16/2011	13:51:49	ALK53105	3.84
007	WGRO5AA	WGRO51125H	04/16/2011	14:18:29	ALK53105	3.84
008	WGRO6AA	WGRO61125H	04/16/2011	14:45:33	ALK53105	3.84
009	WGRO7AA	WGRO71125G	04/16/2011	15:12:22	ALK53105	3.84
010	AA	IBLK	04/16/2011	15:39:20	TPH53105	3.84
011	AA	IBLK	04/16/2011	16:06:14	TPH53105	3.84
012	AA	IBLK	04/16/2011	16:32:58	TPH53105	3.84
013	GICVXCW	GICVX1125J	04/16/2011	16:59:44	ALK53105	3.85
014	GMDLXFO	GMDLX1125K	04/16/2011	17:26:24	ALK53105	3.84

ICAL Dates
 ALK53105 04/16/2011 - 04/16/2011
 GX53105 04/16/2011 - 04/16/2011
 TPH53105 04/16/2011 - 04/16/2011

ICAL RT QC Limits
 TFTF = Trifluorotoluene-F 3.84 (3.81 - 3.88 Minutes)
 TFTF = Trifluorotoluene-F 3.84 (3.81 - 3.87 Minutes)
 TFTF = Trifluorotoluene-F 3.84 (3.81 - 3.87 Minutes)

8-KE89 0232

8D ANALYTICAL SEQUENCE

Sequence: 53152B

Lab Name: Lancaster laboratories

Contract:

Lab Code:

Case No.:

SAS No:

SDG No.:

GC Column: JW DB-VRX

ID: 75

Instrument: 10995F

THIS ANALYTICAL SEQUENCE OF BLANKS, SAMPLES AND STANDARDS IS GIVEN BELOW:

	Sample Code No.	Lab Sample ID	Date Analyzed	Time Analyzed	Calibration File	TFTF
001	AA	GRO MARKER	06/01/2011	17:37:04	ALK53105	3.86
002	WCCPXEM	WCCPX1125DG	06/01/2011	18:03:41	ALK53105	3.84
003	WGCCXXP	WGCCX1125CO	06/01/2011	18:30:31	ALK53105	3.85
004	BLKFC	BLANKA	06/01/2011	18:57:18	ALK53105	3.85
005	LCSDZ	LCSA	06/01/2011	19:24:29	ALK53105	3.85
006	LCSDEH	LCSDA	06/01/2011	19:51:14	ALK53105	3.85
007	LCSE0	LCSB	06/01/2011	20:18:15	ALK53105	3.85
008	LCSDEI	LCSDB	06/01/2011	20:44:46	ALK53105	3.84
009	SEWT1	6297834	06/01/2011	21:11:42	ALK53105	3.84
010	LOATB	6297852	06/01/2011	21:38:29	ALK53105	3.85
011	SHA09	6297820	06/01/2011	22:05:10	ALK53105	3.85
012	SHA10	6297821	06/01/2011	22:31:51	ALK53105	3.84
013	SHA07	6297822	06/01/2011	22:58:30	ALK53105	3.84
014	SHA02	6297823	06/01/2011	23:25:08	ALK53105	3.85
015	SHA08	6297824	06/01/2011	23:51:49	ALK53105	3.85
016	SHA01	6297825	06/02/2011	00:18:48	ALK53105	3.84
017	SHAFD	6297826	06/02/2011	00:45:21	ALK53105	3.85
018	SEW15	6297828	06/02/2011	01:12:20	ALK53105	3.84
019	WCCPXEN	WCCPX1125DG	06/02/2011	01:39:06	ALK53105	3.85
020	WGCCXXO	WGCCX1125CO	06/02/2011	02:05:52	ALK53105	3.85
021	SEW17	6297831	06/02/2011	02:32:59	ALK53105	3.84
022	SEWPR	6297832	06/02/2011	02:59:34	ALK53105	3.84
023	SEWPO	6297833	06/02/2011	03:26:19	ALK53105	3.84
024	LOA01	6297849	06/02/2011	03:53:16	ALK53105	3.85
025	SEW11	6297829	06/02/2011	04:19:56	ALK53105	3.85
026	SEWFD	6297830	06/02/2011	04:46:31	ALK53105	3.85
027	LOA02	6297850	06/02/2011	05:13:33	ALK53105	3.85
028	LOAFD	6297851	06/02/2011	05:40:23	ALK53105	3.84
029	WGCCXXM	WGCCX1125CO	06/02/2011	06:07:26	ALK53105	3.84
030	IBLK	IBLK	06/02/2011	06:34:07	ALK53105	3.84
031	WCCPXER	WCCPX1125DH	06/02/2011	07:00:52	ALK53105	3.84
032	BLKFF	BLANKB	06/02/2011	07:27:48	ALK53105	3.84
033	SSB01	6303288	06/02/2011	07:54:19	ALK53105	3.84

ICAL Dates

ALK53105

04/16/2011 - 04/16/2011

TFTF = Trifluorotoluene-F

ICAL RT QC Limits

3.84

(3.81 - 3.88 Minutes)

AKES9 8833

FORM VIII PEST

8D
ANALYTICAL SEQUENCE

Sequence: 53152B

Lab Name: Lancaster laboratories

Contract:

Lab Code:

Case No.:

SAS No:

SDG No.:

GC Column: JW DB-VRX

ID: 75

Instrument: 10995F

THIS ANALYTICAL SEQUENCE OF BLANKS, SAMPLES AND STANDARDS IS GIVEN BELOW:

	Sample Code No.	Lab Sample ID	Date Analyzed	Time Analyzed	Calibration File	TFTF
034	13101	6303289	06/02/2011	08:21:05	ALK53105	3.84
035	13001	6303290	06/02/2011	08:48:05	ALK53105	3.85
036	76101	6303291	06/02/2011	09:14:50	ALK53105	3.84
037	NES01	6303292	06/02/2011	09:41:31	[8021]53133	
038	AA	IBLK	06/02/2011	10:42:44	ALK53105	3.85
039	SHA08	6297824	06/02/2011	11:09:32	ALK53105	3.84
040	SEW11	6297829	06/02/2011	11:36:20	ALK53105	3.84
041	LOA01	6297849	06/02/2011	12:03:01	ALK53105	3.85
042	WGCCXXV	WGCCX1125CO	06/02/2011	12:29:42	ALK53105	3.85

ALK53105

ICAL Dates

04/16/2011 - 04/16/2011

TFTF = Trifluorotoluene-F

ICAL RT QC Limits

3.84 (3.81 - 3.88 Minutes)

AKB82 8634

FORM VIII PEST

8D ANALYTICAL SEQUENCE

Sequence: 53153B

Lab Name: Lancaster laboratories

Contract:

Lab Code:

Case No.:

SAS No:

SDG No.:

GC Column: JW DB-VRX

ID: 75

Instrument: 10995F

THIS ANALYTICAL SEQUENCE OF BLANKS, SAMPLES AND STANDARDS IS GIVEN BELOW:

	Sample Code No.	Lab Sample ID	Date Analyzed	Time Analyzed	Calibration File	TFTF
001	AA	GRO MARKER	06/02/2011	14:15:34	ALK53105	3.84
002	WCCPXEU	WCCPX1125DH	06/02/2011	14:42:16	ALK53105	3.85
003	BLKFI	BLANKA	06/02/2011	15:09:18	ALK53105	3.84
004	LCSE4	LCSA	06/02/2011	15:36:07	ALK53105	3.85
005	LCSE5	LCSB	06/02/2011	16:02:49	ALK53105	3.85
006	FTA01	6297938	06/02/2011	16:29:42	ALK53105	3.84
007	FTA01MS	6297939	06/02/2011	16:56:30	ALK53105	3.85
008	FTA01MSD	6297940	06/02/2011	17:22:59	ALK53105	3.84
009	FTA01MS	6297939	06/02/2011	17:49:22	ALK53105	3.84
010	FTA01MSD	6297940	06/02/2011	18:16:02	ALK53105	3.85
011	CJTBI	6301432	06/02/2011	18:42:45	ALK53105	3.85
012	GAATB	6299144	06/02/2011	19:09:34	ALK53105	3.85
013	FTATB	6297944	06/02/2011	19:36:21	ALK53105	3.84
014	CJMWS	6301430	06/02/2011	20:03:15	ALK53105	3.84
015	FTEBA	6297949	06/02/2011	20:30:04	ALK53105	3.84
016	WCCPXEV	WCCPX1125DH	06/02/2011	20:56:40	ALK53105	3.85
017	WGCCXX	WGCCX1125CO	06/02/2011	21:23:27	ALK53105	3.84
018	FTTBA	6297950	06/02/2011	21:50:47	ALK53105	3.85
019	FTA02	6297942	06/02/2011	22:17:40	ALK53105	3.84
020	FTAEB	6297943	06/02/2011	22:44:15	ALK53105	3.84
021	GAAM8	6299141	06/02/2011	23:10:56	ALK53105	3.84
022	GAAD1	6299142	06/02/2011	23:37:29	ALK53105	3.84
023	GAAM9	6299143	06/03/2011	00:04:11	ALK53105	3.84
024	CJDU1	6301431	06/03/2011	00:31:19	ALK53105	3.85
025	WGCCXXY	WGCCX1125CO	06/03/2011	00:58:02	ALK53105	3.84

ICAL Dates

04/16/2011 - 04/16/2011

TFTF = Trifluorotoluene-F

ICAL RT QC Limits

3.84 (3.81 - 3.88 Minutes)

AKBB9 8835

FORM VIII PEST

Sample Data



LOQ/MDL Summary
Volatiles by GC

SDG: AKE89

Fraction: Volatiles by GC-GRO

01438: TPH-GRO AK water C6-C10 Analyte Name	Default MDL	Default LOQ	Units
TPH-GRO AK water C6-C10	10	100	ug/l

AKE89 8637

Lancaster Laboratories-Range Data Summary

Sample Name: 6301430 **CJMWS** **Sample ID:** AA **Batchnumber:** 11153A53A
Sample Amount: 1. **Total Volume:** 1. ml **Analyst:** 1991 **SDG:** AKE89 **State:** AK
Analyses: 01440

Injection Summary

Injected on : 6/2/2011 20:03:15
Instrument : CP53--10995F
Result file : 53153B.0014.RAW
Calibration files : ALK53105.cal
Method files : ALK53105.MET
Setting : ALK53105

Surrogate Recoveries

SURR-TFT-F 68.7% (60-120) **Conc.:** 20.612612
SURR-1C3FB

Range	Retention Times	Area	Amount	LOQ	MDL	Flags	Units
<input type="checkbox"/> SURR-TFT-F	3.84 (3.81 - 3.88)	414405	20.6126				ppb
<input type="checkbox"/> SURR-1C3FB	5.11 (5.08 - 5.15)	563133	28.2668				ppb
<input type="checkbox"/> GRO	2.52 - 6.50	9548232	537.0109	100	10		ppb

Comments: _____

Reviewed by: Carrie Miller

Date: JUN 03 2011

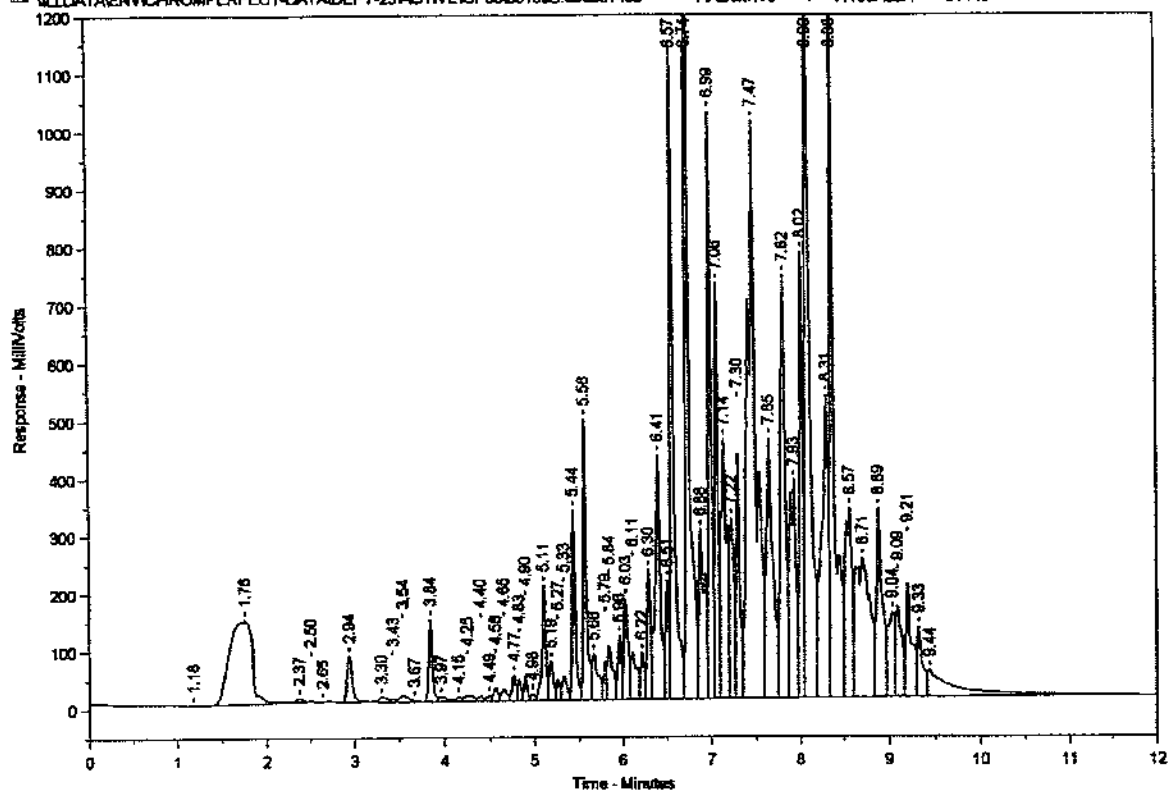
Verified by: M000001

Date: 6-3-11

AKES9 8838

6301430 AACJMW5 T 11153A53A 01440
CP53 FID 10995F 53153B.0014.RAW
Date Acquired: 6/2/2011 8:03:15 PM

\\L\DATA\ENV\CHROMPERFECT-DATA\DEPT-25\ACTIVE\CP53\53153B.001430 AACJMW5 T 11153A53A 01440



6301430 AACJMW5 T 11153A53A 01440
Date Acquired: 6/2/2011 8:03:15 PM Instrument: CP53 10995F
Raw File: 53153B.0014.RAW Units: ug/L
Analyst: Method File: ALK53105.MET
Dilution Factor: 1 Column: 30 M DB-VRX x 0.45mm x 2.55 um

Threshold: 3

Peak Table using calibration : ALK53105.cal- Version 17

Number of Compounds: 3

Component Name	Ret. Time	Exp. Ret Time	Amount ug/L	Peak Area (A)*	Peak Height (H)
SURR-TFT-F	3.84	3.84	20.61	414405	143587.4
SURR-1C3FB	5.11	5.11	28.27	563133	201024.8

RT Start	RT Stop	Unadj GRO	Total Surr.	Adj. GRO
2.52	6.5	9548232	977538	8570694

Surrogate Percent Recovery: 68.70871

Total GRO Area: 8570694.00

Total GRO Concentration: 537.00 ug/L

File: \\L\DATA\ENV\CHROMPERFECT-DATA\DEPT-25\ACTIVE\CP53\53153B.0014.RAW

Lancaster Laboratories-Range Data Summary

Sample Name: 6301431 **CJDU1** **Sample ID:** AA **Batchnumber:** 11153A53A
Sample Amount: 1. **Total Volume:** 1. ml **Analyst:** 1991 **SDG:** AKE89 **State:** AK
Analyses: 01440

Injection Summary

Injected on : 6/3/2011 00:31:19
Instrument : CP53-10995F
Result file : 53153B.0024.RAW
Calibration files : ALK53105.cal
Method files : ALK53105.MET
Setting : ALK53105

Surrogate Recoveries

SURR-TFT-F 72.4% (60-120) **Conc.:** 21.713469
SURR-1C3FB

<u>Range</u>	<u>Retention Times</u>	<u>Area</u>	<u>Amount</u>	<u>LOQ</u>	<u>MDL</u>	<u>Flags</u>	<u>Units</u>
<input type="checkbox"/> SURR-TFT-F	3.85 (3.81 - 3.88)	436537	21.7135				ppb
<input type="checkbox"/> SURR-1C3FB	5.11 (5.08 - 5.15)	578859	29.0562				ppb
<input type="checkbox"/> GRO	2.52 - 8.50	9950828	559.8641	100	10		ppb

Comments: _____

Reviewed by: Carrie Miller

Date: JUN 03 2011

Verified by: MOD001

Date: 6-3-11

AKE89 0040

Chrom Perfect Chromatogram Report

6301431 AACJDU1 T 11153A53A 01440

CP53 FID

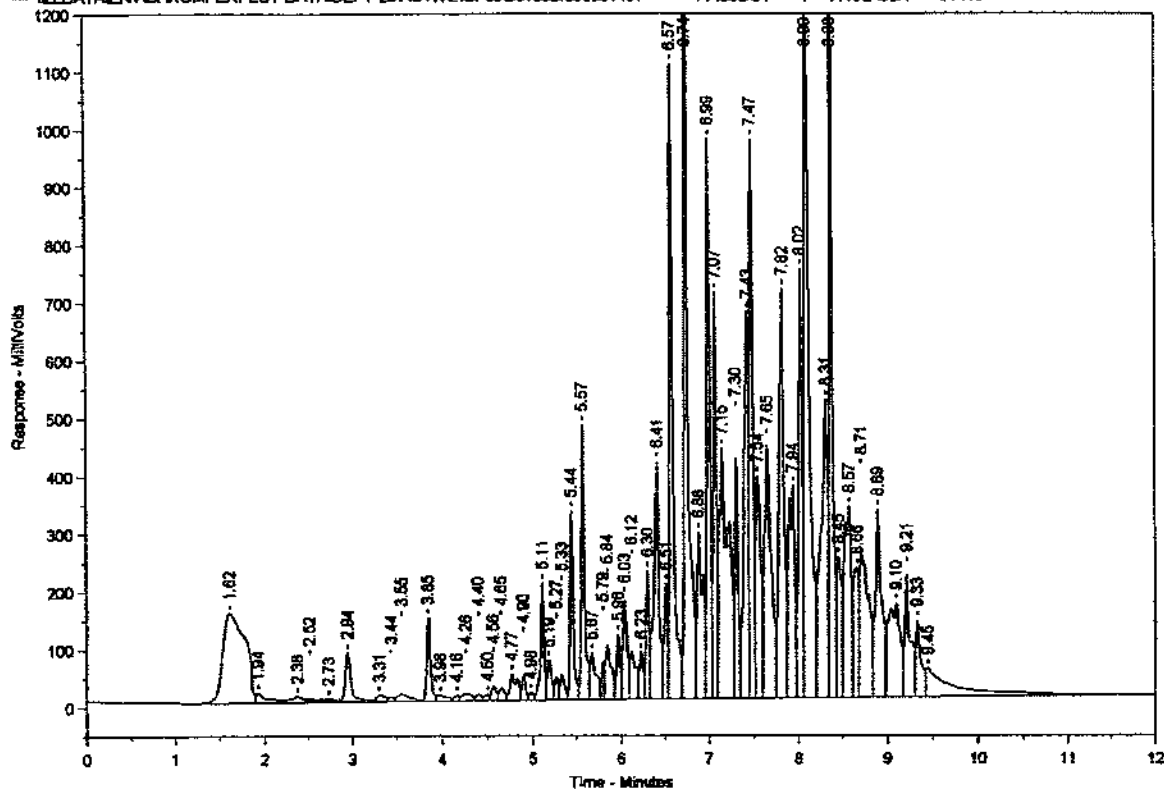
10995F

53153B.0024.RAW

Date Acquired: 6/3/2011 12:31:19 AM

\\L\DATA\ENV\CHROMPERFECT-DATE\DEPT-25\ACTIVE\CP53\53153B.006301431

AACJDU1 T 11153A53A 01440



6301431 AACJDU1 T 11153A53A 01440

Date Acquired: 6/3/2011 12:31:19 AM

Raw File: 53153B.0024.RAW

Analyst:

Dilution Factor: 1

01440

Instrument: CP53 10995F

Units: ug/L

Method File: ALK53105.MET

Column: 30 M DB-VRX x 0.45mm x 2.55 um

Threshold: 3

Peak Table using calibration: ALK53105.cal- Version 17

Number of Compounds: 3

Component Name	Ret. Time	Exp. Ret Time	Amount ug/L	Peak Area (A)*	Peak Height (H)
SURR-TFT-F	3.85	3.84	21.71	436537	145410.4
SURR-1C3FB	5.11	5.11	29.06	578859	203952.9

RT Start	RT Stop	Unadj GRO	Total Surr.	Adj. GRO
2.52	6.5	9950828	1015396	8935432

Surrogate Percent Recovery: 72.37823

Total GRO Area: 8935432.00

Total GRO Concentration: 559.86 ug/L

File: \\L\DATA\ENV\CHROMPERFECT-DATE\DEPT-25\ACTIVE\CP53\53153B.0024.RAW

Lancaster Laboratories-Range Data Summary

Sample Name: 6301432 **CJTBI** **Sample ID:** AA **Batchnumber:** 11153A53A
Sample Amount: 1. **Total Volume:** 1. ml **Analyst:** 1991 **SDG:** AKE89 **State:** AK
Analyses: 01440

Injection Summary

Injected on : 6/2/2011 18:42:45
Instrument : CP53-10995F
Result file : 53153B.0011.RAW
Calibration files : ALK53105.cal
Method files : ALK53105.MET
Setting : ALK53105

Surrogate Recoveries

SURR-TFT-F 68.1% (60-120) **Conc.:** 20.423792
SURR-1C3FB

Range	Retention Times	Area	Amount	LOQ	MDL	Flags	Units
<input type="checkbox"/> SURR-TFT-F	3.85 (3.81 - 3.88)	410809	20.4238				ppb
<input type="checkbox"/> SURR-1C3FB	5.11 (5.08 - 5.15)	414277	20.7949				ppb
<input type="checkbox"/> GRO	2.52 - 6.50	874393	3.1019	<100	<10		ppb

Comments:

Reviewed by:

Carrie E Miller

Date:

JUN 03 2011

Verified by:

M D J 0001

Date:

6-3-11

AKE89 0042

Chrom Perfect Chromatogram Report

6301432 AACJTB1 T 11153A53A 01440

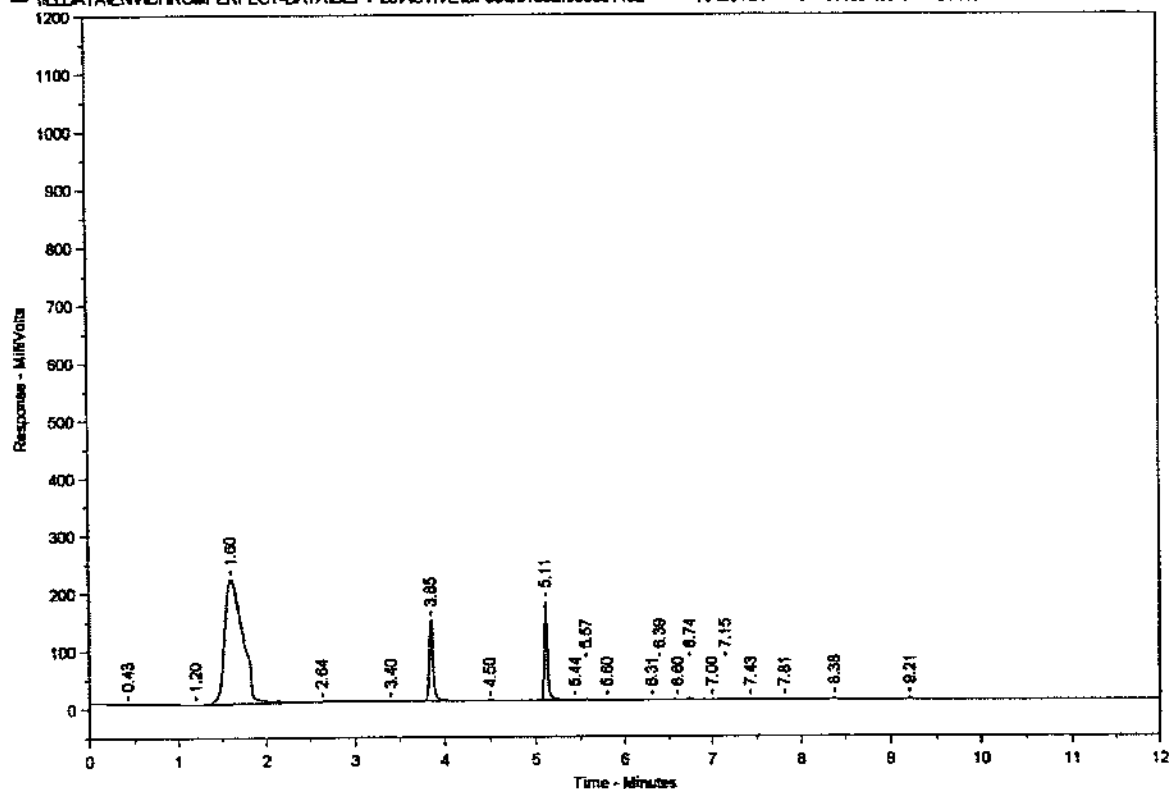
CP53 FID

10995F

53153B.0011.RAW

Date Acquired: 6/2/2011 6:42:45 PM

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6301432 AACJTB1 T 11153A53A 01440

Date Acquired: 6/2/2011 6:42:45 PM

Instrument: CP53 10995F

Raw File: 53153B.0011.RAW

Units: ug/L

Analyst:

Method File: ALK53105.MET

Dilution Factor: 1

Column: 30 M DB-VRX x 0.45mm x 2.55 um

Threshold: 3

Peak Table using calibration: ALK53105.cal- Version 17

Number of Compounds: 3

Component Name	Ret. Time	Exp. Ret Time	Amount ug/L	Peak Area (A)*	Peak Height (H)
SURR-TFT-F	3.85	3.84	20.42	410609	141599.7
SURR-1C3FB	5.11	5.11	20.79	414277	172558.3

RT Start	RT Stop	Unadj GRO	Total Surr.	Adj. GRO
2.52	6.5	874393	824886	49507

Surrogate Percent Recovery: 68.07931

Total GRO Area: 49506.88

Total GRO Concentration: 3.10 ug/L

File: \\LLDATA\EN\CHROMPERFECT-DATA\DEPT-25\ACTIVE\CP53\53153B.0011.RAW

Raw QC Data

Lancaster Laboratories-Range Data Summary

Sample Name: BLANKA **BLKF1** **Sample ID:** AA **Batchnumber:** 11153A53A
Sample Amount: 1. **Total Volume:** 1. ml **Analyst:** 1991 **SDG:** **State:**
Analyses: 01440 02102

Injection Summary

Injected on : 6/2/2011 15:09:18
Instrument : CP53-10995F
Result file : 53153B.0003.RAW
Calibration files : ALK53105.cal
Method files : ALK53105.MET
Setting : ALK53105

Surrogate Recoveries

SURR-TFT-F 68.6% (60-120) **Conc.:** 20.57885
SURR-1C3FB

Range	Retention Times	Area	Amount	LOQ	MDL	Flags	Units
<input type="checkbox"/> SURR-TFT-F	3.84 (3.81 - 3.88)	413726	20.5789				ppb
<input type="checkbox"/> SURR-1C3FB	5.11 (5.08 - 5.15)	417529	20.9581				ppb
<input type="checkbox"/> GRO	2.52 - 6.50	839862	0.5392	<100	<10		ppb

Comments:

Reviewed by:

Date:

cem1991

6/3/11

Verified by: *Carrie Miller*

Date:

JUN 03 2011

MD00001

6-3-11

③ cem1991 6-3-11

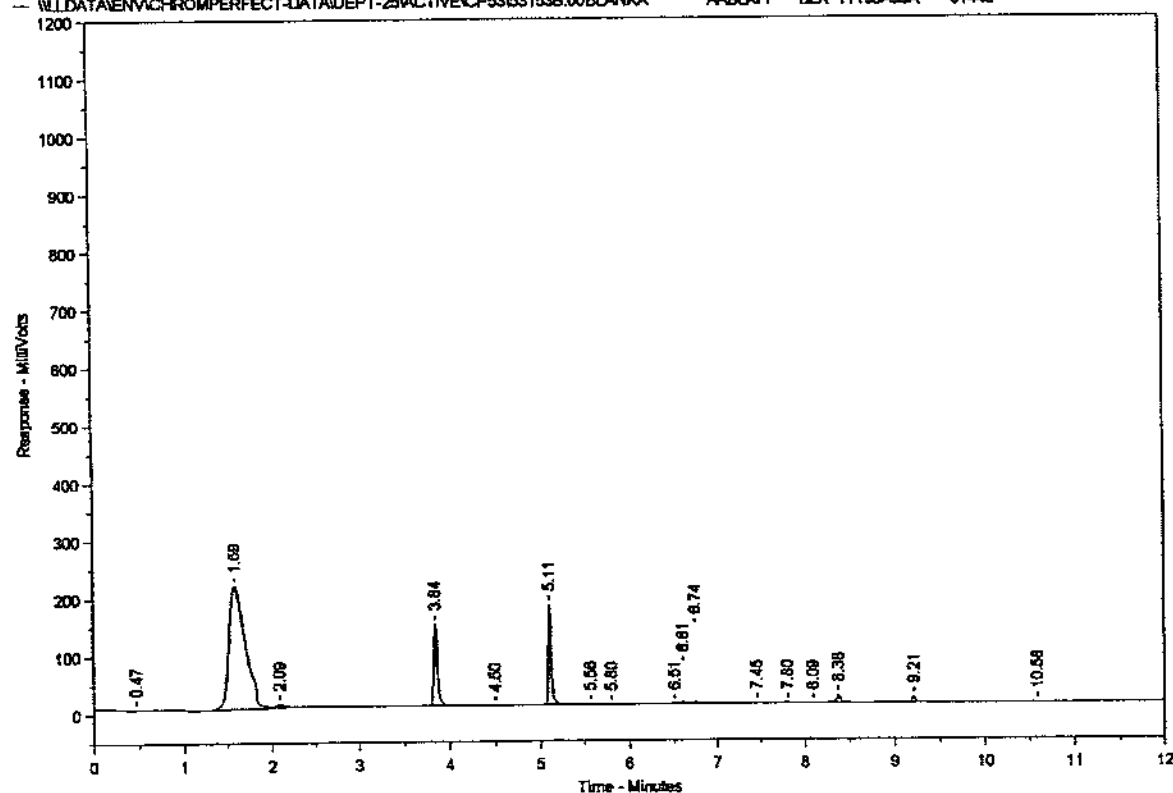
AKZ89 0845

Chrom Perfect Chromatogram Report

BLANKA AABLKFI BLK 11153A53A 01440
CP53 FID 10995F 53153B.0003.RAW

Date Acquired: 6/2/2011 3:09:18 PM

\\L\DATA\ENV\CHROMPERFECT-DATA\DEPT-25\ACTIVE\CP53\53153B.00BLANKA AABLKFI BLK 11153A53A 01440



BLANKA AABLKFI BLK 11153A53A 01440
Date Acquired: 6/2/2011 3:09:18 PM Instrument: CP53 10995F
Raw File: 53153B.0003.RAW Units: ug/L
Analyst: Method File: ALK53105.MET
Dilution Factor: 1 Column: 30 M DB-VRX x 0.45mm x 2.55 um

Threshold: 3

Peak Table using calibration : ALK53105.cal- Version 17

Number of Compounds: 3

Component Name	Ret. Time	Exp. Ret Time	Amount ug/L	Peak Area (A)*	Peak Height (H)
SURR-TFT-F	3.84	3.84	20.58	413728	142423.2
SURR-1C3FB	5.11	5.11	20.96	417529	174302.2

RT Start	RT Stop	Unadj GRO	Total Surr.	Adj. GRO
2.52	6.5	839862	831255	8606

Surrogate Percent Recovery: 68.59617

Total GRO Area: 8606.31

Total GRO Concentration: 0.54 ug/L

File: \\L\DATA\ENV\CHROMPERFECT-DATA\DEPT-25\ACTIVE\CP53\53153B.0003.RAW

CO2 by Headspace Data

Case Narrative Conformance/Non-Conformance Summary



CLIENT: ChevronTexaco
SDG: AKE89

EPH/Miscellaneous GC

Fraction: CO2 by Headspace

CO2 by Headspace

<u>Sample #</u>	<u>Client ID</u>	<u>Matrix</u>		<u>Comments</u>
		<u>Liquid</u>	<u>Solid</u>	
6301430	MW-4-052611 Grab Water Sample	X		
See QC Reference List for Associated Batch QC Samples				

SAMPLE PREPARATION:

No problems were encountered with the preparation of the samples.

ANALYSIS:

There were no dilutions performed for analyses associated with samples in this SDG.

No problems were encountered with the analysis of the samples.

QUALITY CONTROL AND NONCONFORMANCE SUMMARY:

Site specific matrix QC samples were not submitted for this SDG. The batch matrix QC was performed on samples from another project. Therefore the matrix effects would not be relevant and matrix QC is not provided in the data package. Laboratory spike data (LCS) are provided.

All QC is within specification.

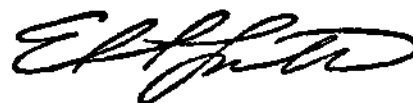
DATA INTERPRETATION:

No further interpretation is necessary for the data submitted.

Abbreviation Key

UNSPK = Unspiked (for MS/MSD)	LOQ = Limit of Quantitation
MS = Matrix Spike	MDL = Method Detection Limit
MSD = Matrix Spike Duplicate	ND = Not Detected
BKG = Background (for Duplicate)	J = Estimated Value
D = Duplicate (DUP)	E = out of calibration range
LCS = Lab Control Sample	
LCSD = Lab Control Sample Duplicate	* = Out of Specification

Narrative Reviewed and Approved 6/28/11 by
(Date)


Elizabeth A. Smith
Specialist

AKE89 0049



CONFORMANCE/NON-CONFORMANCE SUMMARY

SDG: AKE89

Indicate Yes, No, N/A

- | | |
|---|-----|
| 1. Chromatograms labeled / Compounds identified (Field Samples & Method Blanks) | YES |
| 2. Retention times for chromatograms provided | YES |
| 3. Standards summary meet criteria | YES |
| 4. Calibration - Initial calibration performed before sample analysis and continuing calibration performed within 24 hours of sample analysis. | YES |
| 5. Blank contamination
If yes, list compounds and concentrations in each blank: N/A | NO |
| 6. Surrogate recoveries meet criteria
If not met, list those compounds and the recoveries that fall outside the acceptable range: N/A
If not met, were the calculations checked and the results qualified as "estimated"? N/A | YES |
| 7. Matrix Spike / Matrix Spike Duplicate recoveries meet criteria.
If not met, list those compounds and the recoveries that fall outside the acceptable range: | N/A |
| 8. Laboratory Control Sample / Laboratory Control Sample Duplicate meet criteria.
If not met, list those compounds and the recoveries that fall outside the acceptable range: | YES |
| 9. Retention time summaries for primary and confirmation analyses meet criteria | N/A |
| 10. Were samples run on dissimilar columns? | N/A |
| 11. Extraction holding time met
If not met, list number of days exceeded for each sample: N/A | YES |
| 12. Analysis holding time met
If not met, list number of days exceeded for each sample: N/A | YES |

Additional Comments:

Summary reviewed and approved by:


Elizabeth A. Smith

Specialist, Manager Data Deliverables

6/28/11
Date

AKE89. 0050.

Quality Control and Calibration Summary Forms



Quality Control Reference List
EPH/Miscellaneous GC

CLIENT: ChevronTexaco
SDG: AKE89

Fraction: CO2 by Headspace

Analysis
CO2 by Headspace

Batch Number
111520033A

Sample Number
PBLK33152
LCS33152
6301430

Analysis Date
06/02/2011 16:06:00
06/02/2011 16:16:00
06/02/2011 17:19:00

AKE89: 8052

Fraction: CO2 by Headspace

111520033 / PBLK33152 Analyte	Analysis Date	Blank Results	Units	MDL	LOQ
CO2 by Headspace	06/02/11	N.D.	ug/l	4000	12000

~~AKE89 #053~~



Quality Control Summary
Laboratory Control Standard (LCS)
Laboratory Control Standard Duplicate(LCSD)

SDG: AKE89
Matrix: LIQUID

EPH/Miscellaneous GC
Fraction: CO2 by Headspace

LCS: LCS33152	Batch: 111520033A (Sample number(s): 6301430)							
Analyte	Spike Added ug/l	LCS Conc ug/l	LCSD Conc ug/l	LCS %Rec	LCSD %Rec	%Rec Limits	%RPD	%RPD Limits
CO2 by Headspace	36000	34000		94		67-124		

AKE89. 0854

6D

INITIAL CALIBRATION - RETENTION TIME SUMMARY

Lab Name: Lancaster Laboratories

Contract:

Lab Code:

Case No.:

SAS No.:

SDG No.:

Instrument: H3145ACalibration File: 1C20144GC Column (1): CTR1COLUMN ID: 1 (mm)

Update File:

Date(s) Analyzed: 5/24/2011 5/24/2011

COMPOUND	RT OF STANDARDS					MIDPOINT	RT WINDOW	
	LEVEL 1	LEVEL 2	LEVEL 3	LEVEL 4	LEVEL 5	RT	FROM	TO
CARBON DIOXIDE	1.36	1.37	1.38	1.36	1.34	1.37	1.27	1.47

AKES9 0055

6E

INITIAL CALIBRATION - CALIBRATION FACTOR SUMMARY

Lab Name: Lancaster Laboratories

Contract:

Lab Code:

Case No.:

SAS No.:

SDG No.:

Instrument: H3145ACalibration File: 1C20144GC Column (1) : CTR1COLUMN ID: 1 (mm)Date(s) Analyzed: 5/24/2011 5/24/2011

COMPOUND	CALIBRATION FACTORS						%RSD
	LEVEL 1	LEVEL 2	LEVEL 3	LEVEL 4	LEVEL 5	MEAN	
CARBON DIOXIDE	3.98E-01	4.36E-01	4.29E-01	4.84E-01	4.79E-01	4.48E-01	8.1

Average % RSD: 8.1

~~8859 8856~~

6D

INITIAL CALIBRATION - RETENTION TIME SUMMARY

Lab Name: Lancaster Laboratories

Contract:

Lab Code:

Case No.:

SAS No.:

SDG No.:

Instrument: H3145ACalibration File: 2C20144GC Column (1): CTR1COLUMN ID: 1 (mm)Update File: 2CC20144.0003.RAWDate(s) Analyzed: 6/2/2011 6/2/2011

COMPOUND	RT OF STANDARDS					MIDPOINT	RT WINDOW	
	LEVEL 1	LEVEL 2	LEVEL 3	LEVEL 4	LEVEL 5	RT	FROM	TO
CARBON DIOXIDE			1.37			1.37	1.27	1.47

ARE89 0058

6E

INITIAL CALIBRATION - CALIBRATION FACTOR SUMMARY

Lab Name: Lancaster Laboratories

Contract:

Lab Code:

Case No.:

SAS No.:

SDG No.:

Instrument: H3145A

Calibration File: 2C20144

GC Column (1): CTR1COLUMN ID: 1 (mm)

Date(s) Analyzed: 6/2/2011 6/2/2011

COMPOUND	CALIBRATION FACTORS					MEAN	%RSD
	LEVEL 1	LEVEL 2	LEVEL 3	LEVEL 4	LEVEL 5		
CARBON DIOXIDE	3.98E-01	4.36E-01	4.29E-01	4.84E-01	4.79E-01	4.46E-01	8.1

Average % RSD: 8.1

AXES 8859

File Name: Y:\CP20\2C20144.CAL
Version: 1

Creator:
Description:
Reason for change:

*updated ICAK name
to be different from ICAK
in dept. 24*

External standard calibration
Standard injection volume: 1
Standard sample weight: 1
Area reject threshold: 0
Reference peak area reject threshold: 0
Amount units: PPB
No default component

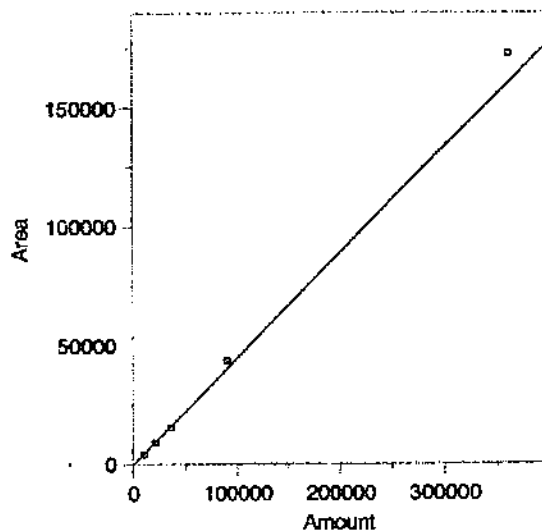
Method of calculating data point averages: Equal weight for all updates
No calibration update report

All levels are normal data points.

✓ TW 6/3/11

EM 6/3/11

1 CARBON DIOXIDE



Expected retention time (frozen): 1.37 minutes
Search window: 0.1 minutes
No retention time reference component
Group number: 0
High alarm limit: 0
Low alarm limit: 0
Component constant: 0
Single peak quantification by area
 $Y = 0.4455497 X + 0$
Average CF fit with equal weighting, forced to origin
Coefficient of determination: 0.9918959
Average error: 6.514%
Average CF: 0.4455497
RSD: 8.101%

Level	Amount	Response	Cal Factor	Error, %	Source	Date and time
1	10800	4303.074	0.3984328	-10.575	\\Uslan-chromperf\Active-data\CP20\1C20144.0005.BND	5/24/2011 3:43:19 PM
2	21600	9425	0.4363426	-2.066	\\Uslan-chromperf\Active-data\CP20\1C20144.0004.BND	5/24/2011 3:53:48 PM
3	36000	15455.46	0.4293183	-3.643	\\Uslan-chromperf\Active-data\CP20\1C20144.0005.BND	5/24/2011 4:04:58 PM
4	90000	43573.96	0.4841551	8.665	\\Uslan-chromperf\Active-data\CP20\1C20144.0006.BND	5/24/2011 4:15:29 PM
5	360000	172619.8	0.4794994	7.620	\\Uslan-chromperf\Active-data\CP20\1C20144.0007.BND	5/24/2011 4:26:03 PM

7E

CALIBRATION VERIFICATION SUMMARY

Lab Name: Lancaster Laboratories

Contract:

Lab Code:

Case No.:

SAS No.:

SDG No.:

Instrument: H3145A

Init. Calib Date(s): 05/24/11

05/24/11

GC Column (1) : CTR1COLUMN ID: 1 (mm)

Date Analyzed: 05/24/11

Lab File ID: 1C20144.0010.RAW

Time Analyzed: 17:19

Lab Standard ID: CO2CXPf

Initial Calibration: 1C20144

COMPOUND	RT	RT WINDOW FROM TO		CALC AMOUNT (ug/l)	NOM AMOUNT (ug/l)	%D
CARBON DIOXIDE	1.36	1.27	1.47	37775.80	35008.00	7.9

Average of %D: 7.9

AKKES 6861

7E

CALIBRATION VERIFICATION SUMMARY

Lab Name: Lancaster Laboratories

Contract:

Lab Code:

Case No.:

SAS No.:

SDG No.:

Instrument: H3145A

Init. Calib Date(s): 06/02/11

06/02/11

GC Column (1) : CTR1COLUMN ID: 1 (mm)

Date Analyzed: 06/02/11

Lab File ID: 2CC20144.0003.RAW

Time Analyzed: 15:56

Lab Standard ID: CO2_3LG

Initial Calibration: 2C20144

COMPOUND	RT	RT WINDOW FROM TO		CALC AMOUNT (ug/l)	NOM AMOUNT (ug/l)	%D
CARBON DIOXIDE	1.37	1.27	1.47	35160.22	36000.00	-2.3

Average of %D: 2.3

AKED9 5862

7E

CALIBRATION VERIFICATION SUMMARY

Lab Name: Lancaster Laboratories

Contract:

Lab Code:

Case No.:

SAS No.:

SDG No.:

Instrument: H3145A

Init. Calib Date(s): 06/02/11

06/02/11

GC Column (1): CTR1COLUMN ID: 1 (mm)

Date Analyzed: 06/02/11

Lab File ID: 2CC20144.0012.RAW

Time Analyzed: 17:30

Lab Standard ID: CO2_3LH

Initial Calibration: 2C20144

COMPOUND	RT	RT WINDOW FROM TO		CALC AMOUNT (ug/l)	NOM AMOUNT (ug/l)	%D
CARBON DIOXIDE	1.37	1.27	1.47	34869.15	36000.00	-3.1

Average of %D: 3.1

AKES9 8863

8D
ANALYTICAL SEQUENCE

Sequence: 1C20144

Lab Name: Lancaster laboratories

Contract:

Lab Code:

Case No.:

SAS No:

SDG No.:

GC Column: CTR1COLUMN

ID: 1

Instrument: H3145A

THIS ANALYTICAL SEQUENCE OF BLANKS, SAMPLES AND STANDARDS IS GIVEN BELOW:

	Sample Code No.	Lab Sample ID	Date Analyzed	Time Analyzed	Calibration File
001		CONDITIONER	05/24/2011	15:21:26	1C20144
002		CONDITIONER	05/24/2011	15:31:14	1C20144
003	CO2_1AA	CO2_11132B	05/24/2011	15:41:12	1C20144
004	CO2_2AA	CO2_21132B	05/24/2011	15:51:40	1C20144
005	CO2_3AA	CO2_31132K	05/24/2011	16:02:56	1C20144
006	CO2_4AA	CO2_41132M	05/24/2011	16:13:28	1C20144
007	CO2_5AA	CO2_51132B	05/24/2011	16:23:58	1C20144
008	AA	IBLK	05/24/2011	16:35:32	1C20144
009	CO2MXBQ	CO2MX1132B	05/24/2011	17:09:02	1C20144
010	CO2CXPF	CO2CX1132B	05/24/2011	17:19:14	1C20144

AKB9 0024

8D
ANALYTICAL SEQUENCE

Sequence: 2CC20144

Lab Name: Lancaster laboratories

Contract:

Lab Code:

Case No.:

SAS No:

SDG No.:

GC Column: CTR1COLUMN

ID: 1

Instrument: H3145A

THIS ANALYTICAL SEQUENCE OF BLANKS, SAMPLES AND STANDARDS IS GIVEN BELOW:

	Sample Code No.	Lab Sample ID	Date Analyzed	Time Analyzed	Calibration File
001		CONDITIONER	06/02/2011	15:36:16	2C20144
002		CONDITIONER	06/02/2011	15:46:24	2C20144
003	CO2_3LG	CO2_31132L	06/02/2011	15:56:19	2C20144
004	PBLK33152	BLANKA	06/02/2011	16:06:24	2C20144
005	LCS33152	LCSA	06/02/2011	16:16:56	2C20144
006	6GW05	6296060	06/02/2011	16:27:23	2C20144
007	6GW05MS	6296060	06/02/2011	16:37:50	2C20144
008	6GW05MSD	6296060	06/02/2011	16:48:19	2C20144
009	3AW05	6296061	06/02/2011	16:58:44	2C20144
010	4RR05	6296062	06/02/2011	17:09:16	2C20144
011	CJMW5	6301430	06/02/2011	17:19:42	2C20144
012	CO2_3LH	CO2_31132L	06/02/2011	17:30:14	2C20144

Sample Data

SDG: AKE89**Fraction: CO2 by Headspace**

08097: CO2 by Headspace Analyte Name	Default MDL	Default LOQ	Units
CO2 by Headspace	4,000	12,000	ug/l

AKE89: 0067

Lancaster Laboratories-Single Component Data Summary

Sample Name: 6301430 **CJMW5** **Sample ID:** AA **Batchnumber:** 111520033A
Sample Amount: 5 **ML** **Total Volume:** 5 **ml** **Analyst:** 2343 **SDG:** AKE89 **State:** AK
Analyses: 08097

Analysis Report (A)

Injected on : JUN 02, 2011 17:19:42
 Instrument : CP20-H3145A
 Result file : 2CC20144.0011.RAW
 Calibration file : 2C20144.CAL
 Method file : CO2.MET

Peak name	Min	R.T.	Max	Area	Amount
CARBON DIOXIDE	1.27	1.36	1.47	39657	89007.568360

Summary Report

Compound Name	Column	Amount Found	LOQ	MDL	Qualifiers	%Difference	Comments
<input checked="" type="checkbox"/> CARBON DIOXIDE	A	89007.568360	12000	4000			

Units: ug/l

Tracy A. Cole

Reviewed by: EM 73613

Verified by: JUN 03 2011

Date: 6/2/11

Date: _____

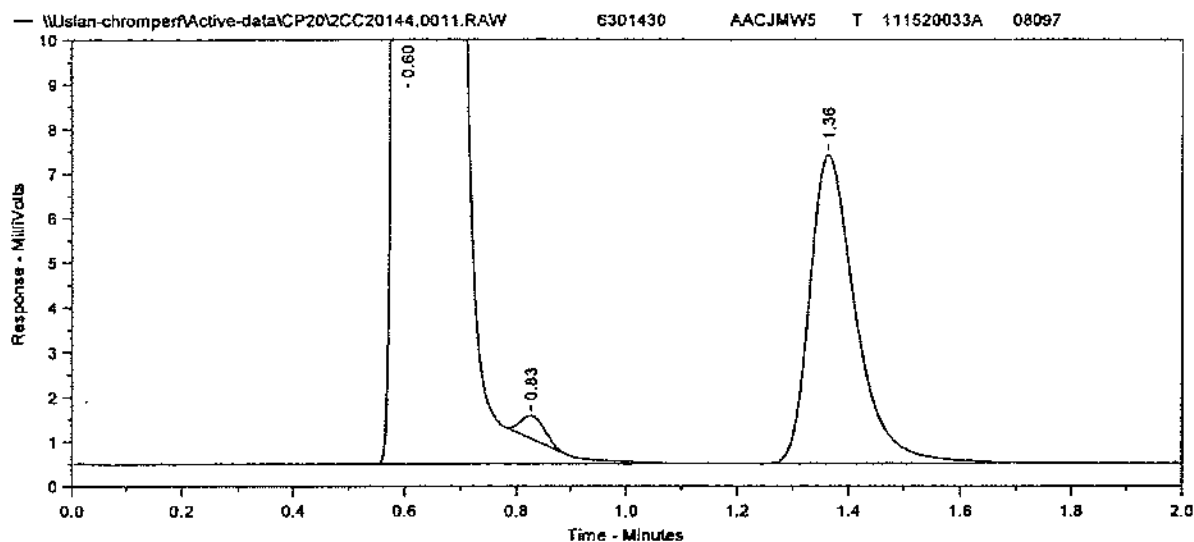
Tracy A. Cole
 Senior Specialist

AKE89 8868

* Recovery outside QC Limits

Printed on: 6/2/2011 17:50:05

Chrom Perfect Chromatogram Report



Sample Name = 6301430 AACJMW5 T 111520033A 08097

Instrument = CP20
Detector = H3145A

Raw File Name = \\Usian-chromperf\Active-data\CP20\2CC20144.0011.RAW

Method File Name = \\USLAN-CHROMPERF\ACTIVE-DATA\CP20\CO2.MET

Calibration File Name = Y:\CP20\2CC20144.CAL

Time Injected = 6/2/2011 5:19:42 PM

Run Time = 2

Dilution Factor = 5

Amount Injected = 1

Peak Threshold = 0

Peak Width = 0.05

Operator = 2343

Incubate 5 minutes at 35C; 40C for .1 minute; 40C/min to 80C; Hold 1 min
Column: CTR 1 COLUMN; 6' X 1/4" X .210 STAINLESS STEEL

	Ret. Time	Amount	Amt %	Area
CARBON DIOXIDE	1.36	89007.57	100.000	39657

AK889 8869

Raw QC Data

Lancaster Laboratories Single Component Data Summary

Sample Name: BLANKA 6/2/11 PBLK33152 Sample ID: AA Batch number: 111520033A
Sample Amount: 5 ML Total Volume: 5 ml Analyst: 2343 SDG: State:
Analyses: 08097

Analysis Report (A)

Injected on : JUN 02, 2011 16:06:24
Instrument : CP20-H3145A
Result file : 2CC20144.0004.RAW
Calibration file : 2C20144.CAL
Method file : CO2.MET

Summary Report

Compound Name	Column	Amount Found	LOQ	MDL	Qualifiers	%Difference	Comments
<input checked="" type="checkbox"/> CARBON DIOXIDE			<12000	<4000			

Units: ug/l

Reviewed by: EM 13413

Date: 6/2/11

Verified by: Tracy A. Cole

Date: _____

JUN 03 2011

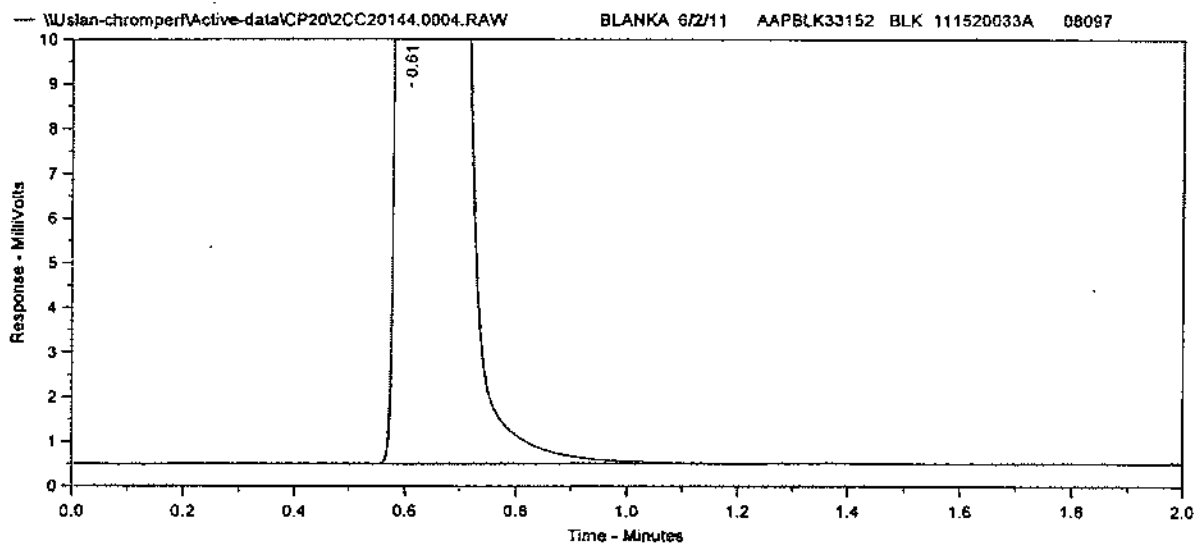
Tracy A. Cole
Senior Specialist

AK869 8871

* Recovery outside QC Limits

Printed on: 6/2/2011 17:48:18

Chrom Perfect Chromatogram Report



Sample Name = BLANKA 6/2/11 AAPBLK33152 BLK 111520033A 08097

Instrument = CP20
Detector = H3145A

Raw File Name = \\Uslan-chromperf\Active-data\CP20\2CC20144.0004.RAW

Method File Name = \\USLAN-CHROMPERF\ACTIVE-DATA\CP20\CO2.MET

Calibration File Name = Y:\CP20\2C20144.CAL

Time Injected = 6/2/2011 4:06:24 PM

Run Time = 2

Dilution Factor = 5

Amount Injected = 1

Peak Threshold = 0

Peak Width = 0.05

Operator = 2343

Incubate 5 minutes at 35C; 40C for .1 minute; 40C/min to 80C; Hold 1 min
Column: CTR 1 COLUMN; 6' X 1/4" X .210 STAINLESS STEEL

Ret. Time	Amount	Amt %	Area
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AKES9 8872

Extraction/Distillation/Digestion Logs

Organic Extraction Batchlog

Assigned to: 2343 Elizabeth Marin

Reviewed by: EM1313Start Date: 6/2/11Start time: 15:30**111520033A**Tech 1: EM1313

Tech 2: _____

Dept: 32 Prep Analysis: 00000

CO2 by Headspace

QC	Sample Code	Amt (mL)	SS/IS Sol.	Amt (mL)	MS Sol.	Amt (mL)	FV (mL)	pH	pH	BC	Comments
6296060MSD	6GW05MSD	5			MS99173B	0.1	5	7	6.5H		
6296060MS	6GW05MS	↓			MS99173B	↓	↓				
BLANKA	PBLK33152										
LCSA	LCS33152	↓			MS99173B	0.1	↓				

Spike Solutions:

MS99173B

Witness:

CO2 SPIKE

Sample #	Sample Code	Amt (mL)	SS/IS Sol.	Amt (mL)	FV (mL)	pH	pH	BC	Comments	Analyses	Due Date	Prio
1	6296060	5			5	NA	6.5H			08097	06/03/2011	P
2	6296061	↓			↓					08097	06/03/2011	P
3	6296062	↓			↓					08097	06/03/2011	P
4	6301430	↓			↓					08097	06/06/2011	P

Rack ID: <u>10</u>	Work Station
Internal Standard	Balance #

DF = Dilution Factor FV = Final Volume

Page 1 of 1

S-bath ID	C	S-bath ID	C	N-Evap	C	M-vap	C	111520033A
-----------	---	-----------	---	--------	---	-------	---	------------

Documented temps are NIST corrected.

B10111520033A

TPH-DRO by GC Data

Case Narrative Conformance/Non-Conformance Summary



CLIENT: ChevronTexaco
SDG: AKE89

EPH/Miscellaneous GC

Fraction: TPH-DRO by GC

TPH-DRO AK water C10-C25

<u>Sample #</u>	<u>Client ID</u>	<u>Matrix</u>		<u>Comments</u>
		<u>Liquid</u>	<u>Solid</u>	
6301430	MW-4-052611 Grab Water Sample	X		DF25
6301431	DUP-1-052611 Grab Water Sample	X		DF50; Field Duplicate Sample

See QC Reference List for Associated Batch QC Samples

SAMPLE PREPARATION:

No problems were encountered with the preparation of the samples.

ANALYSIS:

Dilutions are listed in the table above.

Due to software limitations, form 7's (check standard summary), or form 6s (Initial Calibration and Retention Time Summaries) and the form 8's cannot be automatically generated. Raw data/Chrom Perfect Sequence files containing this information are in the Quality Control and Calibration Summary Forms section of this data package.

QUALITY CONTROL AND NONCONFORMANCE SUMMARY:

Site specific matrix QC samples were not submitted for this SDG. The batch matrix QC was performed on samples from another project. Therefore the matrix effects would not be relevant and matrix QC is not provided in the data package. Laboratory spike data (LCS) are provided.

All QC is within specification.


DATA INTERPRETATION:

No further interpretation is necessary for the data submitted.

Abbreviation Key

UNSPK = Unspiked (for MS/MSD)	LOQ = Limit of Quantitation
MS = Matrix Spike	MDL = Method Detection Limit
MSD = Matrix Spike Duplicate	ND = Not Detected
BKG = Background (for Duplicate)	J = Estimated Value
D = Duplicate (DUP)	E= out of calibration range
LCS = Lab Control Sample	
LCSD = Lab Control Sample Duplicate	* = Out of Specification

Narrative Reviewed and Approved 6/28/11 by
(Date)


Elizabeth A. Smith
Specialist

AKE89 8877



CONFORMANCE/NON-CONFORMANCE SUMMARY

SDG: AKE89

Indicate Yes, No, N/A


- | | |
|---|-----|
| 1. Chromatograms labeled / Compounds identified (Field Samples & Method Blanks) | YES |
| 2. Retention times for chromatograms provided | YES |
| 3. Standards summary meet criteria | YES |
| 4. Calibration - Initial calibration performed before sample analysis and continuing calibration performed within 24 hours of sample analysis. | YES |
| 5. Blank contamination
If yes, list compounds and concentrations in each blank: N/A | NO |
| 6. Surrogate recoveries meet criteria
If not met, list those compounds and the recoveries that fall outside the acceptable range: N/A
If not met, were the calculations checked and the results qualified as "estimated"? N/A | YES |
| 7. Matrix Spike / Matrix Spike Duplicate recoveries meet criteria.
If not met, list those compounds and the recoveries that fall outside the acceptable range: | N/A |
| 8. Laboratory Control Sample / Laboratory Control Sample Duplicate meet criteria.
If not met, list those compounds and the recoveries that fall outside the acceptable range: | YES |
| 9. Retention time summaries for primary and confirmation analyses meet criteria | N/A |
| 10. Were samples run on dissimilar columns? | N/A |
| 11. Extraction holding time met
If not met, list number of days exceeded for each sample: N/A | YES |
| 12. Analysis holding time met
If not met, list number of days exceeded for each sample: N/A | YES |

Additional Comments:

~~Summary reviewed and approved by:~~


Elizabeth A. Smith

~~Specialist~~
Dana Kauffman, Manager Data Deliverables


Date

AKE89 8878

Quality Control and Calibration Summary Forms



Quality Control Reference List
EPH/Miscellaneous GC

CLIENT: ChevronTexaco
SDG: AKE89

Fraction: TPH-DRO by GC

Analysis

TPH-DRO AK water C10-C25

Batch Number

111520013A

Sample Number

PBLK13152

LCS13152

LCSD13152

6301430

6301431

Analysis Date

06/02/2011 18:43:00

06/02/2011 19:11:00

06/02/2011 19:38:00

06/03/2011 22:10:00

06/03/2011 22:37:00

AKE89 0000

Fraction: TPH-DRO by GC

111520013 / PBLK13152					
Analyte	Analysis Date	Blank Results	Units	MDL	LOQ
TPH-DRO AK water C10-C25	06/02/11	N.D.	mg/l	0.050	0.25

~~AKE89 0601~~

Fraction: TPH-DRO by GC

Sample	Orthoterphenyl	
	Spike Added	Limits
111520013A	11.9 mg/l	
6301430	107	50 - 150
6301431	101	50 - 150
LCS13152	109	60 - 120
LCSD13152	110	60 - 120
PBLK13152	102	60 - 120

AKE89 0002



Quality Control Summary
Laboratory Control Standard (LCS)
Laboratory Control Standard Duplicate(LCSD)

SDG: AKE89
Matrix: LIQUID

EPH/Miscellaneous GC
Fraction: TPH-DRO by GC

LCS: LCS13152 LCSD: LCSD13152 Analyte	Batch: 111520013A (Sample number(s): 6301430-6301431)							
	Spike Added mg/l	LCS Conc mg/l	LCSD Conc mg/l	LCS %Rec	LCSD %Rec	%Rec Limits	%RPD	%RPD Limits
TPH-DRO AK water C10-C25	.8	.82	.84	102	105	75-125	2	20

AKE89 8883

Calibration File Name: C:\CPWIN\DATA1\AKDL047A.CAL Version = 3

External standard calibration

No injection volume correction

No sample weight correction

Area reject threshold = 0

Reference peak area reject threshold = 0

Amount units = PPM

4 components with 5 levels each

LLW0007
2/17/11

1 DRO RF C10-<C25

Retention time = 0.001 min., Search window = 0.000 min.

Low alarm amount = 0, High alarm amount = 0

Group number = 0, Component constant = 20482.11

No retention time reference component

Single peak quantification by area

Level	Amount	Area	Area/Amt	Source	Date and time
1	100.000	2097613.0	20976.13	Manual	2/17/2011 11:49:
2	400.000	7984342.0	19960.86	Manual	2/17/2011 11:49:
3	800.000	16525130.0	20656.41	Manual	2/17/2011 11:49:
4	1600.000	32450590.0	20281.62	Manual	2/17/2011 11:49:
5	3200.000	65713620.0	20535.51	Manual	2/17/2011 11:49:

USLO 1047.03-13R
1047.21R MRL
1047.23R 1CV

✓ 1047.24 21711

Calibration formula: $Y = 20482.11 X$

Fit type = Avg CF with equal weighting, forced to origin

Coefficient of determination = 0.9999, Average error = 1.41%

Average CF = 20482.1100 with RSD = 1.87%

2 C10

Retention time = 2.840 min., Search window = 0.050 min.

Low alarm amount = 0, High alarm amount = 0

Group number = 0, Component constant = 0

No retention time reference component

Single peak quantification by area

Level	Amount	Area	Area/Amt	Source	Date and time
1	1.000	0.0	0	Manual	2/17/2011 11:49:
2	-1.000	0		Manual	2/17/2011 11:49:
3	-1.000	0		Manual	2/17/2011 11:49:
4	-1.000	0		Manual	2/17/2011 11:49:
5	-1.000	0		Manual	2/17/2011 11:49:

Calibration formula: No data points to graph

Fit type = Avg CF with equal weighting, forced to origin

Coefficient of determination = 1.0000, Average error = 100.00%

Average CF = 0.0000 with RSD = 0.00%

3 o-Terphenyl SURR

Retention time = 10.130 min., Search window = 0.050 min.

Low alarm amount = 0, High alarm amount = 0

Group number = 0, Component constant = 23783.08

AKD89 8884

No retention time reference component
Single peak quantification by area

Level	Amount	Area	Area/Amt	Source	Date and time
1	2.000	48477.0	24238.5	Manual	2/17/2011 11:49:
2	8.000	191265.0	23908.13	Manual	2/17/2011 11:49:
3	16.000	377747.0	23609.19	Manual	2/17/2011 11:49:
4	20.000	475463.0	23773.15	Manual	2/17/2011 11:49:
5	40.000	935457.0	23386.43	Manual	2/17/2011 11:49:

Calibration formula: $Y = 23783.08 X$

Fit type = Avg CF with equal weighting, forced to origin

Coefficient of determination = 0.9994, Average error = 0.98%

Average CF = 23783.0800 with RSD = 1.35%

4 C25

Retention time = 12.360 min., Search window = 0.100 min.

Low alarm amount = 0, High alarm amount = 0

Group number = 0, Component constant = 0

No retention time reference component

Single peak quantification by area

Level	Amount	Area	Area/Amt	Source	Date and time
1	1.000	0.0	0	Manual	2/17/2011 11:49:
2	-1.000	0		Manual	2/17/2011 11:49:
3	-1.000	0		Manual	2/17/2011 11:49:
4	-1.000	0		Manual	2/17/2011 11:49:
5	-1.000	0		Manual	2/17/2011 11:49:

Calibration formula: No data points to graph

Fit type = Avg CF with equal weighting, forced to origin

Coefficient of determination = 1.0000, Average error = 100.00%

Average CF = 0.0000 with RSD = 0.00%

AKS89 0005

Calibration File Name: C:\CPWIN\DATA1\AKDL047B.CAL Version = 1

External standard calibration

No injection volume correction

No sample weight correction

Area reject threshold = 0

Reference peak area reject threshold = 0

Amount units = PPM

4 components with 5 levels each

1 DRO RF C10-<C25

Retention time = 0.001 min., Search window = 0.000 min.

Low alarm amount = 0, High alarm amount = 0

Group number = 0, Component constant = 20482.11

No retention time reference component

Single peak quantification by area

Level	Amount	Area	Area/Amt	Source	Date and time
1	100.000	2097613.0	20976.13	Manual	2/17/2011 11:49:
2	400.000	7984342.0	19960.86	Manual	2/17/2011 11:49:
3	800.000	16525130.0	20656.41	Manual	2/17/2011 11:49:
4	1600.000	32450590.0	20281.62	Manual	2/17/2011 11:49:
5	3200.000	65713620.0	20535.51	Manual	2/17/2011 11:49:

Calibration formula: $Y = 20482.11 X$

Fit type = Avg CF with equal weighting, forced to origin

Coefficient of determination = 0.9999, Average error = 1.41%

Average CF = 20482.1100 with RSD = 1.87%

2 C10

Retention time = 2.800 min., Search window = 0.050 min.

Low alarm amount = 0, High alarm amount = 0

Group number = 0, Component constant = 0

No retention time reference component

Single peak quantification by area

Level	Amount	Area	Area/Amt	Source	Date and time
1	1.000	0.0	0	Manual	2/17/2011 11:49:
2	-1.000	0		Manual	2/17/2011 11:49:
3	-1.000	0		Manual	2/17/2011 11:49:
4	-1.000	0		Manual	2/17/2011 11:49:
5	-1.000	0		Manual	2/17/2011 11:49:

Calibration formula: No data points to graph

Fit type = Avg CF with equal weighting, forced to origin

Coefficient of determination = 1.0000, Average error = 100.00%

Average CF = 0.0000 with RSD = 0.00%

3 o-Terphenyl SURR

Retention time = 10.090 min., Search window = 0.050 min.

Low alarm amount = 0, High alarm amount = 0

Group number = 0, Component constant = 23783.08

AKDL047B.CAL

No retention time reference component

Single peak quantification by area

Level	Amount	Area	Area/Amt	Source	Date and time
1	2.000	48477.0	24238.5	Manual	2/17/2011 11:49:
2	8.000	191265.0	23908.13	Manual	2/17/2011 11:49:
3	16.000	377747.0	23609.19	Manual	2/17/2011 11:49:
4	20.000	475463.0	23773.15	Manual	2/17/2011 11:49:
5	40.000	935457.0	23386.43	Manual	2/17/2011 11:49:

Calibration formula: $Y = 23783.08 X$

Fit type = Avg CF with equal weighting, forced to origin

Coefficient of determination = 0.9994, Average error = 0.98%

Average CF = 23783.0800 with RSD = 1.35%

4 C25

Retention time = 12.330 min., Search window = 0.100 min.

Low alarm amount = 0, High alarm amount = 0

Group number = 0, Component constant = 0

No retention time reference component

Single peak quantification by area

Level	Amount	Area	Area/Amt	Source	Date and time
1	1.000	0.0	0	Manual	2/17/2011 11:49:
2	-1.000	0		Manual	2/17/2011 11:49:
3	-1.000	0		Manual	2/17/2011 11:49:
4	-1.000	0		Manual	2/17/2011 11:49:
5	-1.000	0		Manual	2/17/2011 11:49:

Calibration formula: No data points to graph

Fit type = Avg CF with equal weighting, forced to origin

Coefficient of determination = 1.0000, Average error = 100.00%

Average CF = 0.0000 with RSD = 0.00%

AKES9 8887

Sample ID: AKCDX1132A BCAKCDXBC CCAL 114699999
 Instrument ID: CP24--H5386A Injected on: 2/17/2011 6:04:40 AM
 Volume Inj. per Column: 1 GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C
 Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN
 Sample Amount: 1 Dilution Factor: 1
 Analyst: 2027

Peak #	Ret Time (min)	Peak Name	Amount PPM	Peak Area	Peak Width (min)	Peak Height
28	2.835	C10		35954	.021	17489
148	10.132	o-Terphenyl SURR	14.849	353155	.022	198072
174	12.35	C25		2194	.036	632

Slice	Start Time	Stop Time	Slice Amount	Amount %	Slice Area	Area %
1	2.740	12.260	14.849	100.000	10944430.0	46.761
2	10.080	10.180	14.849	100.000	353155.1	1.509

Total slice amount= 29.698
 Total slice amount %= 200.0

Total slice area= 11297590.0
 Total slice area %= 48.3

***** RESULTS TABLE *****

C10 - <C25 DRO AMT = 517.0989
 % Level 2 DRO Difference = 29.27474
 % Level 3 DRO Difference = -35.36263
 % Level 4 DRO Difference = -67.68131

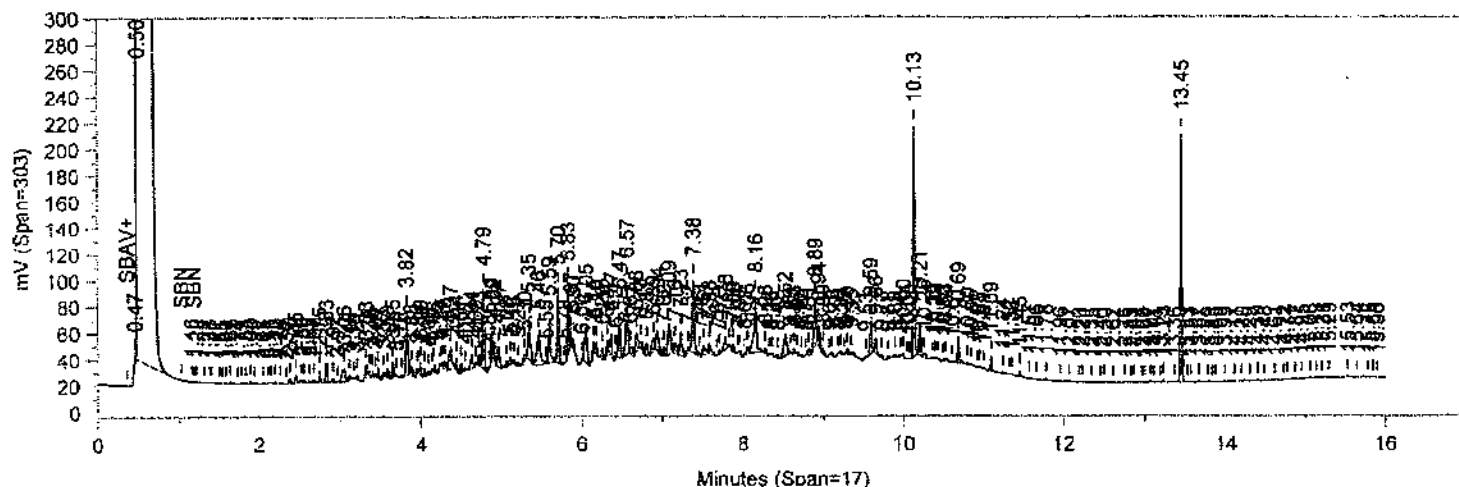
$$\frac{517.0989}{500} - 1 \times 100 = 3.4\% D$$

ELLW200
 2/17/11

FILES:

Area File: C:\CPWIN\DATA\1\047.23A
 Method File: C:\CPWIN\DATA\1\AKDLSTD.MET
 Calibration File: C:\CPWIN\DATA\1\AKDL047A.CAL
 Format File: C:\CPWIN\DATA\1\AKDLSTD.FMT
 Area file created on: 2/17/2011 11:57:14 AM
 File reported on: 2/17/2011 at 11:57:16 AM

AK289 8888



Sample Name: AKCDX1132A BKA KC DXBC CCAL 114699999A

Instrument ID: CP24-H5386A Injected on: 2/17/2011 6:04:40 AM
Volume Inj. per Column: 1 GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C
Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN
Sample Amount: 1 Dilution Factor: 1
Analyst: 2027

Peak #	Ret Time (min)	Peak Name	Amount PPM	Peak Area	Peak Width (min)	Peak Height
30	2.835	C10	.	21819	.021	15304
150	10.132	o-Terphenyl SURR	10.6634	253609	.022	178302
177	12.35	C25	.	1029	.036	440

Slice	Start Time	Stop Time	Slice Amount	Amount %	Slice Area	Area %
1	10.080	10.180	10.663	100.000	253609.5	1.774

Total slice amount= 10.663
Total slice amount %= 100.0

Total slice area= 253609.5
Total slice area %= 1.8

o-Terphenyl Level 2 % Difference = 6.634426 %
o-Terphenyl Level 3 % Difference = -46.68279 %
o-Terphenyl Level 4 % Difference = -73.34139 %

FILES:
Area File: C:\CPWIN\DATA1\047.23A
Method File: C:\CPWIN\DATA1\REAKDLST.MET
Calibration File: C:\CPWIN\DATA1\AKDL047A.CAL
Format File: C:\CPWIN\DATA1\REAKDLST.FMT
Area file created on: 2/17/2011 11:57:28 AM
File reported on: 2/17/2011 at 11:57:29 AM

AKES9 0089

AK 102/103

Sample ID: AKFL21132A SDAKFL2SD CCAL 1115299999
Instrument ID: CP24-H5386A Injected on: 6/2/2011 6:16:16 PM
Volume Inj. per Column: 1 GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C
Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN
Sample Amount: 1 Dilution Factor: 1
Analyst: 2027

Peak #	Ret Time (min)	Peak Name	Amount PPM	Peak Area	Peak Width (min)	Peak Height
30	2.774	C10		75421	.022	42255
137	10.056	o-Terphenyl SURR	13.672	325162	.023	173117
161	12.313	C25		7192	.026	2953

Slice	Start Time	Stop Time	Slice Amount	Amount %	Slice Area	Area %
1	2.700	12.230	13.672	100.000	8471254.0	39.476
2	10.040	10.140	13.672	100.000	467338.7	2.178

Total slice amount= 27.344
Total slice amount %= 200.0

Total slice area= 8938593.0
Total slice area %= 41.7

***** RESULTS TABLE *****

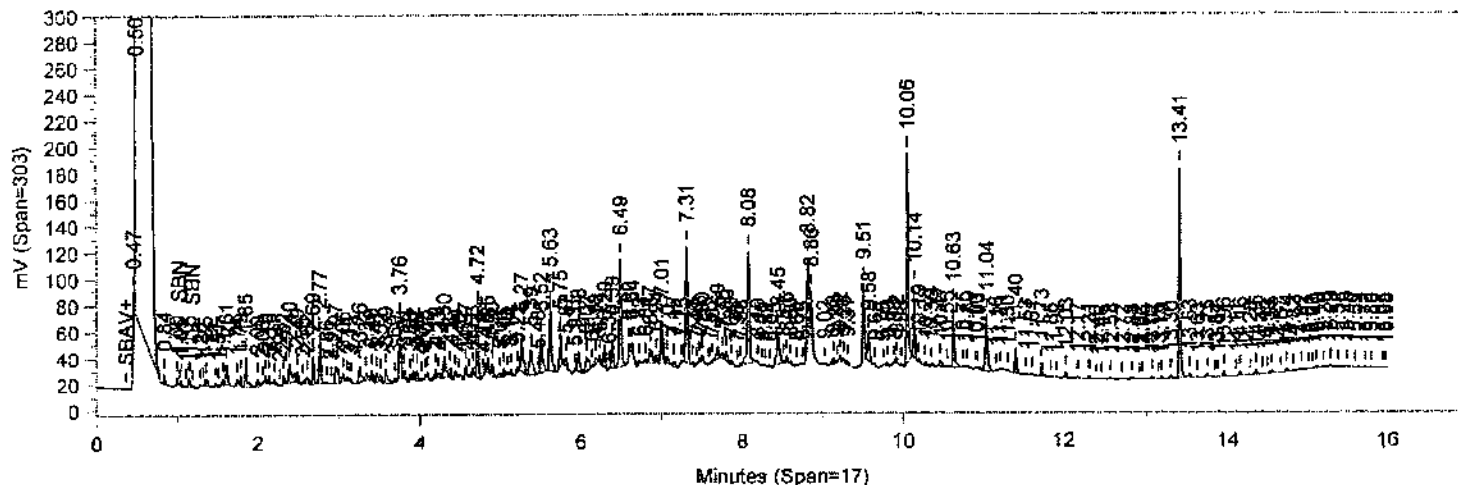
C10 - <C25 DRO AMT = 390.7759
% Level 2 DRO Difference = -2.306014 ✓
% Level 3 DRO Difference = -51.153
% Level 4 DRO Difference = -75.5765

FILES:

Area File: C:\CPWIN\DATA\1\152.41A
Method File: C:\CPWIN\DATA\1\AKDLSTD.MET
Calibration File: C:\CPWIN\DATA\1\AKDL047B.CAL
Format File: C:\CPWIN\DATA\1\AKDLSTD.FMT
Area file created on: 6/2/2011 6:32:24 PM
File reported on: 6/2/2011 at 6:32:26 PM

AKES5 6898

AK 102/103 SURROGATE
 AKFL21132A SDAKFL2SD CCAL 1115299999
 C:\CPWINDATA\1\152.41R



Sample Name: AKFL21132A SDAKFL2SD CCAL 1115299999A

Instrument ID: CP24--H5386A

Injected on: 6/2/2011 6:16:16 PM

Volume Inj. per Column: 1

GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C

Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN

Sample Amount: 1

Dilution Factor: 1

Analyst: 2027

Peak #	Ret Time (min)	Peak Name	Amount PPM	Peak Area	Peak Width (min)	Peak Height
30	2.774	C10	.	58545	.022	39656
138	10.056	o-Terphenyl SURR	10.0851	239855	.023	158835
162	12.313	C25	.	4298	.026	2460

Slice	Start Time	Stop Time	Slice Amount	Amount %	Slice Area	Area %
1	10.040	10.140	10.085	100.000	318419.4	2.161

Total slice amount= 10.085
 Total slice amount %= 100.0

Total slice area= 318419.4
 Total slice area %= 2.2

o-Terphenyl Level 2 % Difference = 0.8509874 %

o-Terphenyl Level 3 % Difference = -49.5745 %

o-Terphenyl Level 4 % Difference = -74.78725 %

FILES:

Area File: C:\CPWINDATA\1\152.41A

Method File: C:\CPWINDATA\1\REAKDLST.MET

Calibration File: C:\CPWINDATA\1\AKDL047B.CAL

Format File: C:\CPWINDATA\1\REAKDLST.FMT

Area file created on: 6/2/2011 6:32:38 PM

File reported on: 6/2/2011 at 6:32:39 PM

AK289 8891

Sample ID: AKFL31132A SIAKFL3SI CCAL 1115299999
 Instrument ID: CP24--H5386A Injected on: 6/2/2011 11:17:25 PM
 Volume Inj. per Column: 1 GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C
 Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN
 Sample Amount: 1 Dilution Factor: 1
 Analyst: 2027

Peak #	Ret Time (min)	Peak Name	Amount PPM	Peak Area	Peak Width (min)	Peak Height
31	2.773	C10	.	158136	.022	89283
143	10.057	o-Terphenyl SURR	28.9906	689485	.023	373150
166	12.309	C25	.	16031	.023	7015

Slice	Start Time	Stop Time	Slice Amount	Amount %	Slice Area	Area %
1	2.700	12.230	28.991	100.000	17679990.0	56.055
2	10.040	10.140	28.991	100.000	982861.3	3.116

Total slice amount= 57.981
 Total slice amount %= 200.0

Total slice area= 18662850.0
 Total slice area %= 59.2

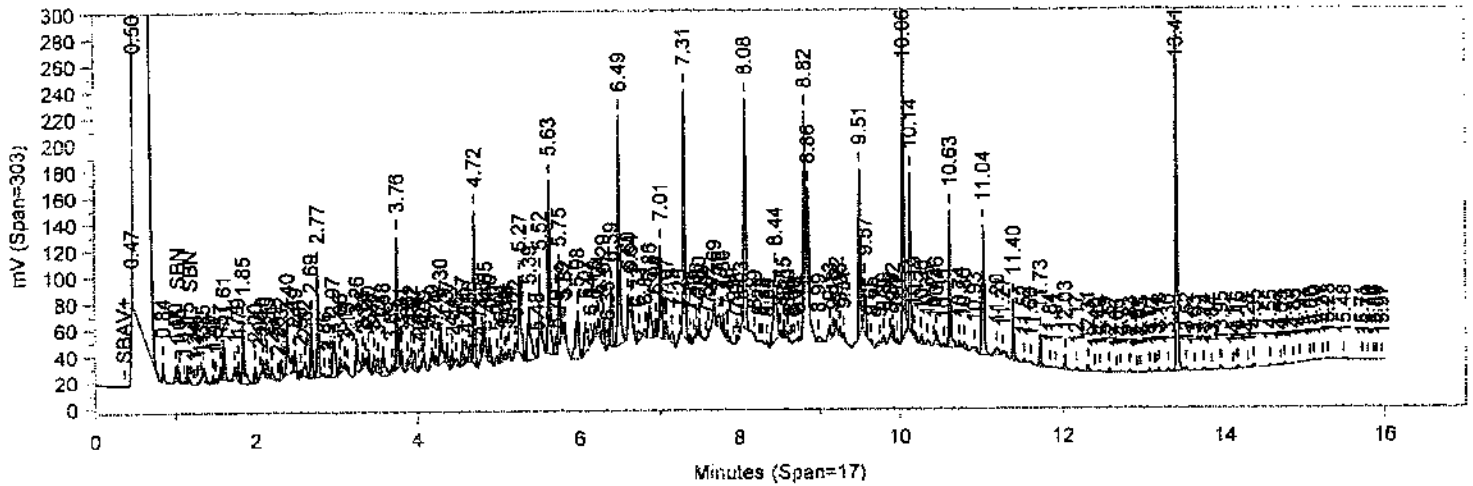
***** RESULTS TABLE *****

C10 - <C25 DRO AMT = 815.2053
 % Level 2 DRO Difference = 103.8013
 % Level 3 DRO Difference = 1.900661 ✓
 % Level 4 DRO Difference = -49.04967

FILES:

Area File: C:\CPWIN\DATA\1\152.52A
 Method File: C:\CPWIN\DATA\1\AKDLSTD.MET
 Calibration File: C:\CPWIN\DATA\1\AKDL047B.CAL
 Format File: C:\CPWIN\DATA\1\AKDLSTD.FMT
 Area file created on: 6/2/2011 11:33:34 PM
 File reported on: 6/2/2011 at 11:33:35 PM

AKES9 0892



Sample Name: AKFL31132A SIAKFL3SI CCAL 1115299999A

Instrument ID: CP24--H5386A Injected on: 6/2/2011 11:17:25 PM
 Volume Inj. per Column: 1 GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C
 Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN
 Sample Amount: 1 Dilution Factor: 1
 Analyst: 2027

Peak #	Ret Time (min)	Peak Name	Amount PPM	Peak Area	Peak Width (min)	Peak Height
31	2.773	C10	.	122986	.022	83878
143	10.057	o-Terphenyl SURR	21.4381	509864	.023	344062
166	12.309	C25	.	8850	.023	5819

Slice	Start Time	Stop Time	Slice Amount	Amount %	Slice Area	Area %
1	10.040	10.140	21.438	100.000	679985.1	3.681

Total slice amount= 21.438 Total slice area= 679985.1
 Total slice amount %= 100.0 Total slice area %= 3.7

o-Terphenyl Level 2 % Difference = 114.3811 %
 o-Terphenyl Level 3 % Difference = 7.190549 %
 o-Terphenyl Level 4 % Difference = -46.40472 %

FILES:
 Area File: C:\CPWINDATA\1\152.52A
 Method File: C:\CPWINDATA\1\REAKDLST.MET
 Calibration File: C:\CPWINDATA\1\AKDL047B.CAL
 Format File: C:\CPWINDATA\1\REAKDLST.FMT
 Area file created on: 6/2/2011 11:33:46 PM
 File reported on: 6/2/2011 at 11:33:49 PM

AKF89 0093

Sample ID: AKFL21132A SEAKFL2SE CCAL 1115399999
 Instrument ID: CP24-H5386A Injected on: 6/3/2011 9:42:46 PM
 Volume Inj. per Column: 1 GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C
 Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN
 Sample Amount: 1 Dilution Factor: 1
 Analyst: 2027

Peak #	Ret Time (min)	Peak Name	Amount PPM	Peak Area	Peak Width (min)	Peak Height
33	2.77	C10	.	81771	.023	44719
142	10.052	o-Terphenyl SURR	14.8361	352848	.023	187687
167	12.308	C25	.	7896	.023	3622

Slice	Start Time	Stop Time	Slice Amount	Amount %	Slice Area	Area %
1	2.700	12.230	14.836	100.000	9114715.0	41.215
2	10.040	10.140	14.836	100.000	511151.8	2.311

Total slice amount= 29.672
 Total slice amount %= 200.0

Total slice area= 9625867.0
 Total slice area %= 43.5

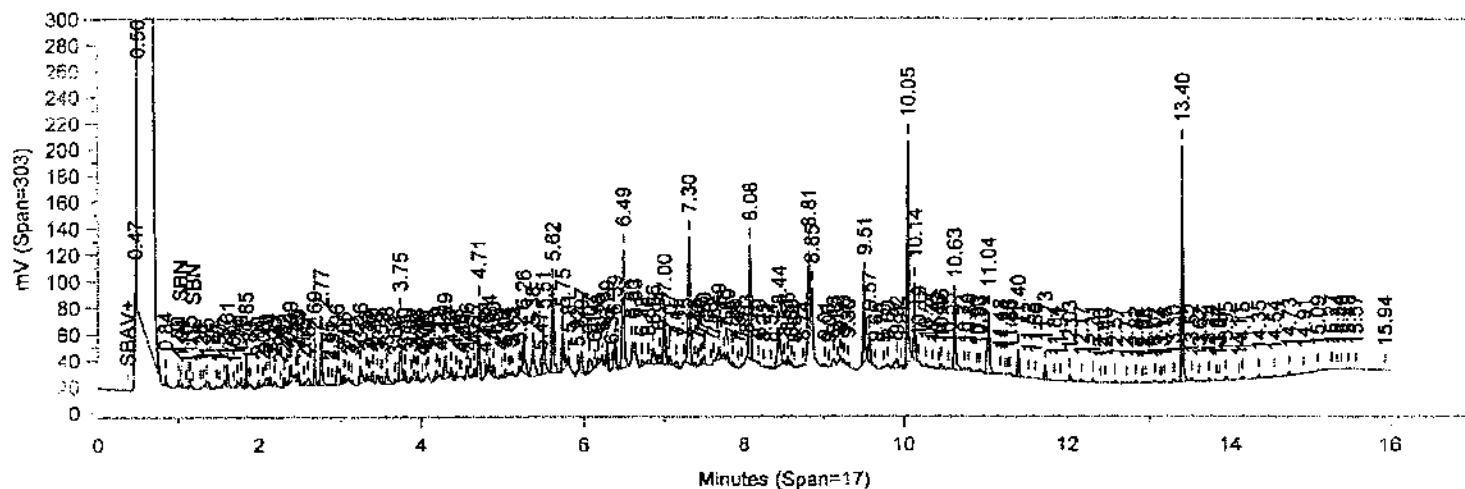
***** RESULTS TABLE *****

C10 - <C25 DRO AMT = 420.0526
 % Level 2 DRO Difference = 5.013144
 % Level 3 DRO Difference = -47.49343
 % Level 4 DRO Difference = -73.74672

FILES:

Area File: C:\CPWINDATA\152.81A
 Method File: C:\CPWINDATA\AKDLSTD.MET
 Calibration File: C:\CPWINDATA\AKDL047B.CAL
 Format File: C:\CPWINDATA\AKDLSTD.FMT
 Area file created on: 6/3/2011 9:58:54 PM
 File reported on: 6/3/2011 at 9:58:57 PM

AKFL21132A



Sample Name: AKFL21132A SEAKFL2SE CCAL 1115399999A

Instrument ID: CP24-H5386A Injected on: 6/3/2011 9:42:46 PM
Volume Inj. per Column: 1 GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C
Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN
Sample Amount: 1 Dilution Factor: 1
Analyst: 2027

Peak #	Ret Time (min)	Peak Name	Amount PPM	Peak Area	Peak Width (min)	Peak Height
33	2.77	C10	.	63423	.023	42018
143	10.052	o-Terphenyl SURR	10.7893	256602	.023	172229
169	12.308	C25	.	4733	.023	3093

Slice	Start Time	Stop Time	Slice Amount	Amount %	Slice Area	Area %
1	10.040	10.140	10.789	100.000	340711.1	2.290

Total slice amount= 10.789 Total slice area= 340711.1
Total slice amount %= 100.0 Total slice area %= 2.3

o-Terphenyl Level 2 % Difference = 7.892847 % ✓
o-Terphenyl Level 3 % Difference = -46.05358 %
o-Terphenyl Level 4 % Difference = -73.02679 %

FILES:

Area File: C:\CPWIN\DATA\1\152.81A
Method File: C:\CPWIN\DATA\1\REAKDLST.MET
Calibration File: C:\CPWIN\DATA\1\AKDL047B.CAL
Format File: C:\CPWIN\DATA\1\REAKDLST.FMT
Area file created on: 6/3/2011 9:59:08 PM
File reported on: 6/3/2011 at 9:59:10 PM

AK69 0095

Sample ID: AKFL31132A SJAKFL3SJ CCAL 1115399999
 Instrument ID: CP24--H5386A Injected on: 6/3/2011 11:04:41 PM
 Volume Inj. per Column: 1 GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C
 Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN
 Sample Amount: 1 Dilution Factor: 1
 Analyst: 2027

Peak #	Ret Time (min)	Peak Name	Amount PPM	Peak Area	Peak Width (min)	Peak Height
31	2.769	C10	.	184939	.023	93230
138	10.054	o-Terphenyl SURR	29.8238	709303	.023	392943
165	12.306	C25	.	18112	.023	7234

Slice	Start Time	Stop Time	Slice Amount	Amount %	Slice Area	Area %
1	2.700	12.230	29.824	100.000	18829220.0	57.591
2	10.040	10.140	29.824	100.000	1020538.0	3.121

Total slice amount= 59.648
 Total slice amount %= 200.0

Total slice area= 19849760.0
 Total slice area %= 60.7

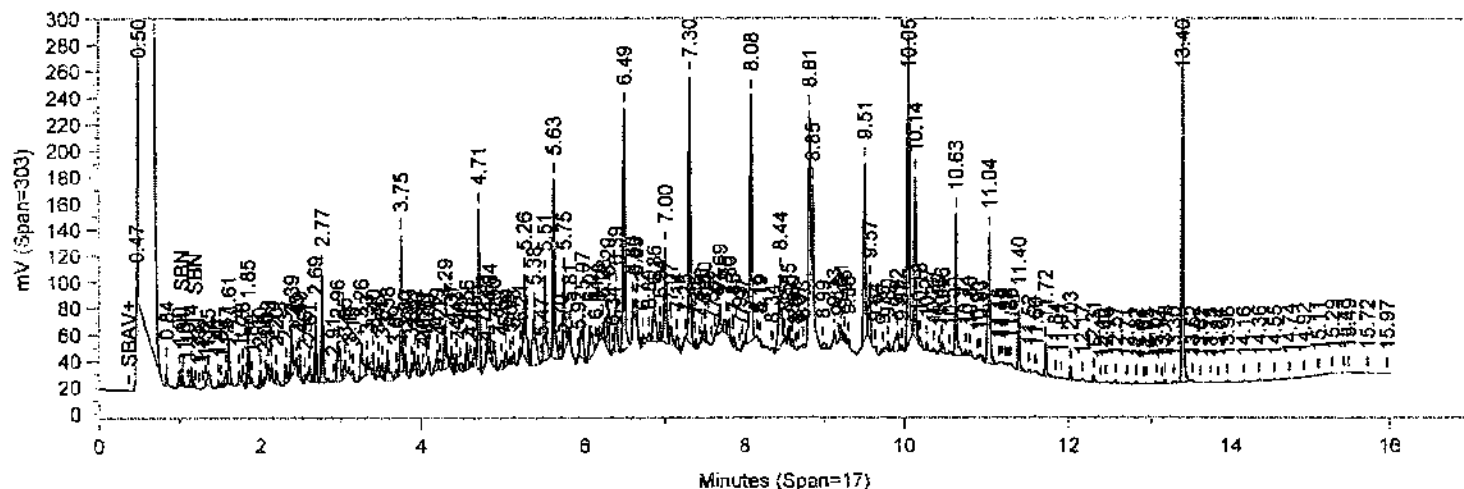
***** RESULTS TABLE *****

C10 - <C25 DRO AMT = 869.475
 % Level 2 DRO Difference = 117.3687
 % Level 3 DRO Difference = 8.684373
 % Level 4 DRO Difference = -45.65781

FILES:

Area File: C:\CPWIN\DATA\1\152.84A
 Method File: C:\CPWIN\DATA\1\AKDLSTD.MET
 Calibration File: C:\CPWIN\DATA\1\AKDL047B.CAL
 Format File: C:\CPWIN\DATA\1\AKDLSTD.FMT
 Area file created on: 6/3/2011 11:20:48 PM
 File reported on: 6/3/2011 at 11:20:51 PM

AK299 2096



Sample Name: AKFL31132A SJAKFL3SJ CCAL 1115399999A

Instrument ID: CP24--H5386A

Injected on: 6/3/2011 11:04:41 PM

Volume Inj. per Column: 1

GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C

Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN

Sample Amount: 1

Dilution Factor: 1

Analyst: 2027

Peak #	Ret Time (min)	Peak Name	Amount PPM	Peak Area	Peak Width (min)	Peak Height
32	2.769	C10	.	137699	.023	87948
139	10.054	o-Terphenyl SURR	22.4633	534247	.023	361048
166	12.306	C25	.	8430	.023	5626

Slice	Start Time	Stop Time	Slice Amount	Amount %	Slice Area	Area %
1	10.040	10.140	22.463	100.000	712925.4	3.812

Total slice amount= 22.463
Total slice amount %= 100.0

Total slice area= 712925.4
Total slice area %= 3.8

o-Terphenyl Level 2 % Difference = 124.633 %

o-Terphenyl Level 3 % Difference = 12.31651 %

o-Terphenyl Level 4 % Difference = -43.84174 %

FILES:

Area File: C:\CPWIN\DATA\1\152.84A

Method File: C:\CPWIN\DATA\1\REAKDLST.MET

Calibration File: C:\CPWIN\DATA\1\AKDL047B.CAL

Format File: C:\CPWIN\DATA\1\REAKDLST.FMT

Area file created on: 6/3/2011 11:21:02 PM

File reported on: 6/3/2011 at 11:21:04 PM

AK09 0097

Lancaster Laboratories
CHROM PERFECT SEQUENCE FILE

Sequence File: \\cp24\c-Drive\CPWIN\DATA1\L047.seq
 Chromatography Directory: \\cp24\c-Drive\CPWIN\DATA1
 Method Directory: \\cp24\c-Drive\CPWIN\DATA1
 Number of Entries: 30

SampleName	Code	ID	FileName	Method	Samp Amt	DF	Int Std	C	Batch Number	Analysis
1 CONDITIONER	MISC	AA	L047.01R	AKDLSTD.MET	1	1	1	0	114699999	
2 CONDITIONER	MISC	AA	L047.02R	AKDLSTD.MET	1	1	1	0	114699999	
3 AKRTX1032D	CCAL	LN	L047.03R	AKRTL.MET	1	1	1	0	114699999	
4 AKSS11032B	ICAL	AA	L047.04R	AKRLSTD.MET	1	1	1	0	114699999	
5 AKSS21032B	ICAL	AA	L047.05R	AKRLSTD.MET	1	1	1	0	114699999	
6 AKSS31032B	ICAL	AA	L047.06R	AKRLSTD.MET	1	1	1	0	114699999	
7 AKSS41032B	ICAL	AA	L047.07R	AKRLSTD.MET	1	1	1	0	114699999	
8 AKSS51032B	ICAL	AA	L047.08R	AKRLSTD.MET	1	1	1	0	114699999	
9 1FUL11132A	ICAL	AA	L047.09R	AKDLSTD.MET	1	1	1	0	114699999	
10 1FUL21132A	ICAL	AA	L047.10R	AKDLSTD.MET	1	1	1	0	114699999	
11 1FUL31132B	ICAL	AA	L047.11R	AKDLSTD.MET	1	1	1	0	114699999	
12 1FUL41132A	ICAL	AA	L047.12R	AKDLSTD.MET	1	1	1	0	114699999	
13 1FUL51132A	ICAL	AA	L047.13R	AKDLSTD.MET	1	1	1	0	114699999	
14 MECL2	MISC	AA	L047.14R	AKDLSTD.MET	1	1	1	0	114699999	
15 AKSW11032B	ICAL	AA	L047.15R	AKRLSTD.MET	1	1	1	0	114699999	
16 AKSW21032B	ICAL	AA	L047.16R	AKRLSTD.MET	1	1	1	0	114699999	
17 AKSW31032B	ICAL	AA	L047.17R	AKRLSTD.MET	1	1	1	0	114699999	
18 AKSW41032B	ICAL	AA	L047.18R	AKRLSTD.MET	1	1	1	0	114699999	
19 AKSW51032B	ICAL	AA	L047.19R	AKRLSTD.MET	1	1	1	0	114699999	
20 MECL2	MISC	AA	L047.20R	AKDLSTD.MET	1	1	1	0	114699999	
21 1MDLX1132A	CCAL	IL	L047.21R	AKDLSTD.MET	1	1	1	0	114699999	
22 AKMDX1132A	CCAL	CA	L047.22R	AKRLSTD.MET	1	1	1	0	114699999	
23 AKCDX1132A	CCAL	BC	L047.23R	AKDLSTD.MET	1	1	1	0	114699999	
24 AKCRX1132A	CCAL	BD	L047.24R	AKRLSTD.MET	1	1	1	0	114699999	
25 AKRTX1032D	CCAL	LO	L047.25R	AKRTL.MET	1	1	1	0	114699999	
26 MS1104732C	CCAL	AN	L047.26R	AKRLSTD.MET	1	1	1	0	114699999	
27 MS1102132C	CCAL	AO	L047.27R	AKRLSTD.MET	1	1	1	0	114699999	
28 MS1103332A	CCAL	AP	L047.28R	AKRLSTD.MET	1	1	1	0	114699999	
29 SS1101932B	CCAL	AL	L047.29R	AKRLSTD.MET	1	1	1	0	114699999	
30 MS1104732A	CCAL	AO	L047.30R	TNLCK.MET	1	1	1	0	114699999	

AKED9 0090

Set-up by:

North Williams

Date:

2/17/11

2/17/11

Page 1 of 1

Lancaster Laboratories

CHROM PERFECT SEQUENCE FILE

Sequence File: \\cp24\C-Drive\CPWIN\DATA1\L124.seq

Chromatography Directory: \\cp24\C-Drive\CPWIN\DATA1

Method Directory: \\cp24\C-Drive\CPWIN\DATA1

Number of Entries: 33

SampleName	Code	ID	FileName	Method	Samp Amt	DF	Int Std	C	Batch Number	Analysis
1 CONDITIONER	MISC	AA	L124.01R	AKDLSTD.MET	1	1	1	0	1112399999	
2 CONDITIONER	MISC	AA	L124.02R	AKDLSTD.MET	1	1	1	0	1112399999	
3 AKRTX1132A	CCAL	NE	L124.03R	AKRTL.MET	1	1	1	0	1112399999	
4 AKFL41032D	CCAL	QH	L124.04R	AKDLSTD.MET	1	1	1	0	1112399999	
5 BLANKA 5/4/11	BLK	AA	L124.05R	AKDLSUM.MET	1000	1	1	0	111230025A	01741
6 LCSA 5/4/11	LCS	AA	L124.06R	AKDLSUM.MET	1000	1	1	0	111230025A	01741
7 LCSDA 5/4/11	LCSD	AA	L124.07R	AKDLSUM.MET	1000	1	1	0	111230025A	01741
8 6272432	T	AA	L124.08R	AKDLSUM.MET	1044	1	1	0	111230025A	01741
9 6272434	T	AA	L124.09R	AKDLSUM.MET	1027	1	1	0	111230025A	01741
10 6272435	T	AA	L124.10R	AKDLSUM.MET	1018	1	1	0	111230025A	01741
11 6272436	T	AA	L124.11R	AKDLSUM.MET	1026	1	1	0	111230025A	01741
12 6272437	T	AA	L124.12R	AKDLSUM.MET	1021	1	1	0	111230025A	01741
13 6272443	T	AA	L124.13R	AKDLSUM.MET	903	1	1	0	111230025A	01741
14 6272444	T	AA	L124.14R	AKDLSUM.MET	1027	1	1	0	111230025A	01741
15 6272445	T	AA	L124.15R	AKDLSUM.MET	1030	1	1	0	111230025A	01741
16 6272433	T	AA	L124.16R	AKDLSUM.MET	1035	1	1	0	111230025A	01741
17 AKFL21032D	CCAL	RV	L124.17R	AKDLSTD.MET	1	1	1	0	1112399999	
18 AKCK21132A	CCAL	OM	L124.18R	AKRLSTD.MET	1	1	1	0	1112399999	
19 BLANKA 5/4/11	BLK	AA	L124.19R	AKRLSUM.MET	25	1	1	0	111240006A	01738
20 LCSA 5/4/11	LCS	AA	L124.20R	AKRLSUM.MET	25	1	1	0	111240006A	01738
21 LCSDA 5/4/11	LCSD	AA	L124.21R	AKRLSUM.MET	25	1	1	0	111240006A	01738
22 6274142	T	AA	L124.22R	AKRLSUM.MET	25	1	1	0	111240006A	01738
23 6274142MS	MS	AA	L124.23R	AKRLSUM.MET	25	1	1	0	111240006A	01738
24 6274142MSD	MSD	AA	L124.24R	AKRLSUM.MET	25	1	1	0	111240006A	01738
25 6274143	T	AA	L124.25R	AKRLSUM.MET	25	1	1	0	111240006A	01738
26 6274144	T	AA	L124.26R	AKRLSUM.MET	25	1	1	0	111240006A	01738
27 6274145	T	AA	L124.27R	AKRLSUM.MET	25	1	1	0	111240006A	01738
28 6274146	T	AA	L124.28R	AKRLSUM.MET	25.1	1	1	0	111240006A	01738
29 6274147	T	AA	L124.29R	AKRLSUM.MET	25.1	1	1	0	111240006A	01738
30 6274148	T	AA	L124.30R	AKRLSUM.MET	25	1	1	0	111240006A	01738
31 MECL2	MISC	AA	L124.31R	AKRLSTD.MET	1	1	1	0	1112399999	
32 AKCK31132A	CCAL	NU	L124.32R	AKRLSTD.MET	1	1	1	0	1112399999	
33 AKRTX1132A	CCAL	NF	L124.33R	AKRTL.MET	1	1	1	0	1112399999	

AKES 0099

Set-up by:

H. W. Williams

Date:

5/5/11

5/5/11

Lancaster Laboratories

CHROM PERFECT SEQUENCE FILE

Sequence File: \\cp24\C-Drive\CPWINDATA1\L152.seq

Chromatography Directory: \\cp24\C-Drive\CPWINDATA1

Method Directory: \\cp24\C-Drive\CPWINDATA1

Number of Entries: 85

SampleName	Code	ID	FileName	Method	Samp Amt	DF	Int Std	C	Batch Number	Analysis
1 CONDITIONER	MISC	AA	L152.01R	WILSTD.MET	1	1	1	0	1115199999	
2 CONDITIONER	MISC	AA	L152.02R	WILSTD.MET	1	1	1	0	1115199999	
3 CONDITIONER	MISC	AA	L152.03R	WILSTD.MET	1	1	1	0	1115199999	
4 WISC11132A	ICAL	AA	L152.04R	WILSTD.MET	1	1	1	0	1115199999	
5 WISC21132A	ICAL	AA	L152.05R	WILSTD.MET	1	1	1	0	1115199999	
6 WISC31132A	ICAL	AA	L152.06R	WILSTD.MET	1	1	1	0	1115199999	
7 WISC41132A	ICAL	AA	L152.07R	WILSTD.MET	1	1	1	0	1115199999	
8 WISC51132A	ICAL	AA	L152.08R	WILSTD.MET	1	1	1	0	1115199999	
9 MECL2	MISC	AA	L152.09R	WILSTD.MET	1	1	1	0	1115199999	
10 TPH_41132F	CCAL	NY	L152.10R	WILSTD.MET	1	1	1	0	1115199999	
11 WISC41132A	CCAL	GP	L152.11R	WILSTD.MET	1	1	1	0	1115199999	
12 BLANKA 5/23/11	BLK	AA	L152.12R	WILSUM.MET	1000	1	1	0	111430013A	07554
13 LCSA 5/23/11	LCS	AA	L152.13R	WILSUM.MET	1000	1	1	0	111430013A	07554
14 6293915	T	AA	L152.14R	WILSUM.MET	806	1	1	0	111430013A	07554
15 MECL2	MISC	AA	L152.15R	WILSTD.MET	1	1	1	0	1115199999	
16 LCSDA 5/23/11	LCSD	AA	L152.16R	WILSUM.MET	1000	1	1	0	111430013A	07554
17 WISC31132A	CCAL	IO	L152.17R	WILSTD.MET	1	1	1	0	1115199999	
18 FLA_21132A	CCAL	VQ	L152.18R	FLALSTD.MET	1	1	1	0	1115199999	
19 6291357S OF50	T	AC	L152.19R	FLALSUM.MET	25.4	50	1	0	111390029A	02100
20 6291357MSS DF50	MS	AC	L152.20R	FLALSUM.MET	25	50	1	0	111390029A	02100
21 6291357MSDS DF50	MSD	AC	L152.21R	FLALSUM.MET	25.5	50	1	0	111390029A	02100
22 6291358S	T	AB	L152.22R	FLALSUM.MET	25.4	1	1	0	111390029A	02100
23 FLA_31132A	CCAL	RY	L152.23R	FLALSTD.MET	1	1	1	0	1115199999	
24 AKRTX1132A	CCAL	NN	L152.24R	AKRTL.MET	1	1	1	0	1115199999	
25 AKCK31132A	CCAL	NX	L152.25R	AKRLSTD.MET	1	1	1	0	1115199999	
26 6297824DF2	T	AB	L152.26R	AKRLSUM.MET	1042	2	1	0	111460032A	02923
27 AKCK41132A	CCAL	MG	L152.27R	AKRLSTD.MET	1	1	1	0	1115199999	
28 AKRTX1132A	CCAL	NO	L152.28R	AKRTL.MET	1	1	1	0	1115199999	
29 CONDITIONER	MISC	AA	L152.29R	WILSTD.MET	1	1	1	0	1115199999	
30 FLA_41132A	CCAL	UG	L152.30R	FLALSTD.MET	1	1	1	0	1115199999	
31 6291357S DF50	T	AC	L152.31R	FLALSUM.MET	25.4	50	1	0	111390029A	02100
32 6291357MSS DF50	MS	AC	L152.32R	FLALSUM.MET	25	50	1	0	111390029A	02100
33 6291357MSDS DF50	MSO	AC	L152.33R	FLALSUM.MET	25.5	50	1	0	111390029A	02100
34 6291358S	T	AB	L152.34R	FLALSUM.MET	25.4	1	1	0	111390029A	02100
35 FLA_21132A	CCAL	VR	L152.35R	FLALSTD.MET	1	1	1	0	1115199999	
36 FLA_21132A	CCAL	VR	L152.36R	FLALSTD.MET	1	1	1	0	1115199999	
37 AKRTX1132A	CCAL	NP	L152.37R	AKRTL.MET	1	1	1	0	1115299999	
38 AKCK21132A	CCAL	OO	L152.38R	AKRLSTD.MET	1	1	1	0	1115299999	
39 6297824DF5	T	AC	L152.39R	AKRLSUM.MET	1042	5	1	0	111460032A	02923
40 AKCK31132A	CCAL	NY	L152.40R	AKRLSTD.MET	1	1	1	0	1115299999	
41 AKFL21132A	CCAL	SD	L152.41R	AKDLSTD.MET	1	1	1	0	1115299999	
42 BLANKA 6/2/11	BLK	AA	L152.42R	AKDLSUM.MET	1000	1	1	0	111520013A	01741
43 LCSA 6/2/11	LCS	AA	L152.43R	AKDLSUM.MET	1000	1	1	0	111520013A	01741
44 LCSDA 6/2/11	LCSD	AA	L152.44R	AKDLSUM.MET	1000	1	1	0	111520013A	01741
45 6297068R	T	AA	L152.45R	AKDLSUM.MET	993	1	1	0	111520013A	01741
46 6299141	T	AA	L152.46R	AKDLSUM.MET	968	1	1	0	111520013A	01741
47 6299142	T	AA	L152.47R	AKDLSUM.MET	975	1	1	0	111520013A	01741
48 6299143	T	AA	L152.48R	AKDLSUM.MET	987	1	1	0	111520013A	01741
49 6299151	T	AA	L152.49R	AKDLSUM.MET	1004	1	1	0	111520013A	01741
50 6299152	T	AA	L152.50R	AKDLSUM.MET	657	1	1	0	111520013A	01741

Lancaster Laboratories

CHROM PERFECT SEQUENCE FILE

Sequence File: \\cp24\c-Drive\CPWIN\DATA1\L152.seq

Chromatography Directory: \\cp24\c-Drive\CPWIN\DATA1

Method Directory: \\cp24\c-Drive\CPWIN\DATA1

Number of Entries: 85

SampleName	Code	ID	FileName	Method	Samp Amt	DF	Int Std	C	Batch Number	Analysis
51 6299153	T	AA	L152.51R	AKDLSUM.MET	906	1	1	0	111520013A	01741
52 AKFL31132A	CCAL	SI	L152.52R	AKDLSTD.MET	1	1	1	0	1115299999	
53 6299156	T	AA	L152.53R	AKDLSUM.MET	1047	1	1	0	111520013A	01741
54 6301430	T	AA	L152.54R	AKDLSUM.MET	1011	1	1	0	111520013A	01741
55 6301431	T	AA	L152.55R	AKDLSUM.MET	1024	1	1	0	111520013A	01741
56 6299154	T	AA	L152.56R	AKDLSUM.MET	1005	1	1	0	111520013A	01741
57 MECL2	MISC	AA	L152.57R	AKDLSTD.MET	1	1	1	0	1115299999	
58 6299155	T	AA	L152.58R	AKDLSUM.MET	1001	1	1	0	111520013A	01741
59 MECL2	MISC	AA	L152.59R	AKDLSTD.MET	1	1	1	0	1115299999	
60 6299157	T	AA	L152.60R	AKDLSUM.MET	988	1	1	0	111520013A	01741
61 MECL2	MISC	AA	L152.61R	AKDLSTD.MET	1	1	1	0	1115299999	
62 MECL2	MISC	AA	L152.62R	AKDLSTD.MET	1	1	1	0	1115299999	
63 AKFL41132A	CCAL	QO	L152.63R	AKDLSTD.MET	1	1	1	0	1115299999	
64 AKRTX1132A	CCAL	NQ	L152.64R	AKRTL.MET	1	1	1	0	1115299999	
65 TPH_41132F	CCAL	NZ	L152.65R	TNLCK.MET	1	1	1	0	1115299999	
66 TNIC41132A	CCAL	IT	L152.66R	TNLCK.MET	1	1	1	0	1115299999	
67 BLANKA 5/27/11	BLK	AA	L152.67R	TNLSUM.MET	1000	1	1	0	111460031A	02784
68 LCSA 5/27/11	LCS	AA	L152.68R	TNLSUM.MET	1000	1	1	0	111460031A	02784
69 LCSDA 5/27/11	LCSD	AA	L152.69R	TNLSUM.MET	1000	1	1	0	111460031A	02784
70 6297963	T	AA	L152.70R	TNLSUM.MET	1036	1	1	0	111460031A	02784
71 6299103	T	AA	L152.71R	TNLSUM.MET	1037	1	1	0	111460031A	02784
72 BLANKA 6/1/11	BLK	AA	L152.72R	TNLSUM.MET	1000	1	1	0	111520006A	02784
73 LCSA 6/1/11	LCS	AA	L152.73R	TNLSUM.MET	1000	1	1	0	111520006A	02784
74 LCSDA 6/1/11	LCSD	AA	L152.74R	TNLSUM.MET	1000	1	1	0	111520006A	02784
75 6297970	T	AA	L152.75R	TNLSUM.MET	1036	1	1	0	111520006A	02784
76 6300755	T	AA	L152.76R	TNLSUM.MET	1027	1	1	0	111520006A	02784
77 TNIC31132A	CCAL	KE	L152.77R	TNLCK.MET	1	1	1	0	1115299999	
78 TPH_31132F	CCAL	HX	L152.78R	TNLCK.MET	1	1	1	0	1115299999	
79 CONDITIONER	MISC	AA	L152.79R	WILSTD.MET	1	1	1	0	1115199998	
80 AKRTX1132A	CCAL	NR	L152.80R	AKRTL.MET	1	1	1	0	1115399999	
81 AKFL21132A	CCAL	SE	L152.81R	AKDLSTD.MET	1	1	1	0	1115399999	
82 6301430DF25	T	AB	L152.82R	AKDLSUM.MET	1011	25	1	0	111520013A	01741
83 6301431DF50	T	AB	L152.83R	AKDLSUM.MET	1024	50	1	0	111520013A	01741
84 AKFL31132A	CCAL	SJ	L152.84R	AKDLSTD.MET	1	1	1	0	1115399999	
85 AKRTX1132A	CCAL	NS	L152.85R	AKRTL.MET	1	1	1	0	1115399999	

30/ASG
6/28/11

AKES9 8181

Set-up by:

Kathleen Villalobos

Date:

6/6/11

6/6/11

Page 2 of 2

Sample Data

SDG: AKE89**Fraction: TPH-DRO by GC**

01741: TPH-DRO AK water C10-C25 Analyte Name	Default MDL	Default LOQ	Units
TPH-DRO AK water C10-C25	0.050	0.25	mg/l

AKE89 0103

Lancaster Laboratories-Range Data Summary

Sample Name: 6301430DF25 **CJMW5** **Sample ID:** AB **Batchnumber:** 111520013A
Sample Amount: 1011. **Total Volume:** 25. ml **Analyst:** 2027 **SDG:** AKE89 **State:** AK
Analyses: 01741

Injection Summary

Injected on : 6/3/2011 22:10:05
Instrument : CP24-H5386A
Result file : L152.82R
Calibration files : AKDL047B.CAL
Method files : AKDLSUM.MET REAKDL.MET
Setting : AKDL047B

Surrogate Recoveries

O-TERPHENYL SURR 106.5% (50-150) **Conc.:** 12.646

Range	Retention Times	Area	Amount	LOQ	MDL	Flags	Units
<input type="checkbox"/> C10-<C25 DRO	2.70 - 12.23	15351820	18.5195	6.182	1.2364		ppm
<input type="checkbox"/> o-Terphenyl SURR	10.05 (10.04 - 10.14)	12163	12.6460				ppb

Comments: _____

Tracy A. Cole

Reviewed by: CJMW5P

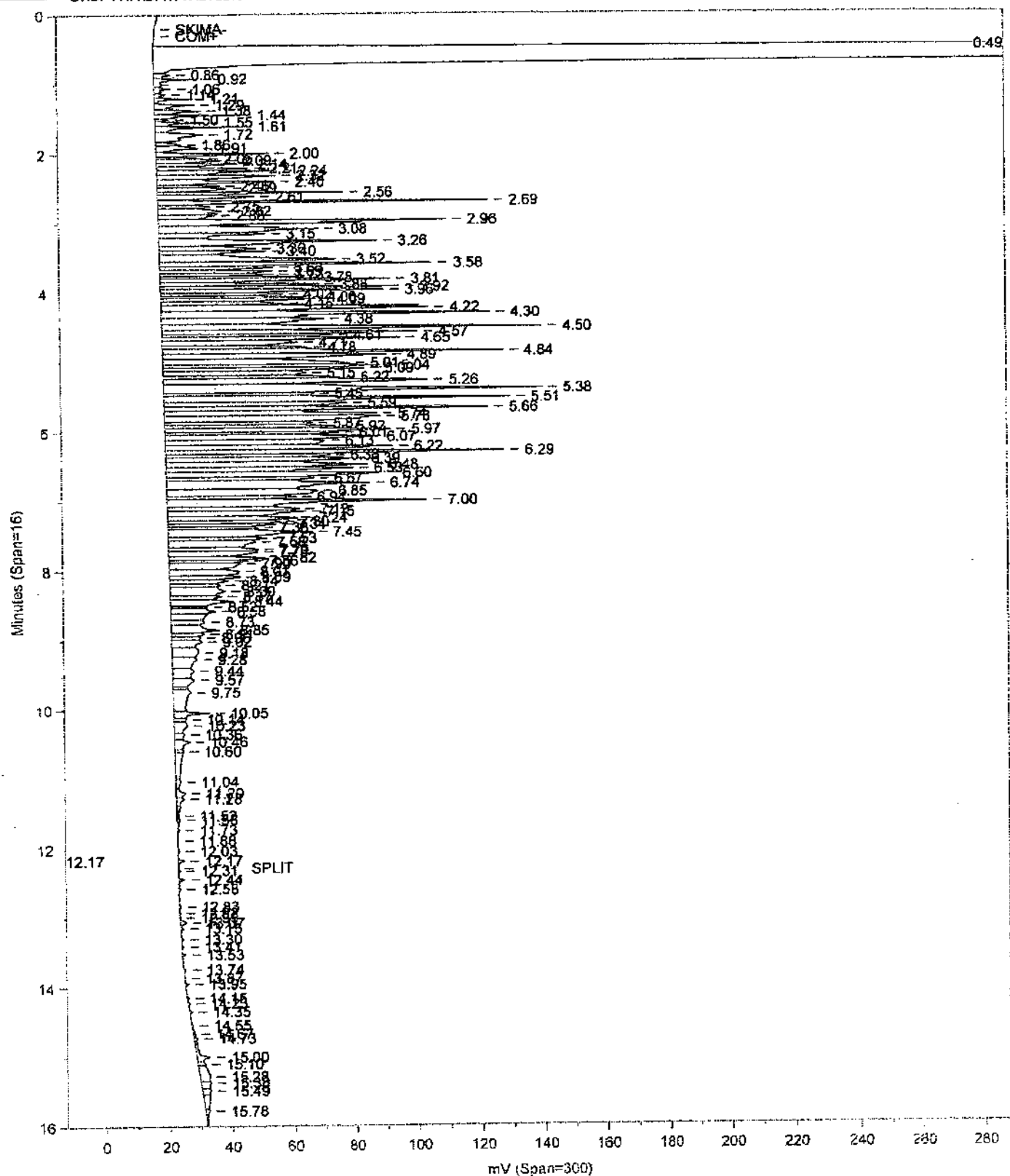
Verified by: _____

Date: 6/16/11

Date: JUN 06 2011

Tracy A. Cole
Senior Specialist

AKES89 0104



Instrument ID: CP24-H5386A

Volume Inj. per Column: 1

Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN

Sample Amount: 1011

Injected on: 6/3/2011 10:10:04 PM

GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C; DET 320C

Dilution Factor: 25

C:\CPWINDATA\1\152.82R

Printed on 6/3/2011 10:26:15 PM

Sample ID: 6301430DF25 ABCJMW5 T 111520013A 01741
 Instrument ID: CP24-H5386A Injected on: 6/3/2011 10:10:04 PM
 Volume Inj. per Column: 1 GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C
 Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN
 Sample Amount: 1011 Dilution Factor: 25
 Analyst: 2027

Peak #	Ret Time (min)	Peak Name	Amount PPB	Peak Area	Peak Width (min)	Peak Height
30	2.819	C10	.	97767	.073	20384
127	10.05	o-Terphenyl SURR	36.0518	34674	.023	11561
143	12.312	C25	.	1663	.035	656

Slice	Start Time	Stop Time	Slice Amount	Amount %	Slice Area	Area %
1	2.700	12.230	36.448	100.000	15351820.0	51.889
2	10.040	10.140	36.448	100.000	45930.4	0.155

Total slice amount= 72.897
 Total slice amount %= 200.0

Total slice area= 15397750.0
 Total slice area %= 52.0

***** RESULTS TABLE *****

C10-<C25 DRO AREA = 1.535182E+07
 C10-<C25 AMT = 18.47876

FILES:

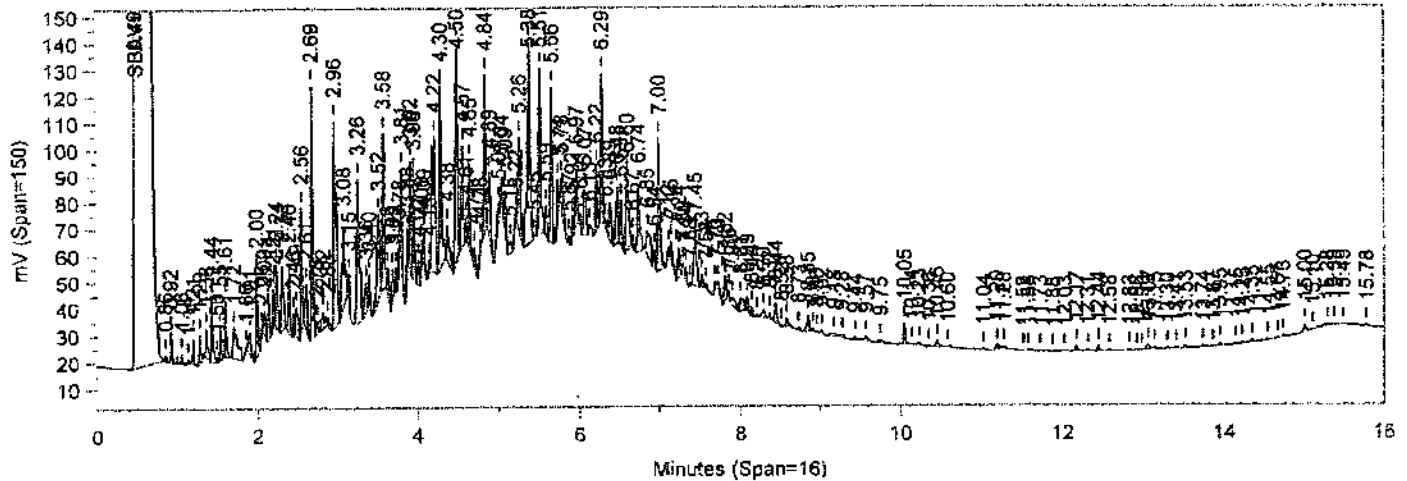
Area File: C:\CPWIN\DATA\1\152.82A
 Method File: C:\CPWIN\DATA\1\AKDLSUM.MET
 Calibration File: C:\CPWIN\DATA\1\AKDL047B.CAL
 Format File: C:\CPWIN\DATA\1\AKDLSUM.FMT
 Area file created on: 6/3/2011 10:26:12 PM
 File reported on: 6/3/2011 at 10:26:14 PM

AK289 8186

AK 102/103

6301430DF25 ABCJMW5 T 111520013A 01741

C:\CPWINDATA\1L152.82R



Sample Name: 6301430DF25 ABCJMW5 T 111520013A 01741A

Instrument ID: CP24--H5386A

Injected on: 6/3/2011 10:10:04 PM

Volume Inj. per Column: 1

GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C

Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN

Sample Amount: 1011

Dilution Factor: 25

Analyst: 2027

Peak #	Ret Time (min)	Peak Name	Amount PPM	Peak Area	Peak Width (min)	Peak Height
30	2.819	C10	.	18938	.073	6567
127	10.05	o-Terphenyl SURR	.0126463	12163	.023	7778
142	12.312	C25	.	1405	.035	608

Slice	Start Time	Stop Time	Slice Amount	Amount %	Slice Area	Area %
-------	------------	-----------	--------------	----------	------------	--------

Total slice amount= 0.000

Total slice area= 0.0

Total slice amount %= 0.0

Total slice area %= 0.0

O-TERPHENYL % RECOVERY = 106.5449 %

FILES:

Area File: C:\CPWINDATA\1L152.82A

Method File: C:\CPWINDATA\1\REAKDL.MET

Calibration File: C:\CPWINDATA\1\AKDL047B.CAL

Format File: C:\CPWINDATA\1\REAKDL.FMT

Area file created on: 6/3/2011 10:26:26 PM

File reported on: 6/3/2011 at 10:26:27 PM

AKES9 0107

Lancaster Laboratories-Range Data Summary

Sample Name: 6301431DF50 **CJDU1** **Sample ID:** AB **Batchnumber:** 111520013A
Sample Amount: 1024. **Total Volume:** 50. ml **Analyst:** 2027 **SDG:** AKE89 **State:** AK
Analyses: 01741

Injection Summary

Injected on : 6/3/2011 22:37:25
 Instrument : CP24-H5386A
 Result file : L152.83R
 Calibration files : AKDL047B.CAL
 Method files : AKDLSUM.MET REAKDL.MET
 Setting : AKDL047B

Surrogate Recoveries

O-TERPHENYL SURR 101.5% (50-150) Conc.: 11.892

Range	Retention Times	Area	Amount	LOQ	MDL	Flags	Units
<input type="checkbox"/> C10-<C25 DRO	2.70 - 12.23	9980843	23.7799	12.207	2.4414		ppm
<input type="checkbox"/> o-Terphenyl SURR	10.05 (10.04 - 10.14)	5792	11.8920				ppb

Comments: _____

Tracy A. Cole

Reviewed by: UWSP

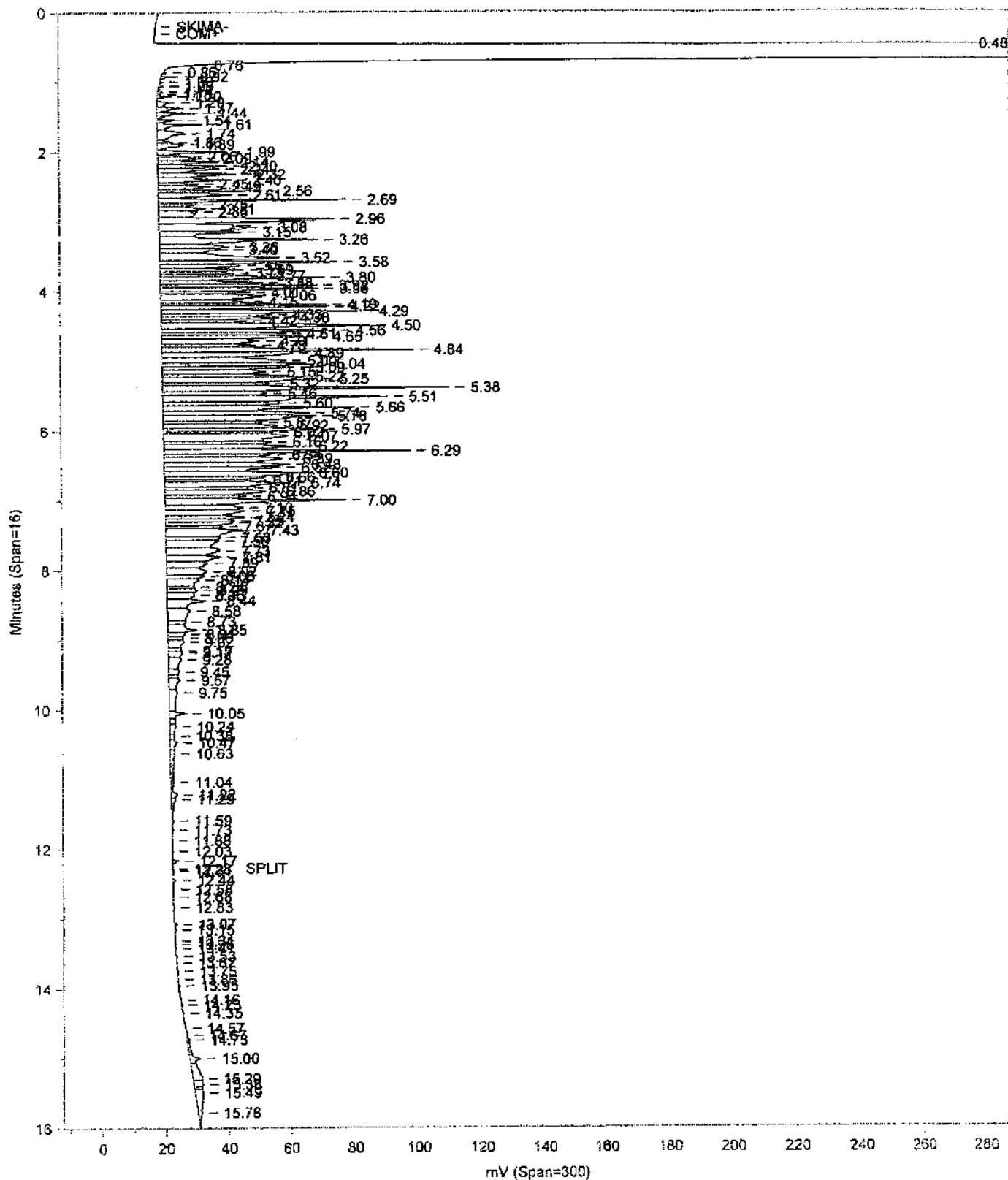
Verified by: _____

Date: 6/6/11

Date: JUN 06 2011

Tracy A. Cole
 Senior Specialist

AKE89 8183



Instrument ID: CP24-H5386A

Volume Inj. per Column: 1

Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN

Sample Amount: 1024

Injected on: 6/3/2011 10:37:24 PM

GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C

Dilution Factor: 50

Sample ID: 6301431DF50 ABCJDU1 T 111520013A 01741
 Instrument ID: CP24-H5386A Injected on: 6/3/2011 10:37:24 PM
 Volume Inj. per Column: 1 GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C
 Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN
 Sample Amount: 1024 Dilution Factor: 50
 Analyst: 2027

Peak #	Ret Time (min)	Peak Name	Amount PPB	Peak Area	Peak Width (min)	Peak Height
32	2.815	C10	.	42871	.03	14917
132	10.051	o-Terphenyl SURR	35.8623	17468	.024	5303
146	12.313	C25	.	1075	.037	407

Slice	Start Time	Stop Time	Slice Amount	Amount %	Slice Area	Area %
1	2.700	12.230	36.723	100.000	9980845.0	89.560
2	10.040	10.140	36.723	100.000	17467.7	0.157

Total slice amount= 73.446
 Total slice amount %= 200.0

Total slice area= 9998313.0
 Total slice area %= 89.7

***** RESULTS TABLE *****

C10-<C25 DRO AREA = 9980845
 C10-<C25 AMT = 23.75209

FILES:

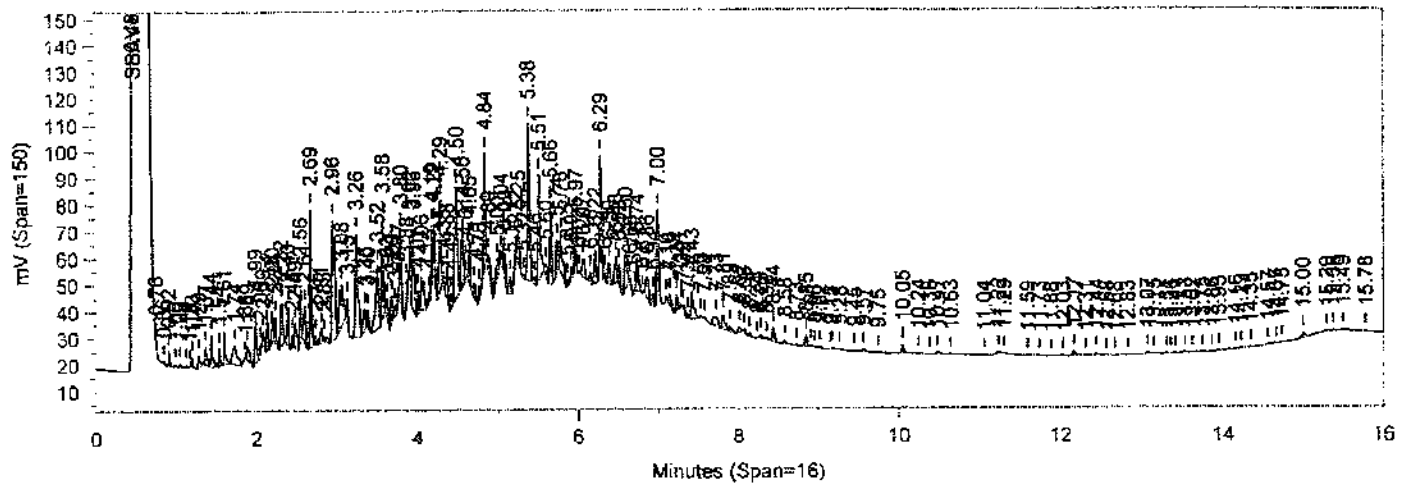
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 Method File: C:\CPWIN\DATA1\AKDLSUM.MET
 Calibration File: C:\CPWIN\DATA1\AKDL047B.CAL
 Format File: C:\CPWIN\DATA1\AKDLSUM.FMT
 Area file created on: 6/3/2011 10:53:32 PM
 File reported on: 6/3/2011 at 10:53:34 PM

AK889 0118

AK 102/103

6301431DF50 ABCJDU1 T 111520013A 01741

C:\CPWIN\DATA\1\152.83R



Sample Name: 6301431DF50 ABCJDU1 T 111520013A 01741A

Instrument ID: CP24--H5386A

Injected on: 6/3/2011 10:37:24 PM

Volume Inj. per Column: 1

GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C

Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN

Sample Amount: 1024

Dilution Factor: 50

Analyst: 2027

Peak #	Ret Time (min)	Peak Name	Amount PPM	Peak Area	Peak Width (min)	Peak Height
32	2.815	C10	.	9151	.03	5342
132	10.051	o-Terphenyl SURR	.0118916	5792	.024	3466
145	12.313	C25	.	934	.037	375

Slice	Start Time	Stop Time	Slice Amount	Amount %	Slice Area	Area %
-------	------------	-----------	--------------	----------	------------	--------

Total slice amount= 0.000

Total slice area= 0.0

Total slice amount %= 0.0

Total slice area %= 0.0

O-TERPHENYL % RECOVERY = 101.4753 %

FILES:

Area File: C:\CPWIN\DATA\1\152.83A

Method File: C:\CPWIN\DATA\1\REAKDL.MET

Calibration File: C:\CPWIN\DATA\1\AKDL047B.CAL

Format File: C:\CPWIN\DATA\1\REAKDL.FMT

Area file created on: 6/3/2011 10:53:46 PM

File reported on: 6/3/2011 at 10:53:48 PM

AK102 8111

Raw QC Data

Lancaster Laboratories-Range Data Summary

Sample Name: BLANKA 6/2/11 **PBLK13152** **Sample ID:** AA **Batchnumber:** 111520013A
Sample Amount: 1000. **Total Volume:** 1. ml **Analyst:** 2027 **SDG:** **State:**
Analyses: 01741

Injection Summary

Injected on : 6/2/2011 18:43:38
Instrument : CP24-H5386A
Result file : L152.42R
Calibration files : AKDL047B.CAL
Method files : AKDLSUM.MET REAKDL.MET
Setting : AKDL047B

Surrogate Recoveries

O-TERPHENYL SURR 101.6% (50-150) **Conc.:** 12.195

Range	Retention Times	Area	Amount	LOQ	MDL	Flags	Units
<input type="checkbox"/> C10-<C25 DRO	2.70 - 12.23	843835	0.0270	<0.25	<0.05		ppm
<input type="checkbox"/> o-Terphenyl SURR	10.06 (10.04 - 10.14)	290043	12.1950				ppb

Comments: _____

Tracy A. Cole

Reviewed by: WMSD

Verified by: JUN 06 2011

Date: 6/3/11

Date: _____

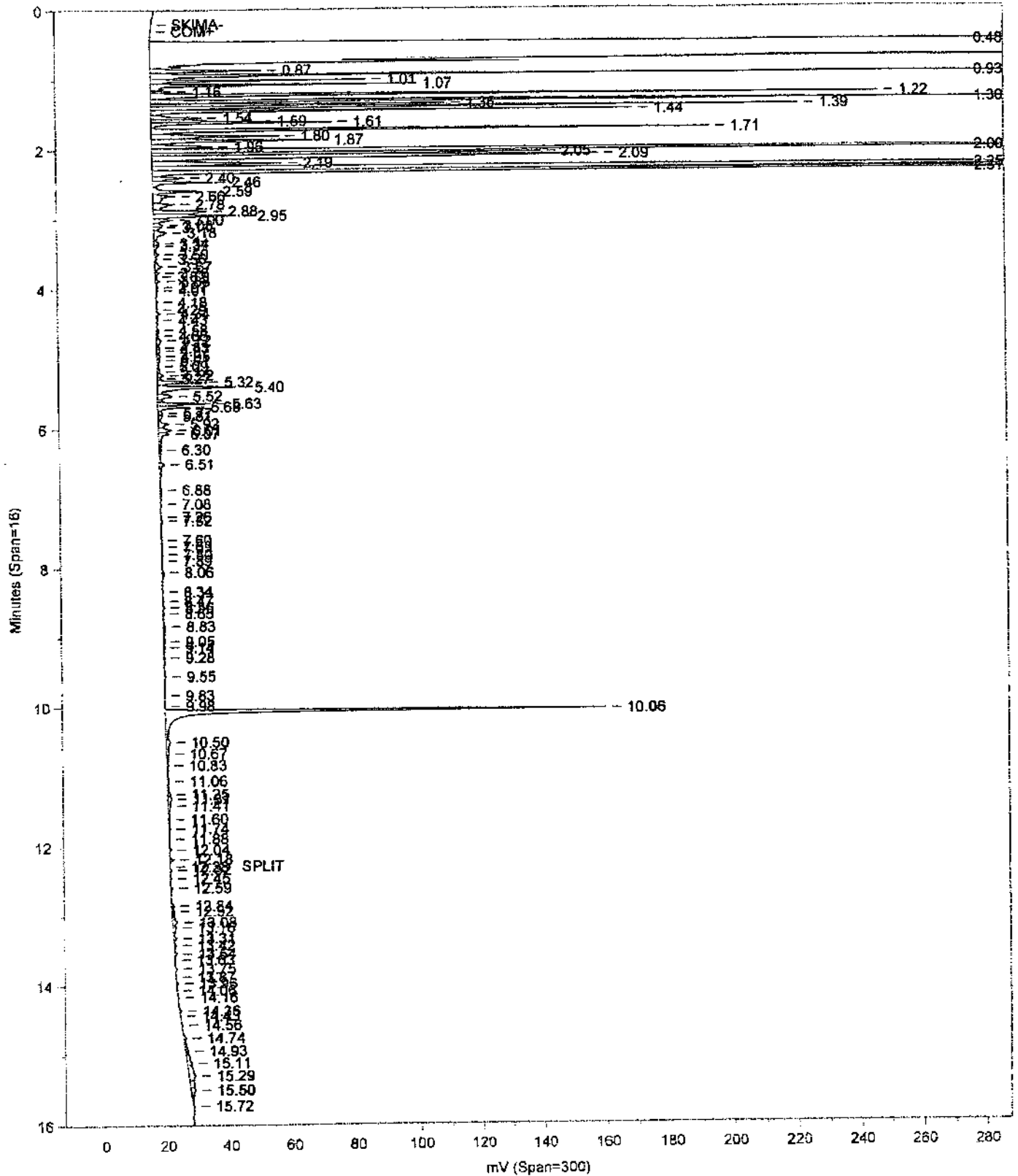
Tracy A. Cole
Senior Specialist

6/3/11 13:31

AK 102/103

BLANKA 6/2/11 AAPBLK13152 BLK 111520013A 01741

C:\CPWIN\DATA\1\152.42R



Instrument ID:CP24-H5386A

Volume Inj. per Column: 1

Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN

Sample Amount: 1000

Injected on: 6/2/2011 6:43:37 PM

GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C

AKES9 0114

Dilution Factor: 1

C:\CPWIN\DATA\1\152.42R

Printed on 6/2/2011 6:59:48 PM

AK 102/103

Sample ID: BLANKA 6/2/11 AAPBLK13152 BLK 111520013A 01741
Instrument ID: CP24--H5386A Injected on: 6/2/2011 6:43:37 PM
Volume Inj. per Column: 1 GC Column: ZB-5 30M x 0.32mm x 0.25um INJ 300C: DET 320C
Oven Parameters: 50C 1MIN; 15C/MIN TO 180C; 30C/MIN TO 340C; HOLD 1MIN
Sample Amount: 1000 Dilution Factor: 1
Analyst: 2027

Peak #	Ret Time (min)	Peak Name	Amount PPB	Peak Area	Peak Width (min)	Peak Height
29	2.778	C10	.	21841	.039	7002
93	10.057	o-Terphenyl SURR	12.5612	298743	.026	140651
107	12.324	C25	.	1444	.037	608

Slice	Start Time	Stop Time	Slice Amount	Amount %	Slice Area	Area %
1	2.700	12.230	12.561	100.000	843834.8	4.142
2	10.040	10.140	12.561	100.000	298743.1	1.466

Total slice amount= 25.122
Total slice amount %= 200.0

Total slice area= 1142578.0
Total slice area %= 5.6

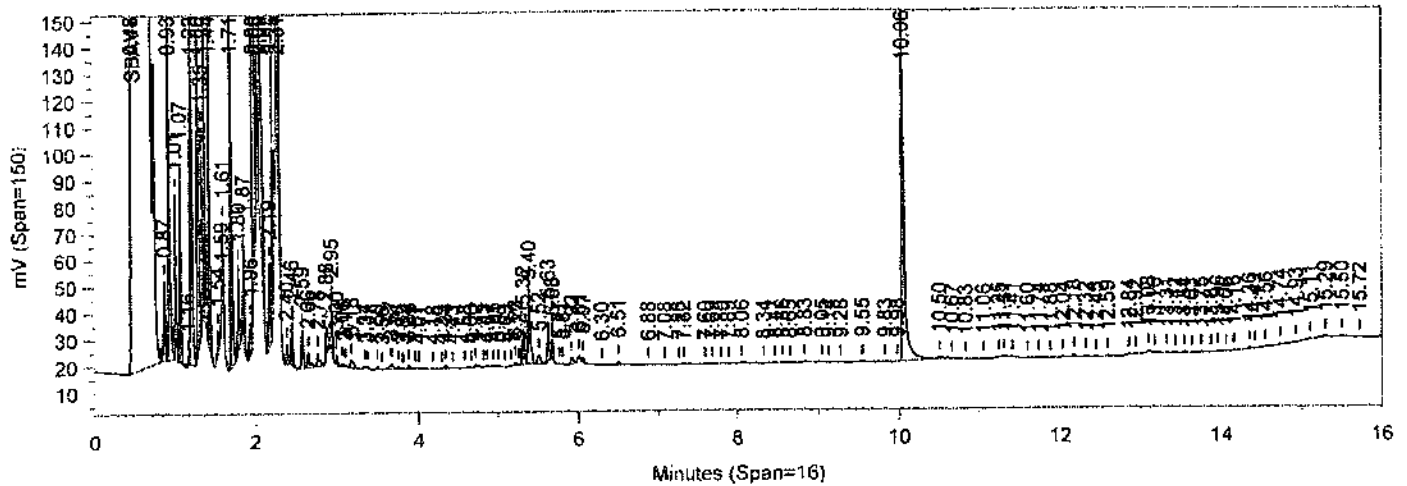
***** RESULTS TABLE *****

C10-<C25 DRO AREA = 843834.8
C10-<C25 AMT = 2.661307E-02

FILES:

Area File: C:\CPWIN\DATA\1\152.42A
Method File: C:\CPWIN\DATA\1\AKDLSUM.MET
Calibration File: C:\CPWIN\DATA\1\AKDL047B.CAL
Format File: C:\CPWIN\DATA\1\AKDLSUM.FMT
Area file created on: 6/2/2011 6:59:46 PM
File reported on: 6/2/2011 at 6:59:47 PM

AK 102/103



Extraction/Distillation/Digestion Logs

Organic Extraction Batchlog

Assigned to: 1785 Sherry Morrow

Reviewed by: LL/11/2011Start Date: 6/2/11Start time: 3:00 PM**111520013A**Tech 1: SLM 1785

Tech 2: _____

Dept: 32 Prep Analysis: 11184 AK DRO Waters Extraction TPH-DRO AK water C10-C25

Solvent Used	Lot No.
1:1 HCl	9026-10
Methylene Chloride	111867
Sodium Sulfate	11153A

QC	Sample Code	Amt (mL)	SS/IS Sol.	Amt (mL)	MS Sol.	Amt (mL)	FV (mL)	pH	pH	BC	Comments
BLANKA	PBLK13152	1000	SS1113732A	1.0			1.0	2	7	N/A	DI H2O
LCSA	LCS13152	1000	SS1113732A	1.0	MS1113632B		1.0	2	7	N/A	DI H2O
LCSDA	LCS13152	1000	SS1113732A	1.0	MS1113632B		1.0	2	7	N/A	DI H2O

Spike Solutions: _____

Witness: N/A

MS1113632B DRO WATER SPIKE

SS1114732A + SS1114732A DRO WATER SURROGATE

Sample #	Sample Code	Amt (mL)	SS/IS Sol.	Amt (mL)	FV (mL)	pH	BC	Comments	Analyses	Due Date	Prio
1	6297068 R	993	SS1113732A	1.0	1.0	2	29A	Cloudy LIGHT GRAY	01741	06/02/2011	P
2	6299141	918	SS1113732A	1.0	1.0	2	29A		01741	06/05/2011	P
3	6298142	975	SS1113732A	1.0	1.0	2	29A		01741	06/06/2011	P
4	6299143	987	SS1113732A	1.0	1.0	2	29A	Cloudy LIGHT BROWN	01741	06/06/2011	P
5	6299151	1004	SS1113732A	1.0	1.0	2	29A	Cloudy	01741	06/06/2011	P
6	6299152	857	SS1113732A	1.0	1.0	2	29A	Cloudy Tan	01741	06/06/2011	P
7	6299153	996	SS1113732A	1.0	1.0	2	29A	Brown LIGHT SEGMENT	01741	06/06/2011	P
8	6299154	1005	SS1113732A	1.0	1.0	2	29A	Cloudy Tan	01741	06/06/2011	P
9	6299155	1601	SS1113732A	1.0	1.0	2	29A	Brown LIGHT SEGMENT	01741	06/06/2011	P
10	6299156	1047	SS1113732A	1.0	1.0	2	29A	Clear	01741	06/06/2011	P
11	6299157	948	SS1113732A	1.0	1.0	2	29A	Brown LIGHT SEGMENT	01741	06/06/2011	P
12	6301430	1011	SS1113732A	1.0	1.0	2	29A	Cloudy	01741	06/08/2011	P
13	6301431	1024	SS1113732A	1.0	1.0	2	29A	Cloudy	01741	06/08/2011	P

Sum 785 6/2/11

Rack ID:

Internal Standard

Work Station

H2O 5

Balance #

—

DF = Dilution Factor FV = Final Volume

Page 1 of 1

S-bath ID 11185 C S-bath ID C N-Evap C M-vap C

111520013A

Documented temps are NIST corrected.

Instrumental Wet Chemistry Data

Case Narrative Conformance/Nonconformance Summary



CLIENT: ChevronTexaco
SDG: AKE89

Instrumental Water Quality

Sample #	Matrix		Comments
	Liquid	Solid	
6301430	X		

ANALYSIS:

Dilutions are listed in the table below:

Samples	Nitrate Nitrogen	Nitrite Nitrogen	Sulfate
6301430			DF5
LCS	DF10	DF5	

QUALITY CONTROL AND NONCONFORMANCE SUMMARY:

Site specific matrix QC samples were not submitted for this SDG. The batch matrix QC was performed on samples from another project. Therefore the matrix effects would not be relevant and matrix QC is not provided in the data package. Laboratory spike data (LCS) are provided.

Method defined actions are taken for any failed matrix QC.

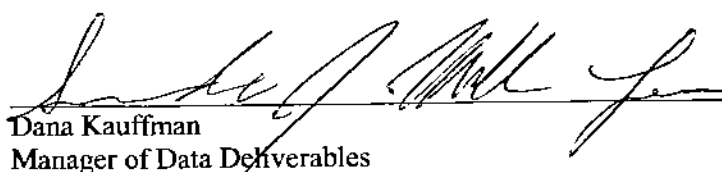
DATA INTERPRETATION:

No further interpretation is necessary for the data submitted.

Abbreviation Key

U = Unspiked (for MS/MSD)	LOQ = Limit of Quantitation
R = Matrix Spike (MS)	MDL = Method Detection Limit
M = Matrix Spike Duplicate (MSD)	ND = Not Detected
BKG = Background (for Duplicate)	J = Estimated Value
D = Duplicate (DUP)	NA = Not Applicable
HS = High Spike	ME = Method
LS = Low Spike	CO = Colorimetric
SS = Soluble Spike	G = Gravimetric
IS = Insoluble Spike	IR = Infrared Spectrophotometry
ISD = Insoluble Spike Duplicate	MTR = Meter
PDS = Post Digestion Spike	OD = Oven Dried
* = Out of Specification	TI = Titration
V = Visual	TOC = Total Organic Carbon
AK = Alpkem	IC = Ion Chromatography
TC = Total Carbon	RA = Rapid Analyzer

Narrative Reviewed and Approved by:


Dana Kauffman
Manager of Data Deliverables

AKE89 8121
Date 6/22/17

Quality Control and Calibration Summary Forms



Quality Control Reference List
Instrumental Water Quality

CLIENT: ChevronTexaco
SDG: AKE89

Analyte	Batch Number	Sample Number
Nitrate Nitrogen	11152106102A	6301430 Blank LCS
Nitrite Nitrogen	11148105101A	6301430 Blank LCS
Sulfate	11152196901A	Blank LCS/LCSD
Sulfate	11152196901C	6301430

Analyte	Analysis Date	Method	Batch Number	Blank Results	Units	MDL	LOQ
Nitrate Nitrogen	06/01/11	AK	11152106102	N.D.	mg/l	0.040	0.10
Nitrite Nitrogen	05/28/11	AK	11148105101	N.D.	mg/l	0.015	0.050
Sulfate	06/02/11	IC	11152196901	N.D.	mg/l	0.30	1.0

AKE89 6124

Comments: The blank is acceptable when the result is less than the limit of quantitation.



Quality Control Summary
Laboratory Control Standard (LCS)
Laboratory Control Standard Duplicate (LCSD)
Instrumental Water Quality
SDG: AKE89
Matrix: LIQUID

Batch #	Analyte	Analysis Date	ME	True LCS/LCSD Value	LCS Results	LCSD Results	Units	Acceptance Range	% RPD Results	% RPD Acceptance <=
11152106102	Nitrate Nitrogen	06/01/11	AK	21	22.2	NA	mg/l	18.8 - 23.2	NA	NA
11148105101	Nitrite Nitrogen	05/28/11	AK	1.92	2.0	NA	mg/l	1.719 - 2.12	NA	NA
11152196901	Sulfate	06/02/11	IC	7.5	8.1	8.1	mg/l	6.7125 - 8.28	1	20

AKE89 8125



SDG: AKE89

Instrument ID: 09106
Calibration Date: 05/28/2011

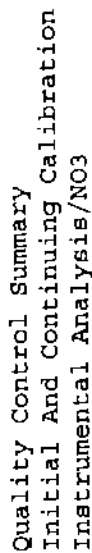
Acceptance Range:
ICV/CCV: +/- 10%
ICB/CCB: < MDL

Concentration units: mg/L

Batch Numbers: 11148105101A
Run Start Dates: 05/28/2011
Run Names: 1114801C02

Analysis	AUTO CAL1	AUTO CAL2	AUTO CAL3	AUTO CAL4	AUTO CAL5	AUTO CAL6	CC
Nitrite-N	2.2425	1.9003	1.1204	0.7066	0.4928	0.4218	0.9996

Sample	Nitrite-N		
	True	Result	%Rec
ICV	0.6	0.60113	100
ICB	0	ND	NA
CCV2	0.6	0.59619	99
CCB 1	0	ND	NA
CCV2	0.6	0.62450	104
CCB 5	0	ND	NA
CCV2	0.6	0.62395	104
CCB 6	0	ND	NA



Quality Control Summary
Initial And Continuing Calibration
Instrumental Analysis/NO3

Analysis	AUTO CAL1	AUTO CAL2	AUTO CAL3	AUTO CAL4	AUTO CAL5	AUTO CAL6	CC
Nitrate-N	3.9274	3.0885	1.2856	0.8103	0.5094	0.4182	0.9998

11153106101B

Sample	Nitrate-N		
	True	Result	%Rec
ICV	2.5	2.56886	103
ICB	0	ND	NA
CCV2	2.5	2.58526	103
CCB 1	0	ND	NA



SDG: AKE89

Instrument ID: 17694
Calibration Date: 05/11/2011

Acceptance Range:
ICV/CCV: 90%-110%
ICB/CCB: < MDL

Concentration units: mg/L

Batch Numbers: 11152196901A, 11152196901C
Run Start Dates: 06/01/2011
Run Names: 1115201D09

Analysis	AUTO CAL1	AUTO CAL2	AUTO CAL3	AUTO CAL4	AUTO CAL5	R ²	CC
Sulfate	0.096547	0.191227	0.463002	0.936085	1.426662	0.9998	0.9999

Sample	Sulfate		
	True	Result	%Rec
ICV	7.5	7.4580	99
ICB	0	ND	NA
CCV2	7.5	7.9866	106
CCB	0	ND	NA
CCV2	7.5	8.0540	107
CCB	0	ND	NA
CCV2	7.5	8.0915	108
CCB	0	ND	NA
CCV2	7.5	8.2479	110
CCB	0	ND	NA

1115201D09

SDG: AKE89

Parameter	Default MDL	Default LOQ	Units
Nitrate Nitrogen	0.040	0.10	mg/l
Nitrite Nitrogen	0.015	0.050	mg/l
Sulfate	0.30	1.0	mg/l

AKE89 8129

Wet Chemistry Data

Case Narrative Conformance/ Non-Conformance Summary



CLIENT: ChevronTexaco
SDG: AKE89

Miscellaneous Wet Chemistry

<u>Sample #</u>	<u>Matrix</u>		<u>Comments</u>
	<u>Liquid</u>	<u>Solid</u>	
6301430	X		

ANALYSIS:

There were no dilutions performed for analyses associated with samples in this SDG.

QUALITY CONTROL AND NONCONFORMANCE SUMMARY:

Site specific matrix QC samples were not submitted for this SDG. The batch matrix QC was performed on samples from another project. Therefore the matrix effects would not be relevant and matrix QC is not provided in the data package. Laboratory spike data (LCS) are provided.

Method defined actions are taken for any failed matrix QC.

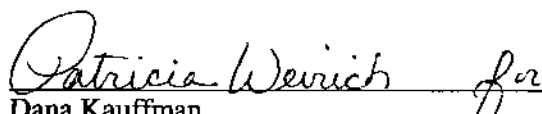
DATA INTERPRETATION:

No further interpretation is necessary for the data submitted.

Abbreviation Key

U = Unspiked (for MS/MSD)	LOQ = Limit of Quantitation
R = Matrix Spike (MS)	MDL = Method Detection Limit
M = Matrix Spike Duplicate (MSD)	ND = Not Detected
BKG = Background (for Duplicate)	J = Estimated Value
D = Duplicate (DUP)	NA = Not Applicable
HS = High Spike	ME = Method
LS = Low Spike	CO = Colorimetric
SS = Soluble Spike	G = Gravimetric
IS = Insoluble Spike	IR = Infrared Spectrophotometry
ISD = Insoluble Spike Duplicate	MTR = Meter
PDS = Post Digestion Spike	OD = Oven Dried
* = Out of Specification	TI = Titration
V = Visual	TOC = Total Organic Carbon
AK = Alpkem	IC = Ion Chromatography
TC = Total Carbon	RA = Rapid Analyzer

Narrative Reviewed and Approved by:


Dana Kauffman
Manager of Data Deliverables

Date 6-16-11

AKE89 6132

QC Summary



Quality Control Reference List
Miscellaneous Wet Chemistry

CLIENT: ChevronTexaco
SDG: AKE89

Batch Number	Sample Number	Alkalinity to pH 4.5	Alkalinity to pH 8.3
11153020201A	Blank	X	
	LCS	X	
11153020201B	6301430	X	X

Analyte	Analysis Date	Method	Batch Number	Blank Results	Units	MDL	LOQ
Alkalinity to pH 4.5	06/02/11	TI	11153020201	N.D.	mg/l as CaCO ₃	0.46	2.0

AKE89 8135

Comments: The blank is acceptable when the result is less than the limit of quantitation.



Quality Control Summary
Laboratory Control Standard (LCS)
Laboratory Control Standard Duplicate (LCSD)
Miscellaneous Wet Chemistry
SDG: AKE89
Matrix: LIQUID

Batch #	Analyte	Analysis Date	ME	True LCS/LCSD Value	LCS Results	LCSD Results	Units	Acceptance Range	% RPD Results	% RPD Acceptance <=
11153020201	Alkalinity to pH 4.5	06/02/11	TI	188	186	NA	mg/l as CaCO3	183.3 - 194.56	NA	NA

AKE89 0136

SDG: AKE89

Parameter	Default MDL	Default LOQ	Units
Alkalinity to pH 4.5	0.46	2.0	mg/l as CaCO ₃
Alkalinity to pH 8.3	0.46	2.0	mg/l as CaCO ₃

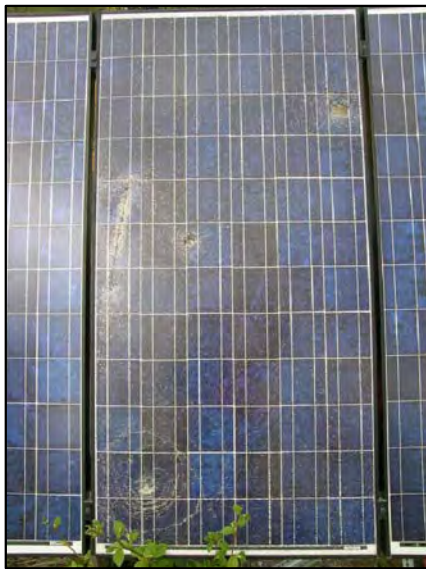
AKE89 6137

ATTACHMENT C

SITE PHOTOS



1. View of MW-4 and ozone injection system looking west.



2. Damaged solar panel.



3. New solar panel.

Appendix A

SITE PHOTOGRAPHS
FORMER DELTA WESTERN/CHEVRON BULK TERMINAL 8-2307
9203 CESSNA DRIVE
Juneau, Alaska



ATTACHMENT D

STANDARD OPERATION PROCEDURES
FOR GROUNDWATER MONITORING AND SAMPLING



**CONESTOGA-ROVERS
& ASSOCIATES**

STANDARD FIELD PROCEDURES FOR GROUNDWATER MONITORING AND SAMPLING

This document presents standard field methods for groundwater monitoring, purging and sampling, and well development. These procedures are designed to comply with Federal, State and local regulatory guidelines. Conestoga-Rovers & Associates' specific field procedures are summarized below.

Groundwater Monitoring

Prior to performing monitoring activities, the historical monitoring and analytical data of each monitoring well shall be reviewed to determine if any of the wells are likely to contain separate phase hydrocarbons (SPH) and to determine the order in which the wells will be monitored (i.e. cleanest to dirtiest). Groundwater monitoring should not be performed when the potential exists for surface water to enter the well (i.e. flooding during a rainstorm).

Prior to monitoring, each well shall be opened and the well cap removed to allow water levels to stabilize and equilibrate. The condition of the well box and well cap shall be observed and recommended repairs noted. Any surface water that may have entered and flooded the well box should be evacuated prior to removing the well cap. In wells with no history of SPH, the static water level and total well depth shall be measured to the nearest 0.01 foot with an electronic water level meter. Wells with the highest contaminant concentrations shall be monitored last. In wells with a history of SPH, the SPH level/thickness and static water level shall be measured to the nearest 0.01 foot using an electronic interface probe. The water level meter and/or interface probe shall be thoroughly cleaned and decontaminated at the beginning of the monitoring event and between each well. Monitoring equipment shall be washed using soapy water consisting of Liqui-noxTM or AlconoxTM followed by one rinse of clean tap water and then two rinses of distilled water.

Groundwater Sampling

Purge Sampling

Prior to groundwater purging and sampling, the historical analytical data of each monitoring well shall be reviewed to determine the order in which the wells should be purged and sampled (i.e. cleanest to dirtiest). No purging or groundwater sampling shall be performed on wells with a measurable thickness of SPH or floating SPH globules. If a sheen is observed, the well should be purged and a groundwater sample collected only if no SPH is present. Wells shall be purged either by hand using a disposal or PVC bailer or by using an aboveground pump (e.g. peristaltic or WatteraTM) or down-hole pump (e.g. GrundfosTM or DC Purger pump).

Groundwater wells shall be purged approximately three to ten well-casing volumes (depending on the regulatory agency requirements) or until a minimum of three groundwater parameters have stabilized for three consecutive readings. Temperature, dissolved oxygen (DO), pH, conductivity, and oxidation-reduction potential (ORP) shall be measured and recorded. The total volume of groundwater removed shall be recorded along with any other notable physical characteristic such as color and odor. If required, turbidity shall also be measured prior to collection of each groundwater sample.

Groundwater samples shall be collected after the well has been purged and allowed to recharge to 80% of the pre-purging static water level, or if the well is slow to recharge, after waiting a minimum of 2 hours. Groundwater samples shall be collected using clean disposable bailers or



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pumps (if an operating remediation system exists on site and the project manager approves of its use for sampling) and shall be decanted into clean containers supplied by the analytical laboratory. New nitrile gloves and disposable tubing or bailers shall be used for sampling each well. If a PVC bailer or down-hole pump is used for groundwater purging, it shall be decontaminated before purging each well by using soapy water consisting of Liqui-nox™ or Alconox™ followed by one rinse of clean tap water and then two rinses of distilled water. If a submersible pump with non-dedicated discharge tubing is used for groundwater purging, both the inside and outside of pump and discharge tubing shall be decontaminated as described above.

No Purge Sampling

Groundwater samples shall be collected using clean disposable bailers (PVC or Teflon) and decanted into clean containers supplied by the analytical laboratory. New nitrile gloves shall be used for sampling each well.

HydraSleeve Sampling

HydraSleeve sampling devices shall be deployed a minimum of 24 hours prior to sampling to allow for equilibration and stabilization.

For wells without a previously deployed HydraSleeve sampling device, the groundwater monitoring SOP shall be followed to determine water column length. For wells with a previously deployed HydraSleeve sampling device, the depth to groundwater shall be measured prior to HydraSleeve removal. The depth to well bottom shall be measured prior to new HydraSleeve deployment.

The top of the HydraSleeve sampling device shall be set no more than 3 feet below static groundwater level. If the length of the water column does not allow for this, a top weight bottom set will be employed. Groundwater samples collected using clean disposable HydraSleeve sampling devices will be decanted into clean containers supplied by the analytical laboratory. New nitrile gloves and HydraSleeve sampling devices will be used for sampling each well.

Following sampling, a new HydraSleeve sampling device shall be deployed for the next sampling event if applicable.

Sample Handling

Except for samples that will be tested in the field, or that require special handling or preservation, samples shall be stored in coolers chilled to 4° C for shipment to the analytical laboratory. Samples shall be labeled, placed in protective foam sleeves or bubble wrap as needed, stored on crushed ice at or below 4° C, and submitted under chain-of-custody (COC) to the laboratory. The laboratory shall be notified of the sample shipment schedule and arrival time. Samples shall be shipped to the laboratory within a time frame to allow for extraction and analysis to be performed within the standard sample holding times.

Sample labels shall be filled out using indelible ink and must contain the site name; field identification number; the date, time, and location of sample collection; notation of the type of sample; identification of preservatives used; remarks; and the signature of the sampler. Field identification must be sufficient to allow easy cross-reference with the field datasheet.

All samples submitted to the laboratory shall be accompanied by a COC record to ensure adequate documentation. One copy of the COC shall be kept in the QA/QC file and another copy shall be retained in the project file. Information on the COC shall consist of the project name and



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number; project location; sample numbers; sampler/recorder's signature; date and time of collection of each sample; sample type; analyses requested; name of person receiving the sample; and date of receipt of sample.

Laboratory-supplied trip blanks shall accompany the samples and be analyzed to check for cross-contamination, if requested by the project manager.

Well Development

Wells shall be developed using a combination of groundwater surging and extraction. A surge block shall be used to swab the well and agitate the groundwater in order to dislodge any fine sediment from the sand pack. After approximately ten minutes of swabbing the well, groundwater shall be extracted from the well using a bailer, pump and/or reverse air-lifting through a pipe to remove the sediments from the well. Alternating surging and extraction shall continue until the sediment volume in the groundwater (i.e. turbidity) is negligible, which typically requires extraction of approximately ten well-casing volumes of groundwater. Preliminary well development usually is performed during well installation prior to placing the sanitary surface seal to ensure sand pack stabilization. Well development that is performed after surface seal installation, should occur 72 hours after seal installation to ensure that the cement has had adequate time to set.

Waste Handling and Disposal

Groundwater extracted during development and sampling shall be stored onsite in sealed U.S. DOT H17 55-gallon drums. Each drum shall be labeled with the contents, date of generation, generator identification and consultant contact. If hydrocarbon concentrations in the purged groundwater are below ADEC cleanup levels or the site is in a remote area (pending ADEC approval) groundwater will be discharged to the ground surface, at least 100 feet from the nearest surface water body.

\\\\DEN-S1\\Shared\\Denver\\Alaska\\AK SOP\\CRA Alaska SOP\\AK Groundwater Monitoring and Sampling SOP - CRA.doc

ATTACHMENT E

ADEC LABORATORY DATA REVIEW CHECKLIST AND MEMORANDUM

Laboratory Data Review Checklist

Completed by:

Title: Date:

CS Report Name: Report Date:

Consultant Firm:

Laboratory Name: Laboratory Report Number:

ADEC File Number: ADEC RecKey Number:

1. Laboratory

- a. Did an ADEC CS approved laboratory receive and perform all of the submitted sample analyses?
X Yes ☐ No ☐ NA (Please explain.) Comments:

- b. If the samples were transferred to another “network” laboratory or sub-contracted to an alternate laboratory, was the laboratory performing the analyses ADEC CS approved?
☐ Yes ☐ No X NA (Please explain.) Comments:

2. Chain of Custody (COC)

- a. COC information completed, signed, and dated (including released/received by)?
X Yes ☐ No ☐ NA (Please explain.) Comments:

- b. Correct analyses requested?
X Yes ☐ No ☐ NA (Please explain.) Comments:

3. Laboratory Sample Receipt Documentation

- a. Sample/cooler temperature documented and within range at receipt ($4^{\circ} \pm 2^{\circ} \text{C}$)?
X Yes ☐ No ☐ NA (Please explain.) Comments:

- b. Sample preservation acceptable – acidified waters, Methanol preserved VOC soil (GRO, BTEX, Volatile Chlorinated Solvents, etc.)?

☒ Yes ☐ No ☐ NA (Please explain.)

Comments:

- c. Sample condition documented – broken, leaking (Methanol), zero headspace (VOC vials)?

☐ Yes ☐ No ☐ NA (Please explain.)

Comments:

- d. If there were any discrepancies, were they documented? For example, incorrect sample containers/preservation, sample temperature outside of acceptable range, insufficient or missing samples, etc.?

☐ Yes ☐ No ☒ NA (Please explain.)

Comments:

No discrepancies

- e. Data quality or usability affected? (Please explain.)

Comments:

None

4. Case Narrative

- a. Present and understandable?

☒ Yes ☐ No ☐ NA (Please explain.)

Comments:

- b. Discrepancies, errors or QC failures identified by the lab?

☒ Yes ☐ No ☐ NA (Please explain.)

Comments:

- c. Were all corrective actions documented?

☐ Yes ☐ No ☒ NA (Please explain.)

Comments:

No corrective actions

- d. What is the effect on data quality/usability according to the case narrative?

Comments:

None

5. Samples Results

- a. Correct analyses performed/reported as requested on COC?

☒ Yes ☐ No ☐ NA (Please explain.)

Comments:

b. All applicable holding times met?

☐ Yes ☒ No ☐ NA (Please explain.)

Comments:

Nitrate was analyzed outside of the 48 hour hold time.

c. All soils reported on a dry weight basis?

☐ Yes ☐ No ☒ NA (Please explain.)

Comments:

No soils

d. Are the reported PQLs less than the Cleanup Level or the minimum required detection level for the project?

☒ Yes ☐ No ☐ NA (Please explain.)

Comments:

e. Data quality or usability affected?

Comments:

The nitrate result for sample MW-4 should be considered estimated.

6. QC Samples

a. Method Blank

i. One method blank reported per matrix, analysis and 20 samples?

☒ Yes ☐ No ☐ NA (Please explain.)

Comments:

ii. All method blank results less than PQL?

☒ Yes ☐ No ☐ NA (Please explain.)

Comments:

iii. If above PQL, what samples are affected?

Comments:

No affected samples

iv. Do the affected sample(s) have data flags and if so, are the data flags clearly defined?

☐ Yes ☐ No ☒ NA (Please explain.)

Comments:

No affected samples

v. Data quality or usability affected? (Please explain.)

Comments:

None

b. Laboratory Control Sample/Duplicate (LCS/LCSD)

- i. Organics – One LCS/LCSD reported per matrix, analysis and 20 samples? (LCS/LCSD required per AK methods, LCS required per SW846)

X Yes ☐ No ☐ NA (Please explain.)

Comments:

- ii. Metals/Inorganics – one LCS and one sample duplicate reported per matrix, analysis and 20 samples?

X Yes ☐ No ☐ NA (Please explain.)

Comments:

- iii. Accuracy – All percent recoveries (%R) reported and within method or laboratory limits? And project specified DQOs, if applicable. (AK Petroleum methods: AK101 60%-120%, AK102 75%-125%, AK103 60%-120%; all other analyses see the laboratory QC pages)

☐ Yes X No ☐ NA (Please explain.)

Comments:

The method 300.0 MS had a high sulfate recovery and was performed on a non-project sample.

- iv. Precision – All relative percent differences (RPD) reported and less than method or laboratory limits? And project specified DQOs, if applicable. RPD reported from LCS/LCSD, MS/MSD, and or sample/sample duplicate. (AK Petroleum methods 20%; all other analyses see the laboratory QC pages)

X Yes ☐ No ☐ NA (Please explain.)

Comments:

- v. If %R or RPD is outside of acceptable limits, what samples are affected?

Comments:

None

- vi. Do the affected sample(s) have data flags? If so, are the data flags clearly defined?

☐ Yes ☐ No X NA (Please explain.)

Comments:

No affected samples

- vii. Data quality or usability affected? (Use comment box to explain.)

Comments:

The sample results would not have been impacted, no qualification of the data was deemed necessary.

c. Surrogates – Organics Only

- i. Are surrogate recoveries reported for organic analyses – field, QC and laboratory samples?

X Yes ☐ No ☐ NA (Please explain.)

Comments:

- ii. Accuracy – All percent recoveries (%R) reported and within method or laboratory limits? And project specified DQOs, if applicable. (AK Petroleum methods 50-150 %R; all other analyses see the laboratory report pages)

X Yes ☐ No ☐ NA (Please explain.)

Comments:

- iii. Do the sample results with failed surrogate recoveries have data flags? If so, are the data flags clearly defined?

☐ Yes ☐ No X NA (Please explain.)

Comments:

No failed surrogates

- iv. Data quality or usability affected? (Use the comment box to explain.)

Comments:

None

d. Trip blank – Volatile analyses only (GRO, BTEX, Volatile Chlorinated Solvents, etc.): Water and Soil

- i. One trip blank reported per matrix, analysis and for each cooler containing volatile samples? (If not, enter explanation below.)

X Yes ☐ No ☐ NA (Please explain.)

Comments:

- ii. Is the cooler used to transport the trip blank and VOA samples clearly indicated on the COC? (If not, a comment explaining why must be entered below)

X Yes ☐ No ☐ NA (Please explain.)

Comments:

- iii. All results less than PQL?

X Yes ☐ No ☐ NA (Please explain.)

Comments:

- iv. If above PQL, what samples are affected?

Comments:

No affected samples

- v. Data quality or usability affected? (Please explain.)

Comments:

None

e. Field Duplicate

i. One field duplicate submitted per matrix, analysis and 10 project samples?

X Yes ☐ No ☐ NA (Please explain.)

Comments:

ii. Submitted blind to lab?

X Yes ☐ No ☐ NA (Please explain.)

Comments:

iii. Precision – All relative percent differences (RPD) less than specified DQOs?

(Recommended: 30% water, 50% soil)

$$\text{RPD (\%)} = \text{Absolute value of: } \frac{(R_1 - R_2)}{((R_1 + R_2)/2)} \times 100$$

Where R_1 = Sample Concentration

R_2 = Field Duplicate Concentration

X Yes ☐ No ☐ NA (Please explain.)

Comments:

iv. Data quality or usability affected? (Use the comment box to explain why or why not.)

Comments:

None

f. Decontamination or Equipment Blank (If not used explain why).

☐ Yes ☐ No X NA (Please explain.)

Comments:

Not collected

i. All results less than PQL?

☐ Yes ☐ No X NA (Please explain.)

Comments:

Not collected

ii. If above PQL, what samples are affected?

Comments:

Not collected

iii. Data quality or usability affected? (Please explain.)

Comments:

Not collected

7. Other Data Flags/Qualifiers (ACOE, AFCEE, Lab Specific, etc.)

a. Defined and appropriate?

☒ Yes ☐ No ☐ NA (Please explain.)

Comments:

--



**CONESTOGA-ROVERS
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MEMORANDUM

TO: ADEC REF. NO.: 622237

FROM: Jeffrey Cloud DATE: November 14, 2011

CC: John Riggi Send via E-Mail and U.S. Mail

RE: QA/QC Review
ChevronTexaco Site # 8-2307
Job #1249125
May 2011

INTRODUCTION

Groundwater samples were submitted to Lancaster Laboratories, located in Lancaster, Pennsylvania. Samples were analyzed for the methods requested on the Chain of Custody.

A full Level III data package was received from Lancaster Laboratories. The final results and supporting quality assurance/quality control (QA/QC) data were reviewed. Evaluation of the data was based on information obtained from the Chain of Custody forms, finished report forms, blank data, and spike recoveries.

QA/QC REVIEW

All samples were prepared and/or analyzed within the required holding times with one exception. Nitrate was analyzed outside of the 48 hour hold time. The nitrate result for sample MW-4 should be considered estimated. All samples were properly preserved and cooled after collection.

All appropriate samples and blanks were spiked with surrogate compounds prior to sample preparation and/or analysis in accordance with the organic methods. All surrogate spike recoveries met the associated method criteria indicating adequate analytical efficiency.

Method blanks were prepared and analyzed with the samples for all parameters. All blank results were non-detect for the analytes of interest.

Laboratory control samples (LCS) were analyzed for all parameters. LCS for methods AK102 and 300.0 were analyzed in duplicate. All recoveries were within required control limits showing adequate analytical accuracy and precision.

Matrix spikes (MS) were prepared and analyzed in duplicate for methods AK101, 8015, 353.2, 300.0 and SM 2320 B. MS for methods AK101 and 8015 were analyzed in duplicate. All recoveries were within required control limits showing adequate analytical accuracy and precision with one exception. The method 300.0

MS had a high sulfate recovery and was performed on a non-project sample. The sample results would not have been impacted, no qualification of the data was deemed necessary.

Trip blanks were collected and analyzed with the investigative samples for all volatile parameters. All trip blank results were non-detect for the compounds of interest.

A field duplicate was collected and submitted blind to the laboratory. The sample ID was MW-4 and its duplicate was DUP-1. A comparison of the results showed good analytical and sampling precision.

CONCLUSION

Based on the QA/QC review, the data submitted were judged to be acceptable for use with the qualification noted.