

# Anion / Cation Summary Report

Lab ID: **0812076-1**

QC Type: SMP

Field ID: Angely WW

Analyte	Final Result	Report Units	mEq
BICARBONATE AS CaCO <sub>3</sub>	57.20488	MG/L	1.14
CHLORIDE	60.66372	MG/L	1.71
FLUORIDE	3.230523	MG/L	0.17
NITRATE AS N	0.4	MG/L	0.00
NITRITE AS N	0.2	MG/L	0.00
SULFATE	1036.584	MG/L	21.58
Anion Result Sum		<b>1158.28</b>	

Analyte	Final Result	Report Units	mEq
CALCIUM	188.9487	MG/L	9.43
IRON	0.1	MG/L	0.00
MAGNESIUM	2.529686	MG/L	0.21
MANGANESE	0.1867702	MG/L	0.01
POTASSIUM	2.167762	MG/L	0.06
SODIUM	264.6501	MG/L	11.51
Cation Result Sum		<b>458.58</b>	

Total Result: **1616.87** MG/LTDS Result: **1570.0001** MG/L

RPD: 2.94%

Anion mEq Sum: **24.61**Cation mEq Sum: **21.21**

RPD: 14.82%

Below is a list of Lab IDs for this Order Number that were logged in for metals analyses. Note: if this area is empty then either no metals analyses were requested or the cations of interest were not requested.

0812076-1



# ALS Paragon



## Dissolved Gasses Case Narrative

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### Colorado Oil & Gas Conservation Commission

Complaint 200199690

Work Order Number: 0812076

1. This report consists of 1 water sample. The sample was received cool and intact by ALS Paragon on 12/09/08. The vial for sample 0812076-1 contained headspace prior to analysis because it was not received headspace free. The sample had a pH > 2 at the time of analysis.
2. The sample was prepared and analyzed according to method RSK-175 procedures and SOP449R0.
3. The preparation batch included a method blank, laboratory control sample, laboratory control sample duplicate, and sample duplicate. Due to insufficient sample volume, a matrix spike was not prepared. The following is the sample used for the matrix QC:

Sample ID	QC Type	Batch ID
0812076-1	DUP	MEE081217-1

Similarity of matrix and therefore relevance of the QC results should not be automatically inferred for any sample other than the native sample selected for QC.

4. All preparation QC results were within the acceptance criteria.
5. All samples are associated with one or more of the following analytical QC: initial calibrations, initial calibration verifications (ICV), and continuing calibration verifications (CCV).
6. All analytical QC were within the acceptance criteria.
7. The sample was analyzed at a dilution in order to bring the target analytes within the calibration range of the instrument. The reporting limits have been adjusted accordingly.
8. The sample was prepared and analyzed within the established holding times.
9. Manual integrations are performed when needed to provide consistent and defensible data following the guidelines in SOP 939 Revision 3.



The data contained in the following report have been reviewed and approved by the personnel listed below. In addition, ALS Paragon certifies that the analyses reported herein are true, complete and correct within the limits of the methods employed.

Emily Knode  
Emily Knode  
Organics Primary Data Reviewer

12-22-08  
Date

Joe Norris  
Organics Final Data Reviewer

12-22-08  
Date

***Paragon Analytics***  
***Data Qualifier Flags***  
***Chromatography and Mass Spectrometry***

- U or ND:** This flag indicates that the compound was analyzed for but not detected.
- J:** This flag indicates an estimated value. This flag is used as follows: (1) when estimating a concentration for tentatively identified compounds (TICs) where a 1:1 response is assumed; (2) when the mass spectral and retention time data indicate the presence of a compound that meets the volatile and semivolatile GC/MS identification criteria, and the result is less than the reporting limit (RL) but greater than the method detection limit (MDL); (3) when the retention time data indicate the presence of a compound that meets the GC identification criteria, and the result is less than the RL but greater than the MDL; and (4) the reported value is estimated.
- B:** This flag is used when the analyte is detected in the associated method blank as well as in the sample. It indicates probable blank contamination and warns the data user. This flag shall be used for a tentatively identified compound (TIC) as well as for a positively identified target compound.
- E:** This flag identifies compounds whose concentration exceeds the upper level of the calibration range.
- A:** This flag indicates that a tentatively identified compound is a suspected aldol-condensation product.
- X:** This flag indicates that the analyte was diluted below an accurate quantitation level.
- \*:** This flag indicates that a spike recovery is equal to or outside the control criteria used.
- +**: This flag indicates that the relative percent difference (RPD) equals or exceeds the control criteria.

# ALS Paragon

## Sample Number(s) Cross-Reference Table

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**Paragon OrderNum:** 0812076

**Client Name:** Colorado Oil & Gas Conservation Commission

**Client Project Name:** Complaint 200199690

**Client Project Number:**

**Client PO Number:** OE PHA 09000000004

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Client Sample Number	Lab Sample Number	COC Number	Matrix	Date Collected	Time Collected
Angely WW	0812076-1		WATER	08-Dec-08	11:58
Trip Blank	0812076-2		WATER	08-Dec-08	

## Paragon Analytics

A Division of DataChem Laboratories, Inc.

225 Commerce Drive Fort Collins, CO 80524  
800-443-1511 or (970) 490-1511 (970) 490-1522 FaxAccession Number (LAB ID) Q812073 CT 12-9-08  
Chain-of-Custody Date 12/08/08 Page 1 of 1 Originator: Retain pink copy!Report To: Peter Grintat  
Phone: 719-466-3071  
Fax:  
E-mail:  
Company: Peter Grintat & Associates, Inc.  
Address: 100, C, 1st & Cross Cons. Lm.

Sampler(s): 6-HA-1001 Turnaround (circle one) Standard or Rush (Due 14 days) Dispose: Date 12/08/08 or Return to Client

Sample ID	Date	Time *	Lab ID	Matrix	Preservative (indicate type... HCl, etc.)	No. of Contaminers	Comments:										
							Oil	Water	Non-soil	Soil	Liquid	Filter	Extrac.	Water	Non-soil	Soil	Liquid
11/14/08-1152	11/14/08	1152	①	W	HCl	3	X										
11/14/08-1153	11/14/08	1153	②	W	HCl	5	X										
11/14/08-1154	11/14/08	1154	③	W	HCl	1											
11/14/08-1155	11/14/08	1155	④	W	HCl	7											
11/14/08-1156	11/14/08	1156	⑤	W	HCl	1											
11/14/08-1157	11/14/08	1157	⑥	W	HCl	3	X										
11/14/08-1158	11/14/08	1158	⑦	W	HCl	1											
11/14/08-1159	11/14/08	1159	⑧	W	HCl	1											
11/14/08-1160	11/14/08	1160	⑨	W	HCl	1											
11/14/08-1161	11/14/08	1161	⑩	W	HCl	1											
11/14/08-1162	11/14/08	1162	⑪	W	HCl	1											
11/14/08-1163	11/14/08	1163	⑫	W	HCl	1											
11/14/08-1164	11/14/08	1164	⑬	W	HCl	1											
11/14/08-1165	11/14/08	1165	⑭	W	HCl	1											
11/14/08-1166	11/14/08	1166	⑮	W	HCl	1											
11/14/08-1167	11/14/08	1167	⑯	W	HCl	1											
11/14/08-1168	11/14/08	1168	⑰	W	HCl	1											
11/14/08-1169	11/14/08	1169	⑱	W	HCl	1											
11/14/08-1170	11/14/08	1170	⑲	W	HCl	1											
11/14/08-1171	11/14/08	1171	⑳	W	HCl	1											
11/14/08-1172	11/14/08	1172	㉑	W	HCl	1											
11/14/08-1173	11/14/08	1173	㉒	W	HCl	1											
11/14/08-1174	11/14/08	1174	㉓	W	HCl	1											
11/14/08-1175	11/14/08	1175	㉔	W	HCl	1											
11/14/08-1176	11/14/08	1176	㉕	W	HCl	1											
11/14/08-1177	11/14/08	1177	㉖	W	HCl	1											
11/14/08-1178	11/14/08	1178	㉗	W	HCl	1											
11/14/08-1179	11/14/08	1179	㉘	W	HCl	1											
11/14/08-1180	11/14/08	1180	㉙	W	HCl	1											
11/14/08-1181	11/14/08	1181	㉚	W	HCl	1											
11/14/08-1182	11/14/08	1182	㉛	W	HCl	1											
11/14/08-1183	11/14/08	1183	㉜	W	HCl	1											
11/14/08-1184	11/14/08	1184	㉝	W	HCl	1											
11/14/08-1185	11/14/08	1185	㉞	W	HCl	1											
11/14/08-1186	11/14/08	1186	㉟	W	HCl	1											
11/14/08-1187	11/14/08	1187	㉟	W	HCl	1											
11/14/08-1188	11/14/08	1188	㉟	W	HCl	1											
11/14/08-1189	11/14/08	1189	㉟	W	HCl	1											
11/14/08-1190	11/14/08	1190	㉟	W	HCl	1											
11/14/08-1191	11/14/08	1191	㉟	W	HCl	1											
11/14/08-1192	11/14/08	1192	㉟	W	HCl	1											
11/14/08-1193	11/14/08	1193	㉟	W	HCl	1											
11/14/08-1194	11/14/08	1194	㉟	W	HCl	1											
11/14/08-1195	11/14/08	1195	㉟	W	HCl	1											
11/14/08-1196	11/14/08	1196	㉟	W	HCl	1											
11/14/08-1197	11/14/08	1197	㉟	W	HCl	1											
11/14/08-1198	11/14/08	1198	㉟	W	HCl	1											
11/14/08-1199	11/14/08	1199	㉟	W	HCl	1											
11/14/08-1200	11/14/08	1200	㉟	W	HCl	1											
11/14/08-1201	11/14/08	1201	㉟	W	HCl	1											
11/14/08-1202	11/14/08	1202	㉟	W	HCl	1											
11/14/08-1203	11/14/08	1203	㉟	W	HCl	1											
11/14/08-1204	11/14/08	1204	㉟	W	HCl	1											
11/14/08-1205	11/14/08	1205	㉟	W	HCl	1											
11/14/08-1206	11/14/08	1206	㉟	W	HCl	1											
11/14/08-1207	11/14/08	1207	㉟	W	HCl	1											
11/14/08-1208	11/14/08	1208	㉟	W	HCl	1											
11/14/08-1209	11/14/08	1209	㉟	W	HCl	1											
11/14/08-1210	11/14/08	1210	㉟	W	HCl	1											
11/14/08-1211	11/14/08	1211	㉟	W	HCl	1											
11/14/08-1212	11/14/08	1212	㉟	W	HCl	1											
11/14/08-1213	11/14/08	1213	㉟	W	HCl	1											
11/14/08-1214	11/14/08	1214	㉟	W	HCl	1											
11/14/08-1215	11/14/08	1215	㉟	W	HCl	1											
11/14/08-1216	11/14/08	1216	㉟	W	HCl	1											
11/14/08-1217	11/14/08	1217	㉟	W	HCl	1											
11/14/08-1218	11/14/08	1218	㉟	W	HCl	1											
11/14/08-1219	11/14/08	1219	㉟	W	HCl	1											
11/14/08-1220	11/14/08	1220	㉟	W	HCl	1											
11/14/08-1221	11/14/08	1221	㉟	W	HCl	1											
11/14/08-1222	11/14/08	1222	㉟	W	HCl	1											
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11/14/08-1224	11/14/08	1224	㉟	W	HCl	1											
11/14/08-1225	11/14/08	1225	㉟	W	HCl	1											
11/14/08-1226	11/14/08	1226	㉟	W	HCl	1											
11/14/08-1227	11/14/08	1227	㉟	W	HCl	1											
11/14/08-1228	11/14/08	1228	㉟	W	HCl	1											
11/14/08-1229	11/14/08	1229	㉟	W	HCl	1											
11/14/08-1230	11/14/08	1230	㉟	W	HCl	1											
11/14/08-1231	11/14/08	1231	㉟	W	HCl	1											
11/14/08-1232	11/14/08	1232	㉟	W	HCl	1											
11/14/08-1233	11/14/08	1233	㉟	W	HCl	1											
11/14/08-1234	11/14/08	1234	㉟	W	HCl	1											
11/14/08-1235	11/14/08	1235	㉟	W	HCl	1											
11/14/08-1236	11/14/08	1236	㉟	W	HCl	1											
11/14/08-1237	11/14/08	1237	㉟	W	HCl	1											
11/14/08-1238	11/14/08	1238	㉟	W	HCl	1											
11/14/08-1239	11/14/08	1239	㉟	W	HCl	1											
11/14/08-1240	11/14/08	1240	㉟	W	HCl	1											
11/14/08-1241	11/14/08	1241	㉟	W	HCl	1											
11/14/08-1242	11/14/08	1242															

## CONDITION OF SAMPLE UPON RECEIPT FORM

Paragon Analytics

Client: COGCCWorkorder No: 0812076Project Manager: AWInitials: CDTDate: 12-9-08

1. Does this project require any <b>special handling</b> in addition to standard Paragon procedures?	YES	NO		
2. Are custody <b>seals on shipping containers</b> intact?	NONE	YES	NO	
3. Are Custody seals on <b>sample containers</b> intact?	(NONE)	YES	NO	
4. Is there a <b>COC (Chain-of-Custody) present</b> or other representative documents?	(YES)	YES	NO	
5. Are the <b>COC and bottle labels complete and legible</b> ?	(YES)	YES	NO	
6. Is the <b>COC in agreement</b> with samples received? (IDs, dates, times, no. of samples, no. of containers, matrix, requested analyses, etc.)	(YES)	YES	NO	
7. Were <b>airbills / shipping documents</b> present and/or removable?	DROP OFF	YES	NO	
8. Are all aqueous <b>samples requiring preservation</b> preserved correctly? (excluding volatiles)	N/A	YES	NO	
9. Are all aqueous <b>non-preserved samples pH 4-9</b> ?	N/A	YES	NO	
10. Is there <b>sufficient sample</b> for the requested analyses?	(YES)	YES	NO	
11. Were all samples placed in the <b>proper containers</b> for the requested analyses?	(YES)	YES	NO	
12. Are all samples within <b>holding times</b> for the requested analyses?	(YES)	YES	NO	
13. Were all sample containers received <b>intact</b> ? (not broken or leaking, etc.)	(YES)	YES	NO	
14. Are all samples requiring <b>no headspace</b> (VOC, GRO, RSK/MEE, Rx CN/S, radon) headspace free? Size of bubble: <u>&lt; green pea</u> ✓ <u>&gt; green pea</u>	N/A	YES	NO	
15. Do perchlorate LCMS-MS samples <b>have headspace</b> ? (at least 1/3 of container required)	(N/A)	YES	NO	
16. Were samples checked for and free from the presence of <b>residual chlorine</b> ? (Applicable when PM has indicated samples are from a chlorinated water source; note if field preservation with sodium thiosulfate was not observed.)	(N/A)	YES	NO	
17. Were the samples <b>shipped on ice</b> ?	(YES)	YES	NO	
18. Were cooler temperatures measured at 0.1-6.0°C?	IR gun used*: #2      #4	RAD ONLY	YES	NO
Cooler #: <u>1</u>				
Temperature (°C): <u>4.8</u>				
No. of custody seals on cooler: <u>2</u>				
DOT Survey/ Acceptance Information	External µR/hr reading:	<u>13</u>		
	Background µR/hr reading:	<u>10</u>		
Were external µR/hr readings ≤ two times background and within DOT acceptance criteria? <u>YES</u> NO / NA (If no. see Form 008.)				

Additional Information: PROVIDE DETAILS BELOW FOR A NO RESPONSE TO ANY QUESTION ABOVE, EXCEPT #1 AND #16.

HEADSPACE ALL &gt; green pea 0812076-1-1 0812076-2-1

↓      ↓      ↓      ↓

-2      -3      -2      -3

-3      -4      -3      -3

-4      -5      -4      -4

-5

If applicable, was the client contacted? YES NO / NA Contact: Peter ConstantasDate/Time: 12/9/08

e-mail

Project Manager Signature / Date: Connelly 12/9/08

# Dissolved Gasses

## Method RSK175

### Method Blank

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: MEE081217-1MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 17-Dec-08

Date Analyzed: 17-Dec-08

Prep Method: METHOD

Prep Batch: MEE081217-1

QCBatchID: MEE081217-1-1

Run ID: MEE081217-1A

Cleanup: NONE

Basis: N/A

File Name: 00837.dat

Sample Aliquot: 38.5 ml

Final Volume: 38.5 ml

Result Units: UG/L

Clean DF: 1

CASNO	Target Analyte	DF	Result	Reporting Limit	Result Qualifier	EPA Qualifier
74-82-8	METHANE	1	1	1	U	
74-85-1	ETHENE	1	1	1	U	
74-84-0	ETHANE	1	2	2	U	

Data Package ID: MEE0812076-1

Date Printed: Monday, December 22, 2008

ALS Paragon

LIMS Version: 6.217A

Page 1 of 1

# Dissolved Gasses

## Method RSK175 Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:	Angely WW
Lab ID:	0812076-1

Sample Matrix: WATER  
% Moisture: N/A  
Date Collected: 08-Dec-08  
Date Extracted: 17-Dec-08  
Date Analyzed: 17-Dec-08  
Prep Method: METHOD  
Prep Batch: MEE081217-1  
QCBatchID: MEE081217-1-1  
Run ID: MEE081217-1A  
Cleanup: NONE  
Basis: As Received  
File Name: 00842.dat  
Sample Aliquot: 38.5 ml  
Final Volume: 38.5 ml  
Result Units: UG/L  
Clean DF: 1

CASNO	Target Analyte	Dilution Factor	Result	Reporting Limit	Result Qualifier	EPA Qualifier
74-82-8	METHANE	10	100000	10		
74-85-1	ETHENE	10	10	10	U	
74-84-0	ETHANE	10	140	20		

Data Package ID: MEE0812076-1

Date Printed: Monday, December 22, 2008

ALS Paragon

LIMS Version: 6.217A

Page 1 of 1

# Dissolved Gasses

## Method RSK175

### Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: MEE081217-1LCS

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 12/17/2008

Date Analyzed: 12/17/2008

Prep Method: METHOD

Prep Batch: MEE081217-1

QCBatchID: MEE081217-1-1

Run ID: MEE081217-1A

Cleanup: NONE

Basis: N/A

File Name: 00836.dat

Sample Aliquot: 38.5 ml

Final Volume: 38.5 ml

Result Units: UG/L

Clean DF: 1

CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
74-82-8	METHANE	140	147	1		105	80 - 120%
74-85-1	ETHENE	245	256	1		104	80 - 120%
74-84-0	ETHANE	262	256	2		98	80 - 120%

Lab ID: MEE081217-1LCSD

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 12/17/2008

Date Analyzed: 12/17/2008

Prep Method: METHOD

Prep Batch: MEE081217-1

QCBatchID: MEE081217-1-1

Run ID: MEE081217-1A

Cleanup: NONE

Basis: N/A

File Name: 00844.dat

Sample Aliquot: 38.5 ml

Final Volume: 38.5 ml

Result Units: UG/L

Clean DF: 1

CASNO	Target Analyte	Spike Added	LCSD Result	Reporting Limit	Result Qualifier	LCSD % Rec.	RPD Limit	RPD
74-82-8	METHANE	140	150	1		107	25	2
74-85-1	ETHENE	245	253	1		103	25	1
74-84-0	ETHANE	262	253	2		97	25	1

Data Package ID: MEE0812076-1

Date Printed: Monday, December 22, 2008

ALS Paragon

LIMS Version: 6.217A

Page 1 of 1

# Dissolved Gasses

## Method RSK175 Duplicate Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:	Angely WW
Lab ID:	0812076-1D

Sample Matrix: WATER  
% Moisture: N/A  
Prep Batch: MEE081217-1  
QCBatchID: MEE081217-1-1  
Date Collected: 12/08/2008  
Run ID: MEE081217-1A  
Date Extracted: 12/17/2008  
Cleanup: NONE  
Date Analyzed: 12/17/2008  
Basis: As Received  
File Name: 00843.dat  
Result Units: UG/L  
Clean DF: 1  
Sample Aliquot: 38.5 ml  
Final Volume: 38.5 ml

CASNO	Target Analyte	Sample Result	Samp Qual	Duplicate Result	Dup Qual	Reporting Limit	Dilution Factor	RPD	RPD Limit
74-82-8	METHANE	100000		92800		10	10	8	25
74-85-1	ETHENE	10	U	10	U	10	10		25
74-84-0	ETHANE	140		131		20	10	7	25

Data Package ID: MEE0812076-1

Date Printed: Monday, December 22, 2008

ALS Paragon

LIMS Version: 6.217A

Page 1 of 1



# ALS Paragon



## Total Extractable Hydrocarbons (Diesel) Case Narrative

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### Colorado Oil & Gas Conservation Commission

Complaint 200199690

**Order Number - 0812076**

1. This report consists of 1 water sample. The sample was received cool and intact by ALS Paragon on 12/09/2008.
2. The water sample was extracted using separatory funnels according to SOP 626 Revision 9 based on Method 3510C.
3. The extract was then analyzed using GC with a DB-5.625 capillary column and a flame ionization detector (FID) according to SOP 406 Revision 13 generally based on SW-846 Method 8000B and Method 8015B and specifically on the California LUFT Field Manual (October 1989 revision). The procedures are based on this general method because SW-846 does not have a specific method for total extractable petroleum hydrocarbons (TEPH) or diesel range organics. The only true modification from this method is that TEPH is a multicomponent mixture and is quantitated by integrating across the entire range, rather than summing areas of individual peaks. All positive results were quantitated using the responses from the initial calibration curve using the external standard technique. Also, a confirmation column is not used, because the analyte is a multicomponent mixture and the specific carbon range of the peaks detected is specified on the individual sample reporting forms.
4. All initial and continuing calibration criteria were met.
5. The method blank associated with this project was below the MDL for diesel range organics.
6. All laboratory control sample and laboratory control sample duplicate recoveries and RPDs were within the acceptance criteria.
7. Matrix spikes and matrix spike duplicates could not be performed because of insufficient sample. A laboratory control sample and laboratory control sample duplicate were performed instead.
8. The sample was extracted and analyzed within the established holding time.
9. All surrogate recoveries were within the acceptance criteria.



11. Manual integrations are performed when needed to provide consistent and defensible data following the guidelines in SOP 939 Revision 3.

The data contained in the following report have been reviewed and approved by the personnel listed below. In addition, ALS Paragon certifies that the analyses reported herein are true, complete and correct within the limits of the methods employed.

Mindy Norton  
Mindy Norton  
Organics Primary Data Reviewer

10-17-08  
Date

Joz Norton  
Organics Final Data Reviewer

12-18-08  
Date



**ALS Paragon**  
**Data Qualifier Flags**  
**Fuels**

- G:** This flag indicates that a pattern resembling gasoline was detected in this sample.
- D:** This flag indicates that a pattern resembling diesel was detected in this sample.
- M:** This flag indicates that a pattern resembling motor oil was detected in this sample.
- C:** This flag indicates that a pattern resembling crude oil was detected in this sample.
- 4:** This flag indicates that a pattern resembling JP-4 was detected in this sample.
- 5:** This flag indicates that a pattern resembling JP-5 was detected in this sample.
- H:** This flag indicates that the fuel pattern was in the heavier end of the retention time window for the analyte of interest.
- L:** This flag indicates that the fuel pattern was in the lighter end of the retention time window for the analyte of interest.
- Z:** This flag indicates that a significant fraction of the reported result did not resemble the patterns of any of the following petroleum hydrocarbon products:  
gasoline  
JP-8  
diesel  
mineral spirits  
motor oil  
Stoddard solvent  
bunker C

Multiple flags may be used to indicate the presence of more than one product or component.

***Paragon Analytics, Inc.***  
***Data Qualifier Flags***  
***Chromatography and Mass Spectrometry***

- U or ND:** This flag indicates that the compound was analyzed for but not detected.
- J:** This flag indicates an estimated value. This flag is used as follows : (1) when estimating a concentration for tentatively identified compounds (TICs) where a 1:1 response is assumed; (2) when the mass spectral and retention time data indicate the presence of a compound that meets the volatile and semivolatile GC/MS identification criteria, and the result is less than the reporting limit (RL) but greater than the method detection limit (MDL); (3) when the data indicate the presence of a compound that meets the identification criteria, and the result is less than the RL but greater than the MDL; and (4) the reported value is estimated.
- B:** This flag is used when the analyte is detected in the associated method blank as well as in the sample. It indicates probable blank contamination and warns the data user. This flag shall be used for a tentatively identified compound (TIC) as well as for a positively identified target compound.
- E:** This flag identifies compounds whose concentration exceeds the upper level of the calibration range.
- A:** This flag indicates that a tentatively identified compound is a suspected aldol-condensation product.
- X:** This flag indicates that the analyte was diluted below an accurate quantitation level.
- \***: This flag indicates that a spike recovery is outside the control criteria.
- +**: This flag indicates that the relative percent difference (RPD) exceeds the control criteria.

# ALS Paragon

## Sample Number(s) Cross-Reference Table

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**Paragon OrderNum:** 0812076

**Client Name:** Colorado Oil & Gas Conservation Commission

**Client Project Name:** Complaint 200199690

**Client Project Number:**

**Client PO Number:** OE PHA 09000000004

---

Client Sample Number	Lab Sample Number	COC Number	Matrix	Date Collected	Time Collected
Angely WW	0812076-1		WATER	08-Dec-08	11:58
Trip Blank	0812076-2		WATER	08-Dec-08	



## CONDITION OF SAMPLE UPON RECEIPT FORM

Paragon Analytics

Client: COGCCWorkorder No: 0812076Project Manager: AWInitials: CDTDate: 12-9-08

1. Does this project require any <b>special handling</b> in addition to standard Paragon procedures?	YES	NO		
2. Are custody <b>seals on shipping containers</b> intact?	NONE	YES	NO	
3. Are Custody seals on <b>sample containers</b> intact?	(NONE)	YES	NO	
4. Is there a <b>COC (Chain-of-Custody) present</b> or other representative documents?	(YES)	YES	NO	
5. Are the <b>COC and bottle labels complete and legible</b> ?	(YES)	YES	NO	
6. Is the <b>COC in agreement</b> with samples received? (IDs, dates, times, no. of samples, no. of containers, matrix, requested analyses, etc.)	(YES)	YES	NO	
7. Were <b>airbills / shipping documents</b> present and/or removable?	DROP OFF	YES	NO	
8. Are all aqueous <b>samples requiring preservation</b> preserved correctly? (excluding volatiles)	N/A	YES	NO	
9. Are all aqueous <b>non-preserved samples pH 4-9</b> ?	N/A	YES	NO	
10. Is there <b>sufficient sample</b> for the requested analyses?	(YES)	YES	NO	
11. Were all samples placed in the <b>proper containers</b> for the requested analyses?	(YES)	YES	NO	
12. Are all samples within <b>holding times</b> for the requested analyses?	(YES)	YES	NO	
13. Were all sample containers received <b>intact</b> ? (not broken or leaking, etc.)	(YES)	YES	NO	
14. Are all samples requiring <b>no headspace</b> (VOC, GRO, RSK/MEE, Rx CN/S, radon) headspace free? Size of bubble: <u>&lt; green pea</u> ✓ <u>&gt; green pea</u>	N/A	YES	NO	
15. Do perchlorate LCMS-MS samples <b>have headspace</b> ? (at least 1/3 of container required)	(N/A)	YES	NO	
16. Were samples checked for and free from the presence of <b>residual chlorine</b> ? (Applicable when PM has indicated samples are from a chlorinated water source; note if field preservation with sodium thiosulfate was not observed.)	(N/A)	YES	NO	
17. Were the samples <b>shipped on ice</b> ?	(YES)	YES	NO	
18. Were cooler temperatures measured at 0.1-6.0°C?	IR gun used*: #2      #4	RAD ONLY	YES	NO
Cooler #: <u>1</u>				
Temperature (°C): <u>4.8</u>				
No. of custody seals on cooler: <u>2</u>				
DOT Survey/ Acceptance Information	External µR/hr reading:	<u>13</u>		
	Background µR/hr reading:	<u>10</u>		
Were external µR/hr readings ≤ two times background and within DOT acceptance criteria? <u>YES</u> NO / NA (If no. see Form 008.)				

Additional Information: PROVIDE DETAILS BELOW FOR A NO RESPONSE TO ANY QUESTION ABOVE, EXCEPT #1 AND #16.

HEADSPACE ALL &gt; green pea 0812076-1-1 0812076-2-1

↓      ↓      ↓      ↓

-2      -3      -2      -3

-3      -4      -3      -3

-4      -5      -4      -4

-5

If applicable, was the client contacted? YES NO / NA Contact: Peter ConstantasDate/Time: 12/9/08

e-mail

Project Manager Signature / Date: Connelly 12/9/08

# Total Extractable Hydrocarbons

## Method SW8015MCALUFTB

### Method Blank

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: EX081211-3MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 11-Dec-08

Date Analyzed: 15-Dec-08

Prep Method: SW3510 Rev C

Prep Batch: EX081211-3

QCBatchID: EX081211-3-2

Run ID: HCD081215-3A

Cleanup: NONE

Basis: N/A

File Name: F3F33074

Sample Aliquot: 1000 ml

Final Volume: 2.5 ml

Result Units: mg/l

Clean DF: 1

CASNO	Target Analyte	DF	Result	Reporting Limit	Result Qualifier	EPA Qualifier
68334-30-5	DIESEL RANGE ORGANICS	1	0.05	0.05	U	

### Surrogate Recovery

CASNO	Surrogate Analyte	Result	Flag	Spike Amount	Percent Recovery	Control Limits
84-15-1	O-TERPHENYL	0.231		0.25	93	60 - 140

Data Package ID: HCD0812076-1

Date Printed: Wednesday, December 17, 2008

ALS Paragon

LIMS Version: 6.216A

Page 1 of 1

# Total Extractable Hydrocarbons

## Method SW8015MCALUFTB Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:	Angely WW
Lab ID:	0812076-1

Sample Matrix: WATER  
% Moisture: N/A  
Prep Batch: EX081211-3  
QCBatchID: EX081211-3-2  
Date Collected: 08-Dec-08  
Run ID: HCD081215-3A  
Date Extracted: 11-Dec-08  
Cleanup: NONE  
Date Analyzed: 15-Dec-08  
Basis: As Received  
Prep Method: SW3510 Rev C  
File Name: F3F33077  
Final Volume: 2.5 ml  
Result Units: mg/l  
Clean DF: 1

CASNO	Target Analyte	Dilution Factor	Result	Reporting Limit	Result Qualifier	EPA Qualifier
68334-30-5	DIESEL RANGE ORGANICS	1	0.048	0.048	U	

## Surrogate Recovery

CASNO	Surrogate Analyte	Result	Flag	Spike Amount	Percent Recovery	Control Limits
84-15-1	O-TERPHENYL	0.21		0.238	88	60 - 140

Data Package ID: HCD0812076-1

Date Printed: Wednesday, December 17, 2008

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Page 1 of 1

# Total Extractable Hydrocarbons

## Method SW8015MCALUFTB

### Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: EX081211-3LCS	Sample Matrix: WATER % Moisture: N/A Date Collected: N/A Date Extracted: 12/11/2008 Date Analyzed: 12/15/2008 Prep Method: SW3510C	Prep Batch: EX081211-3 QCBatchID: EX081211-3-2 Run ID: HCD081215-3A Cleanup: NONE Basis: N/A File Name: F3F33075	Sample Aliquot: 1000 ml Final Volume: 2.5 ml Result Units: mg/l Clean DF: 1
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CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
68334-30-5	DIESEL RANGE ORGANICS	1	0.843	0.05		84	60 - 140%

Lab ID: EX081211-3LCSD	Sample Matrix: WATER % Moisture: N/A Date Collected: N/A Date Extracted: 12/11/2008 Date Analyzed: 12/15/2008 Prep Method: SW3510C	Prep Batch: EX081211-3 QCBatchID: EX081211-3-2 Run ID: HCD081215-3A Cleanup: NONE Basis: N/A File Name: F3F33076	Sample Aliquot: 1000 ml Final Volume: 2.5 ml Result Units: mg/l Clean DF: 1
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CASNO	Target Analyte	Spike Added	LCSD Result	Reporting Limit	Result Qualifier	LCSD % Rec.	RPD Limit	RPD
68334-30-5	DIESEL RANGE ORGANICS	1	0.882	0.05		88	50	5

### Surrogate Recovery LCS/LCSD

CASNO	Target Analyte	Spike Added	LCS % Rec.	LCS Flag	LCSD % Rec.	LCSD Flag	Control Limits
84-15-1	O-TERPHENYL	0.25	94		97		60 - 140

Data Package ID: HCD0812076-1

Date Printed: Wednesday, December 17, 2008

ALS Paragon

LIMS Version: 6.216A

Page 1 of 1



# ALS Paragon



## GC/MS Semivolatiles Case Narrative

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### Colorado Oil & Gas Conservation Commission

Complaint 200199690

**Order Number - 0812076**

1. This report consists of 1 water sample. The sample was received cool and intact by ALS Paragon on 12/09/08.
2. The sample was prepared and analyzed according to SW-846, 3rd Edition protocol utilizing SOP. Specifically, the water sample was extracted using continuous liquid-liquid extractors, according to SW-846 Method 3520C utilizing SOP 617 Revision 13.
3. The extracts were analyzed using GC/MS with a DB-5.625 capillary column according to SOP 506 Revision 15 based on SW-846 Method 8270D. All positive results were quantitated against the initial calibration standards using the internal standard technique. The identification of positive results was achieved by a comparison of the retention time and mass spectrum of the sample versus the daily calibration standard.
4. All initial calibration criteria for SPCC's and CCC's were met. If average response factors were used in the initial calibration, %RSD was  $\leq 15\%$ . If linear or higher order regression calibrations were used in the initial calibration, the coefficient of determination ( $r^2$ )  $\geq 0.99$ .
5. All initial calibration standards are verified by comparing a second source standard initial calibration verification (ICV) against the calibration curve. All target compounds in the second source verification had a %D of less than 25%.
6. All SPCC and CCC criteria were met in each of the daily (continuing) calibration verifications.
7. All method blank criteria were met.
8. All laboratory control sample and laboratory control sample duplicate criteria were met with the following exceptions:

Spiked Compound	QC Sample	Direction
Pyridine	LCS/LCSD	RPD High



Aniline	LCSD	Low
Aniline	LCS/LCSD	RPD High
3,3'-Dichlorobenzidine	LCS/LCSD	RPD High

Because of the large number of target analytes reported by this method, the lab allows for sporadic marginal exceedances. No further action was taken.

Since the recoveries for pyridine and 3,3'-dichlorobenzidine in the laboratory control sample and laboratory control sample duplicate were within control limits, with only the RPD exceeding acceptance criteria, quantitations of target compounds were not compromised. No further action was taken.

9. Since a sample from this order number was not the selected quality control (QC) sample, matrix specific QC results are not included in this report.
10. The sample was extracted and analyzed within the established holding time.
11. All surrogate recoveries were within acceptance criteria.
12. All internal standard recoveries were within acceptance criteria.

The data contained in the following report have been reviewed and approved by the personnel listed below. In addition, ALS Paragon certifies that the analyses reported herein are true, complete and correct within the limits of the methods employed.

SLJ  
Sharon L. Jones  
Organics Primary Data Reviewer

12-16-08  
Date

Joe Marshall Jr.  
Organics Final Data Reviewer

December 16, 2008  
Date



***ALS Paragon***  
***Data Qualifier Flags***  
***Chromatography and Mass Spectrometry***

- U or ND:** This flag indicates that the compound was analyzed for but not detected.
- J:** This flag indicates an estimated value. This flag is used as follows: (1) when estimating a concentration for tentatively identified compounds (TICs) where a 1:1 response is assumed; (2) when the mass spectral and retention time data indicate the presence of a compound that meets the volatile and semivolatile GC/MS identification criteria, and the result is less than the reporting limit (RL) but greater than the method detection limit (MDL); (3) when the retention time data indicate the presence of a compound that meets the GC identification criteria, and the result is less than the RL but greater than the MDL; and (4) the reported value is estimated.
- B:** This flag is used when the analyte is detected in the associated method blank as well as in the sample. It indicates probable blank contamination and warns the data user. This flag shall be used for a tentatively identified compound (TIC) as well as for a positively identified target compound.
- E:** This flag identifies compounds whose concentration exceeds the upper level of the calibration range.
- A:** This flag indicates that a tentatively identified compound is a suspected aldol-condensation product.
- X:** This flag indicates that the analyte was diluted below an accurate quantitation level.
- \*:** This flag indicates that a spike recovery is equal to or outside the control criteria used.
- +:** This flag indicates that the relative percent difference (RPD) equals or exceeds the control criteria.

# ALS Paragon

## Sample Number(s) Cross-Reference Table

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**Paragon OrderNum:** 0812076

**Client Name:** Colorado Oil & Gas Conservation Commission

**Client Project Name:** Complaint 200199690

**Client Project Number:**

**Client PO Number:** OE PHA 09000000004

---

Client Sample Number	Lab Sample Number	COC Number	Matrix	Date Collected	Time Collected
Angely WW	0812076-1		WATER	08-Dec-08	11:58
Trip Blank	0812076-2		WATER	08-Dec-08	



## CONDITION OF SAMPLE UPON RECEIPT FORM

Paragon Analytics

Client: COGCCWorkorder No: 0812076Project Manager: AWInitials: CDTDate: 12-9-08

1. Does this project require any <b>special handling</b> in addition to standard Paragon procedures?	YES	NO	
2. Are custody <b>seals on shipping containers</b> intact?	NONE	YES	NO
3. Are Custody seals on <b>sample containers</b> intact?	(NONE)	YES	NO
4. Is there a <b>COC (Chain-of-Custody) present</b> or other representative documents?	(YES)	YES	NO
5. Are the <b>COC and bottle labels complete and legible</b> ?	(YES)	YES	NO
6. Is the <b>COC in agreement</b> with samples received? (IDs, dates, times, no. of samples, no. of containers, matrix, requested analyses, etc.)	(YES)	YES	NO
7. Were <b>airbills / shipping documents</b> present and/or removable?	DROP OFF	YES	NO
8. Are all aqueous <b>samples requiring preservation</b> preserved correctly? (excluding volatiles)	N/A	YES	NO
9. Are all aqueous <b>non-preserved samples pH 4-9</b> ?	N/A	YES	NO
10. Is there <b>sufficient sample</b> for the requested analyses?	(YES)	YES	NO
11. Were all samples placed in the <b>proper containers</b> for the requested analyses?	(YES)	YES	NO
12. Are all samples within <b>holding times</b> for the requested analyses?	(YES)	YES	NO
13. Were all sample containers received <b>intact</b> ? (not broken or leaking, etc.)	(YES)	YES	NO
14. Are all samples requiring <b>no headspace</b> (VOC, GRO, RSK/MEE, Rx CN/S, radon) headspace free? Size of bubble: <u>&lt; green pea</u> ✓ <u>&gt; green pea</u>	N/A	YES	NO
15. Do perchlorate LCMS-MS samples <b>have headspace</b> ? (at least 1/3 of container required)	(N/A)	YES	NO
16. Were samples checked for and free from the presence of <b>residual chlorine</b> ? (Applicable when PM has indicated samples are from a chlorinated water source; note if field preservation with sodium thiosulfate was not observed.)	(N/A)	YES	NO
17. Were the samples <b>shipped on ice</b> ?	(YES)	YES	NO
18. Were cooler temperatures measured at 0.1-6.0°C?	IR gun used*: #2      #4 RAD ONLY	YES	NO
Cooler #:	<u>1</u>		
Temperature (°C):	<u>4.8</u>		
No. of custody seals on cooler:	<u>2</u>		
External µR/hr reading:	<u>13</u>		
Background µR/hr reading:	<u>10</u>		
Were external µR/hr readings ≤ two times background and within DOT acceptance criteria? <u>YES</u> NO / NA (If no. see Form 008.)			

Additional Information: PROVIDE DETAILS BELOW FOR A NO RESPONSE TO ANY QUESTION ABOVE, EXCEPT #1 AND #16.

HEADSPACE ALL > green pea    0812076-1-1    0812076-2-1

↓      -2  
 ↓      -3  
 ↓      -4  
 ↓      -5

If applicable, was the client contacted? YES NO / NA Contact: Peter ConstantasDate/Time: 12/9/08

e-mail

Project Manager Signature / Date: Connelly 12/9/08

# GC/MS Semi-volatiles

## Method SW8270D

### Method Blank

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: EX081210-1MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 10-Dec-08

Date Analyzed: 12-Dec-08

Prep Method: SW3520 Rev C

Prep Batch: EX081210-1

QCBatchID: EX081210-1-1

Run ID: SV081212-3

Cleanup: NONE

Basis: N/A

File Name: R9981

Sample Aliquot: 1000 ml

Final Volume: 1 ml

Result Units: UG/L

Clean DF: 1

CASNO	Target Analyte	DF	Result	Reporting Limit	Result Qualifier	EPA Qualifier
110-86-1	PYRIDINE	1	10	10	U	
62-75-9	N-NITROSODIMETHYLAMINE	1	10	10	U	
62-53-3	ANILINE	1	10	10	U	
108-95-2	PHENOL	1	10	10	U	
111-44-4	BIS(2-CHLOROETHYL)ETHER	1	10	10	U	
95-57-8	2-CHLOROPHENOL	1	10	10	U	
541-73-1	1,3-DICHLOROBENZENE	1	10	10	U	
106-46-7	1,4-DICHLOROBENZENE	1	10	10	U	
95-50-1	1,2-DICHLOROBENZENE	1	10	10	U	
100-51-6	BENZYL ALCOHOL	1	10	10	U	
108-60-1	BIS(2-CHLOROISOPROPYL)ETHER	1	10	10	U	
95-48-7	2-METHYLPHENOL	1	10	10	U	
621-64-7	N-NITROSO-DI-N-PROPYLAMINE	1	10	10	U	
108-39-4	3+4-METHYLPHENOL	1	10	10	U	
67-72-1	HEXACHLOROETHANE	1	10	10	U	
98-95-3	NITROBENZENE	1	10	10	U	
78-59-1	ISOPHORONE	1	10	10	U	
88-75-5	2-NITROPHENOL	1	10	10	U	
105-67-9	2,4-DIMETHYLPHENOL	1	10	10	U	
111-91-1	BIS(2-CHLOROETHOXY)METHANE	1	10	10	U	
120-83-2	2,4-DICHLOROPHENOL	1	10	10	U	
65-85-0	BENZOIC ACID	1	50	50	U	
120-82-1	1,2,4-TRICHLOROBENZENE	1	10	10	U	
91-20-3	NAPHTHALENE	1	10	10	U	
106-47-8	4-CHLOROANILINE	1	10	10	U	
87-68-3	HEXACHLOROBUTADIENE	1	10	10	U	
59-50-7	4-CHLORO-3-METHYLPHENOL	1	10	10	U	

Data Package ID: SV0812076-1

Date Printed: Tuesday, December 16, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 3

# GC/MS Semi-volatiles

## Method SW8270D

### Method Blank

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: EX081210-1MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 10-Dec-08

Date Analyzed: 12-Dec-08

Prep Method: SW3520 Rev C

Prep Batch: EX081210-1

QCBatchID: EX081210-1-1

Run ID: SV081212-3

Cleanup: NONE

Basis: N/A

File Name: R9981

Sample Aliquot: 1000 ml

Final Volume: 1 ml

Result Units: UG/L

Clean DF: 1

91-57-6	2-METHYLNAPHTHALENE	1	10	10	U	
90-12-0	1-METHYLNAPHTHALENE	1	10	10	U	
77-47-4	HEXACHLOROCYCLOPENTADIENE	1	10	10	U	
88-06-2	2,4,6-TRICHLOROPHENOL	1	10	10	U	
95-95-4	2,4,5-TRICHLOROPHENOL	1	10	10	U	
91-58-7	2-CHLORONAPHTHALENE	1	10	10	U	
88-74-4	2-NITROANILINE	1	20	20	U	
131-11-3	DIMETHYL PHTHALATE	1	10	10	U	
606-20-2	2,6-DINITROTOLUENE	1	10	10	U	
208-96-8	ACENAPHTHYLENE	1	10	10	U	
99-09-2	3-NITROANILINE	1	20	20	U	
83-32-9	ACENAPHTHENE	1	10	10	U	
51-28-5	2,4-DINITROPHENOL	1	20	20	U	
100-02-7	4-NITROPHENOL	1	20	20	U	
132-64-9	DIBENZOFURAN	1	10	10	U	
121-14-2	2,4-DINITROTOLUENE	1	10	10	U	
84-66-2	DIETHYL PHTHALATE	1	10	10	U	
86-73-7	FLUORENE	1	10	10	U	
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	1	10	10	U	
100-01-6	4-NITROANILINE	1	20	20	U	
103-33-3	AZOBENZENE	1	10	10	U	
534-52-1	4,6-DINITRO-2-METHYLPHENOL	1	20	20	U	
86-30-6	N-NITROSODIPHENYLAMINE	1	10	10	U	
101-55-3	4-BROMOPHENYL PHENYL ETHER	1	10	10	U	
118-74-1	HEXACHLOROBENZENE	1	10	10	U	
58-90-2	2,3,4,6-TETRACHLOROPHENOL	1	10	10	U	
87-86-5	PENTACHLOROPHENOL	1	20	20	U	
85-01-8	PHENANTHRENE	1	10	10	U	
120-12-7	ANTHRACENE	1	10	10	U	

Data Package ID: SV0812076-1

Date Printed: Tuesday, December 16, 2008

ALS Paragon

LIMS Version: 6.215A

Page 2 of 3

# GC/MS Semi-volatiles

## Method SW8270D

### Method Blank

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: EX081210-1MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 10-Dec-08

Date Analyzed: 12-Dec-08

Prep Method: SW3520 Rev C

Prep Batch: EX081210-1

QCBatchID: EX081210-1-1

Run ID: SV081212-3

Cleanup: NONE

Basis: N/A

File Name: R9981

Sample Aliquot: 1000 ml

Final Volume: 1 ml

Result Units: UG/L

Clean DF: 1

86-74-8	CARBAZOLE	1	10	10	U	
84-74-2	DI-N-BUTYL PHTHALATE	1	10	10	U	
206-44-0	FLUORANTHENE	1	10	10	U	
129-00-0	PYRENE	1	10	10	U	
85-68-7	BUTYL BENZYL PHTHALATE	1	10	10	U	
56-55-3	BENZO(A)ANTHRACENE	1	10	10	U	
91-94-1	3,3'-DICHLOROBENZIDINE	1	10	10	U	
218-01-9	CHRYSENE	1	10	10	U	
117-81-7	BIS(2-ETHYLHEXYL)PHTHALATE	1	10	10	U	
117-84-0	DI-N-OCTYL PHTHALATE	1	10	10	U	
205-99-2	BENZO(B)FLUORANTHENE	1	10	10	U	
207-08-9	BENZO(K)FLUORANTHENE	1	10	10	U	
50-32-8	BENZO(A)PYRENE	1	10	10	U	
193-39-5	INDENO(1,2,3-CD)PYRENE	1	10	10	U	
53-70-3	DIBENZO(A,H)ANTHRACENE	1	10	10	U	
191-24-2	BENZO(G,H,I)PERYLENE	1	10	10	U	

### Surrogate Recovery

CASNO	Surrogate Analyte	Result	Flag	Spike Amount	Percent Recovery	Control Limits
118-79-6	2,4,6-TRIBROMOPHENOL	46.1		75	61	23 - 100
321-60-8	2-FLUOROBIPHENYL	42.5		50	85	21 - 106
367-12-4	2-FLUOROPHENOL	54.8		75	73	21 - 100
4165-60-0	NITROBENZENE-D5	40.3		50	81	34 - 111
4165-62-2	PHENOL-D5	55.2		75	74	15 - 104
1718-51-0	TERPHENYL-D14	51		50	102	33 - 111

Data Package ID: SV0812076-1

Date Printed: Tuesday, December 16, 2008

ALS Paragon

LIMS Version: 6.215A

Page 3 of 3

# GC/MS Semi-volatiles

## Method SW8270

### Tentatively Identified Compounds

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:   
Lab ID: EX081210-1MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 10-Dec-08

Date Analyzed: 12-Dec-08

Prep Batch: EX081210-1

QCBatchID: EX081210-1-1

Run ID: SV081212-3

Cleanup: NONE

Basis: As Received

Sample Aliquot: 1000 ml

Final Volume: 1 ml

Clean DF: 1

File Name: R9981

CASNO	Retention Time	Target Analyte	Dilution Factor	Result	Units	Qualifier
		NONE DETECTED	1			U

Data Package ID: SV0812076-1

Date Printed: Tuesday, December 16, 2008

ALS Paragon

LIMS Version: 6.215A

Page 2 of 2

# GC/MS Semi-volatiles

## Method SW8270D Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID: Angely WW	Sample Matrix: WATER	Prep Batch: EX081210-1	Sample Aliquot: 1060 ml
Lab ID: 0812076-1	% Moisture: N/A	QCBatchID: EX081210-1-1	Final Volume: 1 ml
	Date Collected: 08-Dec-08	Run ID: SV081212-3	Result Units: UG/L
	Date Extracted: 10-Dec-08	Cleanup: NONE	Clean DF: 1
	Date Analyzed: 12-Dec-08	Basis: As Received	
	Prep Method: SW3520 Rev C	File Name: R9984	

CASNO	Target Analyte	Dilution Factor	Result	Reporting Limit	Result Qualifier	EPA Qualifier
110-86-1	PYRIDINE	1	9.4	9.4	U	
62-75-9	N-NITROSODIMETHYLAMINE	1	9.4	9.4	U	
62-53-3	ANILINE	1	9.4	9.4	U	
108-95-2	PHENOL	1	9.4	9.4	U	
111-44-4	BIS(2-CHLOROETHYL)ETHER	1	9.4	9.4	U	
95-57-8	2-CHLOROPHENOL	1	9.4	9.4	U	
541-73-1	1,3-DICHLOROBENZENE	1	9.4	9.4	U	
106-46-7	1,4-DICHLOROBENZENE	1	9.4	9.4	U	
95-50-1	1,2-DICHLOROBENZENE	1	9.4	9.4	U	
100-51-6	BENZYL ALCOHOL	1	9.4	9.4	U	
108-60-1	BIS(2-CHLOROISOPROPYL)ETHER	1	9.4	9.4	U	
95-48-7	2-METHYLPHENOL	1	9.4	9.4	U	
621-64-7	N-NITROSO-DI-N-PROPYLAMINE	1	9.4	9.4	U	
108-39-4	3+4-METHYLPHENOL	1	9.4	9.4	U	
67-72-1	HEXACHLOROETHANE	1	9.4	9.4	U	
98-95-3	NITROBENZENE	1	9.4	9.4	U	
78-59-1	ISOPHORONE	1	9.4	9.4	U	
88-75-5	2-NITROPHENOL	1	9.4	9.4	U	
105-67-9	2,4-DIMETHYLPHENOL	1	9.4	9.4	U	
111-91-1	BIS(2-CHLOROETHOXY)METHANE	1	9.4	9.4	U	
120-83-2	2,4-DICHLOROPHENOL	1	9.4	9.4	U	
65-85-0	BENZOIC ACID	1	47	47	U	
120-82-1	1,2,4-TRICHLOROBENZENE	1	9.4	9.4	U	
91-20-3	NAPHTHALENE	1	9.4	9.4	U	
106-47-8	4-CHLOROANILINE	1	9.4	9.4	U	
87-68-3	HEXACHLOROBUTADIENE	1	9.4	9.4	U	
59-50-7	4-CHLORO-3-METHYLPHENOL	1	9.4	9.4	U	

Data Package ID: SV0812076-1

Date Printed: Tuesday, December 16, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 3

# GC/MS Semi-volatiles

## Method SW8270D Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:	Angely WW
Lab ID:	0812076-1

Sample Matrix: WATER  
% Moisture: N/A  
Date Collected: 08-Dec-08  
Date Extracted: 10-Dec-08  
Date Analyzed: 12-Dec-08  
Prep Method: SW3520 Rev C

Prep Batch: EX081210-1  
QCBatchID: EX081210-1-1  
Run ID: SV081212-3  
Cleanup: NONE  
Basis: As Received  
File Name: R9984

Sample Aliquot: 1060 ml  
Final Volume: 1 ml  
Result Units: UG/L  
Clean DF: 1

91-57-6	2-METHYLNAPHTHALENE	1	9.4	9.4	U	
90-12-0	1-METHYLNAPHTHALENE	1	9.4	9.4	U	
77-47-4	HEXACHLOROCYCLOPENTADIENE	1	9.4	9.4	U	
88-06-2	2,4,6-TRICHLOROPHENOL	1	9.4	9.4	U	
95-95-4	2,4,5-TRICHLOROPHENOL	1	9.4	9.4	U	
91-58-7	2-CHLORONAPHTHALENE	1	9.4	9.4	U	
88-74-4	2-NITROANILINE	1	19	19	U	
131-11-3	DIMETHYL PHTHALATE	1	9.4	9.4	U	
606-20-2	2,6-DINITROTOLUENE	1	9.4	9.4	U	
208-96-8	ACENAPHTHYLENE	1	9.4	9.4	U	
99-09-2	3-NITROANILINE	1	19	19	U	
83-32-9	ACENAPHTHENE	1	9.4	9.4	U	
51-28-5	2,4-DINITROPHENOL	1	19	19	U	
100-02-7	4-NITROPHENOL	1	19	19	U	
132-64-9	DIBENZOFURAN	1	9.4	9.4	U	
121-14-2	2,4-DINITROTOLUENE	1	9.4	9.4	U	
84-66-2	DIETHYL PHTHALATE	1	9.4	9.4	U	
86-73-7	FLUORENE	1	9.4	9.4	U	
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	1	9.4	9.4	U	
100-01-6	4-NITROANILINE	1	19	19	U	
103-33-3	AZOBENZENE	1	9.4	9.4	U	
534-52-1	4,6-DINITRO-2-METHYLPHENOL	1	19	19	U	
86-30-6	N-NITROSODIPHENYLAMINE	1	9.4	9.4	U	
101-55-3	4-BROMOPHENYL PHENYL ETHER	1	9.4	9.4	U	
118-74-1	HEXACHLOROBENZENE	1	9.4	9.4	U	
58-90-2	2,3,4,6-TETRACHLOROPHENOL	1	9.4	9.4	U	
87-86-5	PENTACHLOROPHENOL	1	19	19	U	
85-01-8	PHENANTHRENE	1	9.4	9.4	U	
120-12-7	ANTHRACENE	1	9.4	9.4	U	

Data Package ID: SV0812076-1

Date Printed: Tuesday, December 16, 2008

ALS Paragon

LIMS Version: 6.215A

Page 2 of 3

# GC/MS Semi-volatiles

## Method SW8270D Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID: Angely WW	Sample Matrix: WATER	Prep Batch: EX081210-1	Sample Aliquot: 1060 ml
Lab ID: 0812076-1	% Moisture: N/A	QCBatchID: EX081210-1-1	Final Volume: 1 ml
	Date Collected: 08-Dec-08	Run ID: SV081212-3	Result Units: UG/L
	Date Extracted: 10-Dec-08	Cleanup: NONE	Clean DF: 1
	Date Analyzed: 12-Dec-08	Basis: As Received	
	Prep Method: SW3520 Rev C	File Name: R9984	

86-74-8	CARBAZOLE	1	9.4	9.4	U	
84-74-2	DI-N-BUTYL PHTHALATE	1	9.4	9.4	U	
206-44-0	FLUORANTHENE	1	9.4	9.4	U	
129-00-0	PYRENE	1	9.4	9.4	U	
85-68-7	BUTYL BENZYL PHTHALATE	1	9.4	9.4	U	
56-55-3	BENZO(A)ANTHRACENE	1	9.4	9.4	U	
91-94-1	3,3'-DICHLOROBENZIDINE	1	9.4	9.4	U	
218-01-9	CHRYSENE	1	9.4	9.4	U	
117-81-7	BIS(2-ETHYLHEXYL)PHTHALATE	1	9.4	9.4	U	
117-84-0	DI-N-OCTYL PHTHALATE	1	9.4	9.4	U	
205-99-2	BENZO(B)FLUORANTHENE	1	9.4	9.4	U	
207-08-9	BENZO(K)FLUORANTHENE	1	9.4	9.4	U	
50-32-8	BENZO(A)PYRENE	1	9.4	9.4	U	
193-39-5	INDENO(1,2,3-CD)PYRENE	1	9.4	9.4	U	
53-70-3	DIBENZO(A,H)ANTHRACENE	1	9.4	9.4	U	
191-24-2	BENZO(G,H,I)PERYLENE	1	9.4	9.4	U	

## Surrogate Recovery

CASNO	Surrogate Analyte	Result	Flag	Spike Amount	Percent Recovery	Control Limits
118-79-6	2,4,6-TRIBROMOPHENOL	44.6		70.8	63	23 - 100
321-60-8	2-FLUOROBIPHENYL	30.9		47.2	65	21 - 106
367-12-4	2-FLUOROPHENOL	40.6		70.8	57	21 - 100
4165-60-0	NITROBENZENE-D5	29.4		47.2	62	34 - 111
4165-62-2	PHENOL-D5	41.4		70.8	59	15 - 104
1718-51-0	TERPHENYL-D14	36.3		47.2	77	33 - 111

Data Package ID: SV0812076-1

Date Printed: Tuesday, December 16, 2008

ALS Paragon

LIMS Version: 6.215A

Page 3 of 3

# GC/MS Semi-volatiles

## Method SW8270

### Tentatively Identified Compounds

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:	Angely WW
Lab ID:	0812076-1

Sample Matrix: WATER

% Moisture: N/A

Date Collected: 08-Dec-08

Date Extracted: 10-Dec-08

Date Analyzed: 12-Dec-08

Prep Batch: EX081210-1

QCBatchID: EX081210-1-1

Run ID: SV081212-3

Cleanup: NONE

Basis: As Received

Sample Aliquot: 1060 ml

Final Volume: 1 ml

Clean DF: 1

File Name: R9984

CASNO	Retention Time	Target Analyte	Dilution Factor	Result	Units	Qualifier
		NONE DETECTED	1			U

Data Package ID: SV0812076-1

Date Printed: Tuesday, December 16, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 2

# GC/MS Semi-volatiles

## Method SW8270D

### Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: EX081210-1LCS	Sample Matrix: WATER % Moisture: N/A Date Collected: N/A Date Extracted: 12/10/2008 Date Analyzed: 12/12/2008 Prep Method: SW3520C	Prep Batch: EX081210-1 QCBatchID: EX081210-1-1 Run ID: SV081212-3 Cleanup: NONE Basis: N/A File Name: R9982	Sample Aliquot: 1000 ml Final Volume: 1 ml Result Units: UG/L Clean DF: 1
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CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
110-86-1	PYRIDINE	60	38.6	10		64	10 - 108%
62-75-9	N-NITROSODIMETHYLAMINE	60	51.5	10		86	26 - 110%
62-53-3	ANILINE	60	44.4	10		74	25 - 125%
108-95-2	PHENOL	60	50.4	10		84	49 - 101%
111-44-4	BIS(2-CHLOROETHYL)ETHER	60	49.7	10		83	37 - 110%
95-57-8	2-CHLOROPHENOL	60	51.7	10		86	37 - 106%
541-73-1	1,3-DICHLOROBENZENE	60	46.3	10		77	32 - 98%
106-46-7	1,4-DICHLOROBENZENE	60	47.3	10		79	32 - 98%
95-50-1	1,2-DICHLOROBENZENE	60	48.3	10		81	33 - 102%
100-51-6	BENZYL ALCOHOL	60	50.4	10		84	30 - 112%
108-60-1	BIS(2-CHLOROISOPROPYL)ETHER	60	55.1	10		92	26 - 131%
95-48-7	2-METHYLPHENOL	60	50.2	10		84	38 - 109%
621-64-7	N-NITROSO-DI-N-PROPYLAMINE	60	54.8	10		91	34 - 128%
108-39-4	3+4-METHYLPHENOL	60	47.8	10		80	32 - 110%
67-72-1	HEXACHLOROETHANE	60	49.7	10		83	28 - 94%
98-95-3	NITROBENZENE	60	42.5	10		71	44 - 109%
78-59-1	ISOPHORONE	60	44.9	10		75	50 - 112%
88-75-5	2-NITROPHENOL	60	47.6	10		79	39 - 113%
105-67-9	2,4-DIMETHYLPHENOL	60	44	10		73	28 - 109%
111-91-1	BIS(2-CHLOROETHOXY)METHANE	60	43.3	10		72	46 - 107%
120-83-2	2,4-DICHLOROPHENOL	60	44.1	10		74	48 - 105%
65-85-0	BENZOIC ACID	100	51	50		51	10 - 125%
120-82-1	1,2,4-TRICHLOROBENZENE	60	39.9	10		67	37 - 107%
91-20-3	NAPHTHALENE	60	43.7	10		73	39 - 102%
106-47-8	4-CHLOROANILINE	60	39.8	10		66	15 - 109%
87-68-3	HEXACHLOROBUTADIENE	60	41.4	10		69	27 - 103%

Data Package ID: SV0812076-1

Date Printed: Tuesday, December 16, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 7

# GC/MS Semi-volatiles

## Method SW8270D

### Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: Paragon Analytics

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: EX081210-1LCS

Sample Matrix: WATER

% Moisture: N/A

Prep Batch: EX081210-1

Sample Aliquot: 1000 ml

QCBatchID: EX081210-1-1

Final Volume: 1 ml

Date Collected: N/A

Run ID: SV081212-3

Result Units: UG/L

Date Extracted: 12/10/2008

Cleanup: NONE

Clean DF: 1

Date Analyzed: 12/12/2008

Basis: N/A

Prep Method: SW3520C

File Name: R9982

CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
59-50-7	4-CHLORO-3-METHYLPHENOL	60	46.4	10		77	47 - 111%
91-57-6	2-METHYLNAPHTHALENE	60	41.2	10		69	46 - 104%
90-12-0	1-METHYLNAPHTHALENE	60	42.4	10		71	46 - 104%
77-47-4	HEXACHLOROCYCLOPENTADIENE	60	11	10		18	10 - 125%
88-06-2	2,4,6-TRICHLOROPHENOL	60	59.3	10		99	49 - 113%
95-95-4	2,4,5-TRICHLOROPHENOL	60	56.1	10		94	49 - 111%
91-58-7	2-CHLORONAPHTHALENE	60	52.9	10		88	36 - 137%
88-74-4	2-NITROANILINE	60	59.8	20		100	48 - 115%
131-11-3	DIMETHYL PHTHALATE	60	54.5	10		91	25 - 127%
606-20-2	2,6-DINITROTOLUENE	60	50.9	10		85	49 - 117%
208-96-8	ACENAPHTHYLENE	60	54.7	10		91	50 - 107%
99-09-2	3-NITROANILINE	60	51	20		85	19 - 126%
83-32-9	ACENAPHTHENE	60	53.2	10		89	47 - 108%
51-28-5	2,4-DINITROPHENOL	60	55.6	20		93	14 - 138%
100-02-7	4-NITROPHENOL	60	57.6	20		96	21 - 119%
132-64-9	DIBENZOFURAN	60	51.2	10		85	54 - 107%
121-14-2	2,4-DINITROTOLUENE	60	52.5	10		87	51 - 118%
84-66-2	DIETHYL PHTHALATE	60	57.6	10		96	41 - 118%
86-73-7	FLUORENE	60	56.2	10		94	50 - 112%
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	60	52.5	10		87	50 - 111%
100-01-6	4-NITROANILINE	60	52.2	20		87	36 - 118%
103-33-3	AZOBENZENE	60	56.8	10		95	21 - 137%
534-52-1	4,6-DINITRO-2-METHYLPHENOL	60	59.1	20		99	40 - 130%
86-30-6	N-NITROSODIPHENYLAMINE	60	45	10		75	48 - 111%
101-55-3	4-BROMOPHENYL PHENYL ETHER	60	52.6	10		88	52 - 113%
118-74-1	HEXACHLOROBENZENE	60	54.1	10		90	52 - 112%
58-90-2	2,3,4,6-TETRACHLOROPHENOL	100	94.4	10		94	23 - 112%

Data Package ID: SV0812076-1

Date Printed: Tuesday, December 16, 2008

ALS Paragon

LIMS Version: 6.215A

Page 2 of 7

# GC/MS Semi-volatiles

## Method SW8270D

### Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: Paragon Analytics

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: EX081210-1LCS

Sample Matrix: WATER

% Moisture: N/A

Prep Batch: EX081210-1

Sample Aliquot: 1000 ml

QCBatchID: EX081210-1-1

Final Volume: 1 ml

Date Collected: N/A

Run ID: SV081212-3

Result Units: UG/L

Date Extracted: 12/10/2008

Cleanup: NONE

Clean DF: 1

Date Analyzed: 12/12/2008

Basis: N/A

Prep Method: SW3520C

File Name: R9982

CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
87-86-5	PENTACHLOROPHENOL	60	57.6	20		96	38 - 117%
85-01-8	PHENANTHRENE	60	55.4	10		92	51 - 117%
120-12-7	ANTHRACENE	60	55.2	10		92	54 - 112%
86-74-8	CARBAZOLE	60	55.3	10		92	48 - 117%
84-74-2	DI-N-BUTYL PHTHALATE	60	64.1	10		107	54 - 116%
206-44-0	FLUORANTHENE	60	59.1	10		98	54 - 116%
129-00-0	PYRENE	60	42.2	10		70	49 - 128%
85-68-7	BUTYL BENZYL PHTHALATE	60	52.1	10		87	46 - 116%
56-55-3	BENZO(A)ANTHRACENE	60	50	10		83	56 - 109%
91-94-1	3,3'-DICHLOROBENZIDINE	60	41.9	10		70	19 - 111%
218-01-9	CHRYSENE	60	49.2	10		82	55 - 109%
117-81-7	BIS(2-ETHYLHEXYL)PHTHALATE	60	50.2	10		84	42 - 126%
117-84-0	DI-N-OCTYL PHTHALATE	60	55.9	10		93	37 - 137%
205-99-2	BENZO(B)FLUORANTHENE	60	53.3	10		89	45 - 118%
207-08-9	BENZO(K)FLUORANTHENE	60	51.6	10		86	45 - 124%
50-32-8	BENZO(A)PYRENE	60	48.6	10		81	53 - 110%
193-39-5	INDENO(1,2,3-CD)PYRENE	60	48.5	10		81	43 - 125%
53-70-3	DIBENZO(A,H)ANTHRACENE	60	51.4	10		86	42 - 127%
191-24-2	BENZO(G,H,I)PERYLENE	60	45.7	10		76	38 - 123%

Data Package ID: SV0812076-1

Date Printed: Tuesday, December 16, 2008

ALS Paragon

LIMS Version: 6.215A

Page 3 of 7

# GC/MS Semi-volatiles

## Method SW8270D

### Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: Paragon Analytics

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: EX081210-1LCSD

Sample Matrix: WATER

% Moisture: N/A

Prep Batch: EX081210-1

Sample Aliquot: 1000 ml

QCBatchID: EX081210-1-1

Final Volume: 1 ml

Date Collected: N/A

Run ID: SV081212-3

Result Units: UG/L

Date Extracted: 12/10/2008

Cleanup: NONE

Clean DF: 1

Date Analyzed: 12/12/2008

Basis: N/A

Prep Method: SW3520C

File Name: R9983

CASNO	Target Analyte	Spike Added	LCSD Result	Reporting Limit	Result Qualifier	LCSD % Rec.	RPD Limit	RPD
110-86-1	PYRIDINE	60	16.2	10	+	27	20	82
62-75-9	N-NITROSODIMETHYLAMINE	60	46.9	10		78	20	9
62-53-3	ANILINE	60	6.39	10	J*+	11	20	150
108-95-2	PHENOL	60	46.8	10		78	20	7
111-44-4	BIS(2-CHLOROETHYL)ETHER	60	45.2	10		75	20	10
95-57-8	2-CHLOROPHENOL	60	47.4	10		79	20	9
541-73-1	1,3-DICHLOROBENZENE	60	42.7	10		71	20	8
106-46-7	1,4-DICHLOROBENZENE	60	43.8	10		73	20	8
95-50-1	1,2-DICHLOROBENZENE	60	44.2	10		74	20	9
100-51-6	BENZYL ALCOHOL	60	46.9	10		78	20	7
108-60-1	BIS(2-CHLORoisopropyl)ETHER	60	50.7	10		84	20	8
95-48-7	2-METHYLPHENOL	60	45.3	10		76	20	10
621-64-7	N-NITROSO-DI-N-PROPYLAMINE	60	50.6	10		84	20	8
108-39-4	3+4-METHYLPHENOL	60	43.5	10		73	20	9
67-72-1	HEXACHLOROETHANE	60	45.4	10		76	20	9
98-95-3	NITROBENZENE	60	40.4	10		67	20	5
78-59-1	ISOPHORONE	60	43.1	10		72	20	4
88-75-5	2-NITROPHENOL	60	44.4	10		74	20	7
105-67-9	2,4-DIMETHYLPHENOL	60	40.1	10		67	20	9
111-91-1	BIS(2-CHLOROETHOXY)METHANE	60	40.8	10		68	20	6
120-83-2	2,4-DICHLOROPHENOL	60	41.3	10		69	20	7
65-85-0	BENZOIC ACID	100	54.7	50		55	20	7
120-82-1	1,2,4-TRICHLOROBENZENE	60	37.4	10		62	20	7
91-20-3	NAPHTHALENE	60	40.7	10		68	20	7
106-47-8	4-CHLOROANILINE	60	33.5	10		56	20	17
87-68-3	HEXACHLOROBUTADIENE	60	38.4	10		64	20	8
59-50-7	4-CHLORO-3-METHYLPHENOL	60	45.3	10		75	20	2

Data Package ID: SV0812076-1

Date Printed: Tuesday, December 16, 2008

ALS Paragon

LIMS Version: 6.215A

Page 4 of 7

# GC/MS Semi-volatiles

## Method SW8270D

### Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: Paragon Analytics

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: EX081210-1LCSD

Sample Matrix: WATER

% Moisture: N/A

Prep Batch: EX081210-1

QCBatchID: EX081210-1-1

Sample Aliquot: 1000 ml

Final Volume: 1 ml

Date Collected: N/A

Run ID: SV081212-3

Result Units: UG/L

Date Extracted: 12/10/2008

Cleanup: NONE

Clean DF: 1

Date Analyzed: 12/12/2008

Basis: N/A

Prep Method: SW3520C

File Name: R9983

CASNO	Target Analyte	Spike Added	LCSD Result	Reporting Limit	Result Qualifier	LCSD % Rec.	RPD Limit	RPD
91-57-6	2-METHYLNAPHTHALENE	60	39.1	10		65	20	5
90-12-0	1-METHYLNAPHTHALENE	60	40.7	10		68	20	4
77-47-4	HEXACHLOROCYCLOPENTADIENE	60	11.2	10		19	20	2
88-06-2	2,4,6-TRICHLOROPHENOL	60	57	10		95	20	4
95-95-4	2,4,5-TRICHLOROPHENOL	60	54.2	10		90	20	3
91-58-7	2-CHLORONAPHTHALENE	60	49.5	10		82	20	7
88-74-4	2-NITROANILINE	60	59.6	20		99	20	0
131-11-3	DIMETHYL PHTHALATE	60	54.3	10		90	20	0
606-20-2	2,6-DINITROTOLUENE	60	50.7	10		85	20	0
208-96-8	ACENAPHTHYLENE	60	52.2	10		87	20	5
99-09-2	3-NITROANILINE	60	51.1	20		85	20	0
83-32-9	ACENAPHTHENE	60	51.9	10		87	20	2
51-28-5	2,4-DINITROPHENOL	60	60.2	20		100	20	8
100-02-7	4-NITROPHENOL	60	58.3	20		97	20	1
132-64-9	DIBENZOFURAN	60	49.7	10		83	20	3
121-14-2	2,4-DINITROTOLUENE	60	51.9	10		87	20	1
84-66-2	DIETHYL PHTHALATE	60	56.7	10		95	20	2
86-73-7	FLUORENE	60	55.2	10		92	20	2
7005-72-3	4-CHLOROPHENYL PHENYL ETHER	60	50.9	10		85	20	3
100-01-6	4-NITROANILINE	60	51.6	20		86	20	1
103-33-3	AZOBENZENE	60	55.7	10		93	20	2
534-52-1	4,6-DINITRO-2-METHYLPHENOL	60	60.3	20		100	20	2
86-30-6	N-NITROSODIPHENYLAMINE	60	41.5	10		69	20	8
101-55-3	4-BROMOPHENYL PHENYL ETHER	60	51.9	10		87	20	1
118-74-1	HEXACHLOROBENZENE	60	52.7	10		88	20	3
58-90-2	2,3,4,6-TETRACHLOROPHENOL	100	92.9	10		93	20	2
87-86-5	PENTACHLOROPHENOL	60	56.7	20		94	20	2

Data Package ID: SV0812076-1

Date Printed: Tuesday, December 16, 2008

ALS Paragon

LIMS Version: 6.215A

Page 5 of 7

# GC/MS Semi-volatiles

## Method SW8270D

### Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: Paragon Analytics

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: EX081210-1LCSD

Sample Matrix: WATER

% Moisture: N/A

Prep Batch: EX081210-1

QCBatchID: EX081210-1-1

Sample Aliquot: 1000 ml

Final Volume: 1 ml

Date Collected: N/A

Run ID: SV081212-3

Result Units: UG/L

Date Extracted: 12/10/2008

Cleanup: NONE

Clean DF: 1

Date Analyzed: 12/12/2008

Basis: N/A

Prep Method: SW3520C

File Name: R9983

CASNO	Target Analyte	Spike Added	LCSD Result	Reporting Limit	Result Qualifier	LCSD % Rec.	RPD Limit	RPD
85-01-8	PHENANTHRENE	60	53.8	10		90	20	3
120-12-7	ANTHRACENE	60	53	10		88	20	4
86-74-8	CARBAZOLE	60	53.1	10		89	20	4
84-74-2	DI-N-BUTYL PHTHALATE	60	61.5	10		102	20	4
206-44-0	FLUORANTHENE	60	57.3	10		95	20	3
129-00-0	PYRENE	60	42.9	10		71	20	2
85-68-7	BUTYL BENZYL PHTHALATE	60	52.9	10		88	20	2
56-55-3	BENZO(A)ANTHRACENE	60	48.3	10		81	20	3
91-94-1	3,3'-DICHLOROBENZIDINE	60	26.6	10	+	44	20	45
218-01-9	CHRYSENE	60	47.4	10		79	20	4
117-81-7	BIS(2-ETHYLHEXYL)PHTHALATE	60	50.2	10		84	20	0
117-84-0	DI-N-OCTYL PHTHALATE	60	54.4	10		91	20	3
205-99-2	BENZO(B)FLUORANTHENE	60	51.2	10		85	20	4
207-08-9	BENZO(K)FLUORANTHENE	60	52.6	10		88	20	2
50-32-8	BENZO(A)PYRENE	60	45.4	10		76	20	7
193-39-5	INDENO(1,2,3-CD)PYRENE	60	47.8	10		80	20	1
53-70-3	DIBENZO(A,H)ANTHRACENE	60	50.8	10		85	20	1
191-24-2	BENZO(G,H,I)PERYLENE	60	45.3	10		76	20	1

Data Package ID: SV0812076-1

Date Printed: Tuesday, December 16, 2008

ALS Paragon

LIMS Version: 6.215A

Page 6 of 7

# GC/MS Semi-volatiles

Method SW8270D

## Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: Paragon Analytics

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

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### Surrogate Recovery LCS/LCSD

CASNO	Target Analyte	Spike Added	LCS % Rec.	LCS Flag	LCSD % Rec.	LCSD Flag	Control Limits
118-79-6	2,4,6-TRIBROMOPHENOL	75	82		82		23 - 100
321-60-8	2-FLUOROBIPHENYL	50	89		83		21 - 106
367-12-4	2-FLUOROPHENOL	75	80		73		21 - 100
4165-60-0	NITROBENZENE-D5	50	74		69		34 - 111
4165-62-2	PHENOL-D5	75	82		76		15 - 104
1718-51-0	TERPHENYL-D14	50	72		73		33 - 111

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Data Package ID: SV0812076-1

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Date Printed: Tuesday, December 16, 2008

ALS Paragon

LIMS Version: 6.215A

Page 7 of 7



# ALS Paragon



## GC/MS Volatiles Case Narrative

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### Colorado Oil & Gas Conservation Commission

Complaint 200199690

**Order Number - 0812076**

1. This report consists of 2 water samples. The samples were received cool and intact by ALS Paragon on 12/09/08. The vials for samples 08012076-1 and -2 contained headspace prior to analysis because they were not received headspace free into the volatiles laboratory.
2. These samples were prepared according to SW-846, 3rd Edition procedures. Specifically, the water samples were prepared by purging 10 mL using purge and trap procedures based on Method 5030C.
3. The samples were analyzed using GC/MS with an RTX-624, RTX-VMS, or equivalent capillary column according to SOP 525 Revision 12 based on SW-846 Method 8260B. All positive results were quantitated against the initial calibration standards using the internal standard technique. The identification of positive results was achieved by a comparison of the retention time and mass spectrum of the sample versus the daily calibration standard.
4. All initial calibration criteria for SPCC's and CCC's were met. If average response factors were used in the initial calibration, %RSD was  $\leq 15\%$ . If linear or higher order regression calibrations were used in the initial calibration, the coefficient of determination ( $r^2$ )  $\geq 0.99$ .
5. All initial calibrations are verified by comparing a second source standard calibration verification (ICV) against the calibration curve. All target compounds in the second source verification had a %D of less than 25%.
6. All criteria for SPCC's and CCC's were met in daily (continuing) calibration verifications (CCV).
7. Methylene chloride, acetone and 2-butanone are common laboratory contaminants. In order to minimize the levels of these compounds detected in the gc/ms analysis, ALS Paragon has designated its volatile laboratory as a restricted access area. In addition, the laboratory has been equipped with a dedicated, air intake and exhaust system that operates under positive pressure in order to minimize cross contamination of these compounds. Due to fluctuations in ambient laboratory conditions, reported sample values for common laboratory contaminants may be due to lab contamination even if the compound in question is not detected in the associated method blank.



All method blank criteria were met.

8. All laboratory control sample and laboratory control sample duplicate recoveries and RPDs were within the acceptance criteria with the following exception:

Spiked Compound	QC Sample	Direction
Bromoform	LCSD	High

The high recovery of this spike compound suggests that the quantitations of target analytes may be biased high. This analyte was not detected above the reporting limit in the associated samples. The reporting limits are defensible because the elevated recovery demonstrates an increase in sensitivity. No further action was taken.

9. Since a sample from this order number was not the selected quality control (QC) sample, matrix specific QC results are not included in this report.
10. The samples were analyzed within the established holding time.
11. All surrogate recoveries were within acceptance criteria.
12. All internal standard recoveries were within acceptance criteria.
13. Manual integrations are performed when needed to provide consistent and defensible data following the guidelines in SOP 939 Revision 3.

The data contained in the following report have been reviewed and approved by the personnel listed below. In addition, ALS Paragon certifies that the analyses reported herein are true, complete and correct within the limits of the methods employed.

Sel  
\_\_\_\_\_  
Sharon L. Jones  
Organics Primary Data Reviewer

12-15-08  
Date

S. D. White  
\_\_\_\_\_  
Organics Final Data Reviewer

12-15-08  
Date



**ALS Paragon**  
**Data Qualifier Flags**  
**Chromatography and Mass Spectrometry**

- U or ND:** This flag indicates that the compound was analyzed for but not detected.
- J:** This flag indicates an estimated value. This flag is used as follows : (1) when estimating a concentration for tentatively identified compounds (TICs) where a 1:1 response is assumed; (2) when the mass spectral and retention time data indicate the presence of a compound that meets the volatile and semivolatile GC/MS identification criteria, and the result is less than the reporting limit (RL) but greater than the method detection limit (MDL); (3) when the retention time data indicate the presence of a compound that meets the GC identification criteria, and the result is less than the RL but greater than the MDL; and (4) the reported value is estimated.
- B:** This flag is used when the analyte is detected in the associated method blank as well as in the sample. It indicates probable blank contamination and warns the data user. This flag shall be used for a tentatively identified compound (TIC) as well as for a positively identified target compound.
- E:** This flag identifies compounds whose concentration exceeds the upper level of the calibration range.
- A:** This flag indicates that a tentatively identified compound is a suspected aldol-condensation product.
- X:** This flag indicates that the analyte was diluted below an accurate quantitation level.
- \*:** This flag indicates that a spike recovery is equal to or outside the control criteria used.
- +:** This flag indicates that the relative percent difference (RPD) equals or exceeds the control criteria.

# ALS Paragon

## Sample Number(s) Cross-Reference Table

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**Paragon OrderNum:** 0812076

**Client Name:** Colorado Oil & Gas Conservation Commission

**Client Project Name:** Complaint 200199690

**Client Project Number:**

**Client PO Number:** OE PHA 09000000004

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Client Sample Number	Lab Sample Number	COC Number	Matrix	Date Collected	Time Collected
Angely WW	0812076-1		WATER	08-Dec-08	11:58
Trip Blank	0812076-2		WATER	08-Dec-08	



## CONDITION OF SAMPLE UPON RECEIPT FORM

Paragon Analytics

Client: COGCCWorkorder No: 0812076Project Manager: AWInitials: CDTDate: 12-9-08

1. Does this project require any <b>special handling</b> in addition to standard Paragon procedures?	YES	NO	
2. Are custody <b>seals on shipping containers</b> intact?	NONE	YES	NO
3. Are Custody seals on <b>sample containers</b> intact?	(NONE)	YES	NO
4. Is there a <b>COC (Chain-of-Custody) present</b> or other representative documents?	(YES)	YES	NO
5. Are the <b>COC and bottle labels complete and legible</b> ?	(YES)	YES	NO
6. Is the <b>COC in agreement</b> with samples received? (IDs, dates, times, no. of samples, no. of containers, matrix, requested analyses, etc.)	(YES)	YES	NO
7. Were <b>airbills / shipping documents</b> present and/or removable?	DROP OFF	YES	NO
8. Are all aqueous <b>samples requiring preservation</b> preserved correctly? (excluding volatiles)	N/A	YES	NO
9. Are all aqueous <b>non-preserved samples pH 4-9</b> ?	N/A	YES	NO
10. Is there <b>sufficient sample</b> for the requested analyses?	(YES)	YES	NO
11. Were all samples placed in the <b>proper containers</b> for the requested analyses?	(YES)	YES	NO
12. Are all samples within <b>holding times</b> for the requested analyses?	(YES)	YES	NO
13. Were all sample containers received <b>intact</b> ? (not broken or leaking, etc.)	(YES)	YES	NO
14. Are all samples requiring <b>no headspace</b> (VOC, GRO, RSK/MEE, Rx CN/S, radon) headspace free? Size of bubble: <u>&lt; green pea</u> ✓ <u>&gt; green pea</u>	N/A	YES	NO
15. Do perchlorate LCMS-MS samples <b>have headspace</b> ? (at least 1/3 of container required)	(N/A)	YES	NO
16. Were samples checked for and free from the presence of <b>residual chlorine</b> ? (Applicable when PM has indicated samples are from a chlorinated water source; note if field preservation with sodium thiosulfate was not observed.)	(N/A)	YES	NO
17. Were the samples <b>shipped on ice</b> ?	(YES)	YES	NO
18. Were cooler temperatures measured at 0.1-6.0°C?	IR gun used*: #2      #4 RAD ONLY	YES	NO
Cooler #:	<u>1</u>		
Temperature (°C):	<u>4.8</u>		
No. of custody seals on cooler:	<u>2</u>		
External µR/hr reading:	<u>13</u>		
Background µR/hr reading:	<u>10</u>		
Were external µR/hr readings ≤ two times background and within DOT acceptance criteria? <u>YES</u> NO / NA (If no. see Form 008.)			

Additional Information: PROVIDE DETAILS BELOW FOR A NO RESPONSE TO ANY QUESTION ABOVE, EXCEPT #1 AND #16.

HEADSPACE ALL > green pea    0812076-1-1    0812076-2-1

↓      -2  
 ↓      -3  
 ↓      -4  
 ↓      -5

If applicable, was the client contacted? YES NO / NA Contact: Peter ConstantasDate/Time: 12/9/08

e-mail

Project Manager Signature / Date: Connelly 12/9/08

# GC/MS Volatiles

## Method SW8260\_25B

### Method Blank

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: VL081210-4MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 10-Dec-08

Date Analyzed: 10-Dec-08

Prep Method: SW5030 Rev C

Prep Batch: VL081210-4

QCBatchID: VL081210-4-1

Run ID: VL081210-4A

Cleanup: NONE

Basis: N/A

File Name: D24936

Sample Aliquot: 10 ml

Final Volume: 10 ml

Result Units: UG/L

Clean DF: 1

CASNO	Target Analyte	DF	Result	Reporting Limit	Result Qualifier	EPA Qualifier
75-71-8	DICHLORODIFLUOROMETHANE	1	1	1	U	
74-87-3	CHLOROMETHANE	1	1	1	U	
75-01-4	VINYL CHLORIDE	1	1	1	U	
74-83-9	BROMOMETHANE	1	1	1	U	
75-00-3	CHLOROETHANE	1	1	1	U	
75-69-4	TRICHLOROFLUOROMETHANE	1	1	1	U	
75-35-4	1,1-DICHLOROETHENE	1	1	1	U	
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROET	1	1	1	U	
67-64-1	ACETONE	1	10	10	U	
74-88-4	IODOMETHANE	1	1	1	U	
75-15-0	CARBON DISULFIDE	1	1	1	U	
75-09-2	METHYLENE CHLORIDE	1	1	1	U	
156-60-5	TRANS-1,2-DICHLOROETHENE	1	1	1	U	
1634-04-4	METHYL TERTIARY BUTYL ETHER	1	1	1	U	
75-34-3	1,1-DICHLOROETHANE	1	1	1	U	
108-05-4	VINYL ACETATE	1	2	2	U	
156-59-2	CIS-1,2-DICHLOROETHENE	1	1	1	U	
78-93-3	2-BUTANONE	1	10	10	U	
74-97-5	BROMOCHLOROMETHANE	1	1	1	U	
67-66-3	CHLOROFORM	1	1	1	U	
71-55-6	1,1,1-TRICHLOROETHANE	1	1	1	U	
594-20-7	2,2-DICHLOROPROPANE	1	1	1	U	
56-23-5	CARBON TETRACHLORIDE	1	1	1	U	
563-58-6	1,1-DICHLOROPROPENE	1	1	1	U	
107-06-2	1,2-DICHLOROETHANE	1	1	1	U	
71-43-2	BENZENE	1	1	1	U	
79-01-6	TRICHLOROETHENE	1	1	1	U	

Data Package ID: VL0812076-1

Date Printed: Thursday, December 11, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 3

# GC/MS Volatiles

## Method SW8260\_25B

### Method Blank

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: VL081210-4MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 10-Dec-08

Date Analyzed: 10-Dec-08

Prep Method: SW5030 Rev C

Prep Batch: VL081210-4

QCBatchID: VL081210-4-1

Run ID: VL081210-4A

Cleanup: NONE

Basis: N/A

File Name: D24936

Sample Aliquot: 10 ml

Final Volume: 10 ml

Result Units: UG/L

Clean DF: 1

78-87-5	1,2-DICHLOROPROPANE	1	1	1	U	
74-95-3	DIBROMOMETHANE	1	1	1	U	
75-27-4	BROMODICHLOROMETHANE	1	1	1	U	
10061-01-5	CIS-1,3-DICHLOROPROPENE	1	1	1	U	
108-10-1	4-METHYL-2-PENTANONE	1	10	10	U	
108-88-3	TOLUENE	1	1	1	U	
10061-02-6	TRANS-1,3-DICHLOROPROPENE	1	1	1	U	
79-00-5	1,1,2-TRICHLOROETHANE	1	1	1	U	
591-78-6	2-HEXANONE	1	10	10	U	
127-18-4	TETRACHLOROETHENE	1	1	1	U	
142-28-9	1,3-DICHLOROPROPANE	1	1	1	U	
124-48-1	DIBROMOCHLOROMETHANE	1	1	1	U	
106-93-4	1,2-DIBROMOETHANE	1	1	1	U	
544-10-5	1-CHLOROHEXANE	1	1	1	U	
108-90-7	CHLOROBENZENE	1	1	1	U	
630-20-6	1,1,1,2-TETRACHLOROETHANE	1	1	1	U	
100-41-4	ETHYLBENZENE	1	1	1	U	
136777-61-2	M+P-XYLENE	1	1	1	U	
95-47-6	O-XYLENE	1	1	1	U	
100-42-5	STYRENE	1	1	1	U	
75-25-2	BROMOFORM	1	1	1	U	
98-82-8	ISOPROPYLBENZENE	1	1	1	U	
96-18-4	1,2,3-TRICHLOROPROPANE	1	1	1	U	
79-34-5	1,1,2,2-TETRACHLOROETHANE	1	1	1	U	
108-86-1	BROMOBENZENE	1	1	1	U	
103-65-1	N-PROPYLBENZENE	1	1	1	U	
95-49-8	2-CHLOROTOLUENE	1	1	1	U	
108-67-8	1,3,5-TRIMETHYLBENZENE	1	1	1	U	
106-43-4	4-CHLOROTOLUENE	1	1	1	U	

Data Package ID: VL0812076-1

Date Printed: Thursday, December 11, 2008

ALS Paragon

LIMS Version: 6.215A

Page 2 of 3

# GC/MS Volatiles

## Method SW8260\_25B

### Method Blank

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: VL081210-4MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 10-Dec-08

Date Analyzed: 10-Dec-08

Prep Method: SW5030 Rev C

Prep Batch: VL081210-4

QCBatchID: VL081210-4-1

Run ID: VL081210-4A

Cleanup: NONE

Basis: N/A

File Name: D24936

Sample Aliquot: 10 ml

Final Volume: 10 ml

Result Units: UG/L

Clean DF: 1

98-06-6	TERT-BUTYLBENZENE	1	1	1	U	
95-63-6	1,2,4-TRIMETHYLBENZENE	1	1	1	U	
135-98-8	SEC-BUTYLBENZENE	1	1	1	U	
541-73-1	1,3-DICHLOROBENZENE	1	1	1	U	
99-87-6	P-ISOPROPYLtoluene	1	1	1	U	
106-46-7	1,4-DICHLOROBENZENE	1	1	1	U	
104-51-8	N-BUTYLBENZENE	1	1	1	U	
95-50-1	1,2-DICHLOROBENZENE	1	1	1	U	
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	1	2	2	U	
120-82-1	1,2,4-TRICHLOROBENZENE	1	1	1	U	
87-68-3	HEXAChlorobutadiene	1	1	1	U	
91-20-3	NAPHTHALENE	1	1	1	U	
87-61-6	1,2,3-TRICHLOROBENZENE	1	1	1	U	

### Surrogate Recovery

CASNO	Surrogate Analyte	Result	Flag	Spike Amount	Percent Recovery	Control Limits
460-00-4	4-BROMOFLUOROBENZENE	26.8		25	107	78 - 129
1868-53-7	DIBROMOFLUOROMETHANE	23.9		25	95	80 - 124
2037-26-5	TOLUENE-D8	24.2		25	97	81 - 119

Data Package ID: VL0812076-1

Date Printed: Thursday, December 11, 2008

ALS Paragon

LIMS Version: 6.215A

Page 3 of 3

# GC/MS Volatiles

## Method SW8260\_25 Tentatively Identified Compounds

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:	
Lab ID:	VL081210-4MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 10-Dec-08

Date Analyzed: 10-Dec-08

Prep Batch: VL081210-4

QCBatchID: VL081210-4-1

Run ID: VL081210-4A

Cleanup: NONE

Basis: As Received

Sample Aliquot: 10 ml

Final Volume: 10 ml

Clean DF: 1

File Name: D24936

CASNO	Retention Time	Target Analyte	Dilution Factor	Result	Units	Qualifier
		NONE DETECTED	1			U

Data Package ID: VL0812076-1

Date Printed: Friday, December 19, 2008

ALS Paragon

LIMS Version: 6.217A

Page 3 of 3

# GC/MS Volatiles

## Method SW8260\_25B

### Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:	Angely WW
Lab ID:	0812076-1

Sample Matrix: WATER  
% Moisture: N/A  
Prep Batch: VL081210-4  
QCBatchID: VL081210-4-1  
Date Collected: 08-Dec-08  
Run ID: VL081210-4A  
Date Extracted: 10-Dec-08  
Cleanup: NONE  
Date Analyzed: 10-Dec-08  
Basis: As Received  
Prep Method: SW5030 Rev C  
File Name: D24938  
Final Volume: 10 ml  
Result Units: UG/L  
Clean DF: 1

CASNO	Target Analyte	Dilution Factor	Result	Reporting Limit	Result Qualifier	EPA Qualifier
75-71-8	DICHLORODIFLUOROMETHANE	1	1	1	U	
74-87-3	CHLOROMETHANE	1	1	1	U	
75-01-4	VINYL CHLORIDE	1	1	1	U	
74-83-9	BROMOMETHANE	1	1	1	U	
75-00-3	CHLOROETHANE	1	1	1	U	
75-69-4	TRICHLOROFUOROMETHANE	1	1	1	U	
75-35-4	1,1-DICHLOROETHENE	1	1	1	U	
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	1	1	1	U	
67-64-1	ACETONE	1	10	10	U	
74-88-4	IODOMETHANE	1	1	1	U	
75-15-0	CARBON DISULFIDE	1	1	1	U	
75-09-2	METHYLENE CHLORIDE	1	1	1	U	
156-60-5	TRANS-1,2-DICHLOROETHENE	1	1	1	U	
1634-04-4	METHYL TERTIARY BUTYL ETHER	1	1	1	U	
75-34-3	1,1-DICHLOROETHANE	1	1	1	U	
108-05-4	VINYL ACETATE	1	2	2	U	
156-59-2	CIS-1,2-DICHLOROETHENE	1	1	1	U	
78-93-3	2-BUTANONE	1	10	10	U	
74-97-5	BROMOCHLOROMETHANE	1	1	1	U	
67-66-3	CHLOROFORM	1	1	1	U	
71-55-6	1,1,1-TRICHLOROETHANE	1	1	1	U	
594-20-7	2,2-DICHLOROPROPANE	1	1	1	U	
56-23-5	CARBON TETRACHLORIDE	1	1	1	U	
563-58-6	1,1-DICHLOROPROPENE	1	1	1	U	
107-06-2	1,2-DICHLOROETHANE	1	1	1	U	
71-43-2	BENZENE	1	1	1	U	
79-01-6	TRICHLOROETHENE	1	1	1	U	

Data Package ID: VL0812076-1

Date Printed: Thursday, December 11, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 6

# GC/MS Volatiles

## Method SW8260\_25B

### Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID: Angely WW	Sample Matrix: WATER	Prep Batch: VL081210-4	Sample Aliquot: 10 ml
Lab ID: 0812076-1	% Moisture: N/A	QCBatchID: VL081210-4-1	Final Volume: 10 ml
	Date Collected: 08-Dec-08	Run ID: VL081210-4A	Result Units: UG/L
	Date Extracted: 10-Dec-08	Cleanup: NONE	Clean DF: 1
	Date Analyzed: 10-Dec-08	Basis: As Received	
	Prep Method: SW5030 Rev C	File Name: D24938	

78-87-5	1,2-DICHLOROPROPANE	1	1	1	U	
74-95-3	DIBROMOMETHANE	1	1	1	U	
75-27-4	BROMODICHLOROMETHANE	1	1	1	U	
10061-01-5	CIS-1,3-DICHLOROPROPENE	1	1	1	U	
108-10-1	4-METHYL-2-PENTANONE	1	10	10	U	
108-88-3	TOLUENE	1	1	1	U	
10061-02-6	TRANS-1,3-DICHLOROPROPENE	1	1	1	U	
79-00-5	1,1,2-TRICHLOROETHANE	1	1	1	U	
591-78-6	2-HEXANONE	1	10	10	U	
127-18-4	TETRACHLOROETHENE	1	1	1	U	
142-28-9	1,3-DICHLOROPROPANE	1	1	1	U	
124-48-1	DIBROMOCHLOROMETHANE	1	1	1	U	
106-93-4	1,2-DIBROMOETHANE	1	1	1	U	
544-10-5	1-CHLOROHEXANE	1	1	1	U	
108-90-7	CHLOROBENZENE	1	1	1	U	
630-20-6	1,1,1,2-TETRACHLOROETHANE	1	1	1	U	
100-41-4	ETHYLBENZENE	1	1	1	U	
136777-61-2	M+P-XYLENE	1	1	1	U	
95-47-6	O-XYLENE	1	1	1	U	
100-42-5	STYRENE	1	1	1	U	
75-25-2	BROMOFORM	1	1	1	U	
98-82-8	ISOPROPYLBENZENE	1	1	1	U	
96-18-4	1,2,3-TRICHLOROPROPANE	1	1	1	U	
79-34-5	1,1,2,2-TETRACHLOROETHANE	1	1	1	U	
108-86-1	BROMOBENZENE	1	1	1	U	
103-65-1	N-PROPYLBENZENE	1	1	1	U	
95-49-8	2-CHLOROTOLUENE	1	1	1	U	
108-67-8	1,3,5-TRIMETHYLBENZENE	1	1	1	U	
106-43-4	4-CHLOROTOLUENE	1	1	1	U	

Data Package ID: VL0812076-1

Date Printed: Thursday, December 11, 2008

ALS Paragon

LIMS Version: 6.215A

Page 2 of 6

# GC/MS Volatiles

## Method SW8260\_25B

### Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:	Angely WW
Lab ID:	0812076-1

Sample Matrix: WATER  
% Moisture: N/A  
Prep Batch: VL081210-4  
QCBatchID: VL081210-4-1  
Date Collected: 08-Dec-08  
Run ID: VL081210-4A  
Date Extracted: 10-Dec-08  
Cleanup: NONE  
Date Analyzed: 10-Dec-08  
Basis: As Received  
Prep Method: SW5030 Rev C  
File Name: D24938  
Sample Aliquot: 10 ml  
Final Volume: 10 ml  
Result Units: UG/L  
Clean DF: 1

98-06-6	TERT-BUTYLBENZENE	1	1	1	U	
95-63-6	1,2,4-TRIMETHYLBENZENE	1	1	1	U	
135-98-8	SEC-BUTYLBENZENE	1	1	1	U	
541-73-1	1,3-DICHLOROBENZENE	1	1	1	U	
99-87-6	P-ISOPROPYLtolUENE	1	1	1	U	
106-46-7	1,4-DICHLOROBENZENE	1	1	1	U	
104-51-8	N-BUTYLBENZENE	1	1	1	U	
95-50-1	1,2-DICHLOROBENZENE	1	1	1	U	
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	1	2	2	U	
120-82-1	1,2,4-TRICHLOROBENZENE	1	1	1	U	
87-68-3	HEXACHLOROBUTADIENE	1	1	1	U	
91-20-3	NAPHTHALENE	1	1	1	U	
87-61-6	1,2,3-TRICHLOROBENZENE	1	1	1	U	

### Surrogate Recovery

CASNO	Surrogate Analyte	Result	Flag	Spike Amount	Percent Recovery	Control Limits
460-00-4	4-BROMOFLUOROBENZENE	26.6		25	106	78 - 129
1868-53-7	DIBROMOFLUOROMETHANE	24.1		25	96	80 - 124
2037-26-5	TOLUENE-D8	25		25	100	81 - 119

Data Package ID: VL0812076-1

Date Printed: Thursday, December 11, 2008

ALS Paragon

LIMS Version: 6.215A

Page 3 of 6

# GC/MS Volatiles

## Method SW8260\_25 Tentatively Identified Compounds

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:	Angely WW
Lab ID:	0812076-1

Sample Matrix: WATER

% Moisture: N/A

Date Collected: 08-Dec-08

Date Extracted: 10-Dec-08

Date Analyzed: 10-Dec-08

Prep Batch: VL081210-4

QCBatchID: VL081210-4-1

Run ID: VL081210-4A

Cleanup: NONE

Basis: As Received

Sample Aliquot: 10 ml

Final Volume: 10 ml

Clean DF: 1

File Name: D24938

CASNO	Retention Time	Target Analyte	Dilution Factor	Result	Units	Qualifier
		NONE DETECTED	1			U

Data Package ID: VL0812076-1

Date Printed: Friday, December 19, 2008

ALS Paragon

LIMS Version: 6.217A

Page 1 of 3

# GC/MS Volatiles

## Method SW8260\_25B

### Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID: Trip Blank	Sample Matrix: WATER	Prep Batch: VL081210-4	Sample Aliquot: 10 ml
Lab ID: 0812076-2	% Moisture: N/A	QCBatchID: VL081210-4-1	Final Volume: 10 ml
	Date Collected: 08-Dec-08	Run ID: VL081210-4A	Result Units: UG/L
	Date Extracted: 10-Dec-08	Cleanup: NONE	Clean DF: 1
	Date Analyzed: 10-Dec-08	Basis: As Received	
	Prep Method: SW5030 Rev C	File Name: D24937	

CASNO	Target Analyte	Dilution Factor	Result	Reporting Limit	Result Qualifier	EPA Qualifier
75-71-8	DICHLORODIFLUOROMETHANE	1	1	1	U	
74-87-3	CHLOROMETHANE	1	1	1	U	
75-01-4	VINYL CHLORIDE	1	1	1	U	
74-83-9	BROMOMETHANE	1	1	1	U	
75-00-3	CHLOROETHANE	1	1	1	U	
75-69-4	TRICHLOROFUOROMETHANE	1	1	1	U	
75-35-4	1,1-DICHLOROETHENE	1	1	1	U	
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	1	1	1	U	
67-64-1	ACETONE	1	10	10	U	
74-88-4	IODOMETHANE	1	1	1	U	
75-15-0	CARBON DISULFIDE	1	1	1	U	
75-09-2	METHYLENE CHLORIDE	1	0.62	1	J	
156-60-5	TRANS-1,2-DICHLOROETHENE	1	1	1	U	
1634-04-4	METHYL TERTIARY BUTYL ETHER	1	1	1	U	
75-34-3	1,1-DICHLOROETHANE	1	1	1	U	
108-05-4	VINYL ACETATE	1	2	2	U	
156-59-2	CIS-1,2-DICHLOROETHENE	1	1	1	U	
78-93-3	2-BUTANONE	1	10	10	U	
74-97-5	BROMOCHLOROMETHANE	1	1	1	U	
67-66-3	CHLOROFORM	1	1	1	U	
71-55-6	1,1,1-TRICHLOROETHANE	1	1	1	U	
594-20-7	2,2-DICHLOROPROPANE	1	1	1	U	
56-23-5	CARBON TETRACHLORIDE	1	1	1	U	
563-58-6	1,1-DICHLOROPROPENE	1	1	1	U	
107-06-2	1,2-DICHLOROETHANE	1	1	1	U	
71-43-2	BENZENE	1	1	1	U	
79-01-6	TRICHLOROETHENE	1	1	1	U	

Data Package ID: VL0812076-1

Date Printed: Thursday, December 11, 2008

ALS Paragon

LIMS Version: 6.215A

Page 4 of 6

# GC/MS Volatiles

## Method SW8260\_25B

### Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:	Trip Blank
Lab ID:	0812076-2

Sample Matrix: WATER  
% Moisture: N/A  
Date Collected: 08-Dec-08  
Date Extracted: 10-Dec-08  
Date Analyzed: 10-Dec-08  
Prep Method: SW5030 Rev C

Prep Batch: VL081210-4  
QCBatchID: VL081210-4-1  
Run ID: VL081210-4A  
Cleanup: NONE  
Basis: As Received  
File Name: D24937

Sample Aliquot: 10 ml  
Final Volume: 10 ml  
Result Units: UG/L  
Clean DF: 1

78-87-5	1,2-DICHLOROPROPANE	1	1	1	U	
74-95-3	DIBROMOMETHANE	1	1	1	U	
75-27-4	BROMODICHLOROMETHANE	1	1	1	U	
10061-01-5	CIS-1,3-DICHLOROPROPENE	1	1	1	U	
108-10-1	4-METHYL-2-PENTANONE	1	10	10	U	
108-88-3	TOLUENE	1	1	1	U	
10061-02-6	TRANS-1,3-DICHLOROPROPENE	1	1	1	U	
79-00-5	1,1,2-TRICHLOROETHANE	1	1	1	U	
591-78-6	2-HEXANONE	1	10	10	U	
127-18-4	TETRACHLOROETHENE	1	1	1	U	
142-28-9	1,3-DICHLOROPROPANE	1	1	1	U	
124-48-1	DIBROMOCHLOROMETHANE	1	1	1	U	
106-93-4	1,2-DIBROMOETHANE	1	1	1	U	
544-10-5	1-CHLOROHEXANE	1	1	1	U	
108-90-7	CHLOROBENZENE	1	1	1	U	
630-20-6	1,1,1,2-TETRACHLOROETHANE	1	1	1	U	
100-41-4	ETHYLBENZENE	1	1	1	U	
136777-61-2	M+P-XYLENE	1	1	1	U	
95-47-6	O-XYLENE	1	1	1	U	
100-42-5	STYRENE	1	1	1	U	
75-25-2	BROMOFORM	1	1	1	U	
98-82-8	ISOPROPYLBENZENE	1	1	1	U	
96-18-4	1,2,3-TRICHLOROPROPANE	1	1	1	U	
79-34-5	1,1,2,2-TETRACHLOROETHANE	1	1	1	U	
108-86-1	BROMOBENZENE	1	1	1	U	
103-65-1	N-PROPYLBENZENE	1	1	1	U	
95-49-8	2-CHLOROTOLUENE	1	1	1	U	
108-67-8	1,3,5-TRIMETHYLBENZENE	1	1	1	U	
106-43-4	4-CHLOROTOLUENE	1	1	1	U	

Data Package ID: VL0812076-1

Date Printed: Thursday, December 11, 2008

ALS Paragon

LIMS Version: 6.215A

Page 5 of 6

# GC/MS Volatiles

## Method SW8260\_25B

### Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:	Trip Blank
Lab ID:	0812076-2

Sample Matrix: WATER  
% Moisture: N/A  
Date Collected: 08-Dec-08  
Date Extracted: 10-Dec-08  
Date Analyzed: 10-Dec-08  
Prep Method: SW5030 Rev C

Prep Batch: VL081210-4  
QCBatchID: VL081210-4-1  
Run ID: VL081210-4A  
Cleanup: NONE  
Basis: As Received  
File Name: D24937

Sample Aliquot: 10 ml  
Final Volume: 10 ml  
Result Units: UG/L  
Clean DF: 1

98-06-6	TERT-BUTYLBENZENE	1	1	1	U	
95-63-6	1,2,4-TRIMETHYLBENZENE	1	1	1	U	
135-98-8	SEC-BUTYLBENZENE	1	1	1	U	
541-73-1	1,3-DICHLOROBENZENE	1	1	1	U	
99-87-6	P-ISOPROPYLtolUENE	1	1	1	U	
106-46-7	1,4-DICHLOROBENZENE	1	1	1	U	
104-51-8	N-BUTYLBENZENE	1	1	1	U	
95-50-1	1,2-DICHLOROBENZENE	1	1	1	U	
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	1	2	2	U	
120-82-1	1,2,4-TRICHLOROBENZENE	1	1	1	U	
87-68-3	HEXACHLOROBUTADIENE	1	1	1	U	
91-20-3	NAPHTHALENE	1	1	1	U	
87-61-6	1,2,3-TRICHLOROBENZENE	1	1	1	U	

### Surrogate Recovery

CASNO	Surrogate Analyte	Result	Flag	Spike Amount	Percent Recovery	Control Limits
460-00-4	4-BROMOFLUOROBENZENE	26.4		25	105	78 - 129
1868-53-7	DIBROMOFLUOROMETHANE	24.2		25	97	80 - 124
2037-26-5	TOLUENE-D8	24.8		25	99	81 - 119

Data Package ID: VL0812076-1

Date Printed: Thursday, December 11, 2008

ALS Paragon

LIMS Version: 6.215A

Page 6 of 6

# GC/MS Volatiles

## Method SW8260\_25 Tentatively Identified Compounds

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:	Trip Blank
Lab ID:	0812076-2

Sample Matrix: WATER

% Moisture: N/A

Date Collected: 08-Dec-08

Date Extracted: 10-Dec-08

Date Analyzed: 10-Dec-08

Prep Batch: VL081210-4

QCBatchID: VL081210-4-1

Run ID: VL081210-4A

Cleanup: NONE

Basis: As Received

Sample Aliquot: 10 ml

Final Volume: 10 ml

Clean DF: 1

File Name: D24937

CASNO	Retention Time	Target Analyte	Dilution Factor	Result	Units	Qualifier
		NONE DETECTED	1			U

Data Package ID: VL0812076-1

Date Printed: Friday, December 19, 2008

ALS Paragon

LIMS Version: 6.217A

Page 2 of 3

# GC/MS Volatiles

## Method SW8260\_25B

### Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: VL081210-4LCS	Sample Matrix: WATER % Moisture: N/A Date Collected: N/A Date Extracted: 12/10/2008 Date Analyzed: 12/10/2008 Prep Method: SW5030C	Prep Batch: VL081210-4 QCBatchID: VL081210-4-1 Run ID: VL081210-4A Cleanup: NONE Basis: N/A File Name: D24934	Sample Aliquot: 10 ml Final Volume: 10 ml Result Units: UG/L Clean DF: 1
-----------------------	---	--	---

CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
75-71-8	DICHLORODIFLUOROMETHANE	10	11	1		110	38 - 131%
74-87-3	CHLOROMETHANE	10	10.1	1		101	62 - 141%
75-01-4	VINYL CHLORIDE	10	10.4	1		104	77 - 124%
74-83-9	BROMOMETHANE	10	10.3	1		103	76 - 133%
75-00-3	CHLOROETHANE	10	11.4	1		114	81 - 130%
75-69-4	TRICHLORODIFLUOROMETHANE	10	11	1		110	84 - 146%
75-35-4	1,1-DICHLOROETHENE	10	10	1		100	75 - 126%
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	10	9.75	1		98	71 - 144%
67-64-1	ACETONE	40	39.4	10		99	50 - 150%
74-88-4	IODOMETHANE	10	11.3	1		113	76 - 116%
75-15-0	CARBON DISULFIDE	10	10.4	1		104	68 - 129%
75-09-2	METHYLENE CHLORIDE	10	9.87	1		99	22 - 146%
156-60-5	TRANS-1,2-DICHLOROETHENE	10	10.3	1		103	76 - 135%
1634-04-4	METHYL TERTIARY BUTYL ETHER	20	18.6	1		93	75 - 125%
75-34-3	1,1-DICHLOROETHANE	10	9.7	1		97	77 - 131%
108-05-4	VINYL ACETATE	10	9.63	2		96	56 - 151%
156-59-2	CIS-1,2-DICHLOROETHENE	10	10.1	1		101	81 - 121%
78-93-3	2-BUTANONE	40	33.8	10		85	50 - 150%
74-97-5	BROMOCHLOROMETHANE	10	10.1	1		101	85 - 126%
67-66-3	CHLOROFORM	10	9.83	1		98	84 - 125%
71-55-6	1,1,1-TRICHLOROETHANE	10	10.1	1		101	82 - 129%
594-20-7	2,2-DICHLOROPROPANE	10	9.94	1		99	79 - 130%
56-23-5	CARBON TETRACHLORIDE	10	10.1	1		101	83 - 135%
563-58-6	1,1-DICHLOROPROPENE	10	9.88	1		99	85 - 127%
107-06-2	1,2-DICHLOROETHANE	10	9.41	1		94	84 - 126%
71-43-2	BENZENE	10	9.73	1		97	82 - 122%

Data Package ID: VL0812076-1

Date Printed: Thursday, December 11, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 6

# GC/MS Volatiles

## Method SW8260\_25B

### Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: Paragon Analytics

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: VL081210-4LCS

Sample Matrix: WATER

% Moisture: N/A

Prep Batch: VL081210-4

QCBatchID: VL081210-4-1

Sample Aliquot: 10 ml

Final Volume: 10 ml

Date Collected: N/A

Run ID: VL081210-4A

Result Units: UG/L

Date Extracted: 12/10/2008

Cleanup: NONE

Clean DF: 1

Date Analyzed: 12/10/2008

Basis: N/A

Prep Method: SW5030C

File Name: D24934

CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
79-01-6	TRICHLOROETHENE	10	10.1	1		101	82 - 121%
78-87-5	1,2-DICHLOROPROPANE	10	9.39	1		94	81 - 121%
74-95-3	DIBROMOMETHANE	10	9.42	1		94	81 - 125%
75-27-4	BROMODICHLOROMETHANE	10	9.54	1		95	82 - 120%
10061-01-5	CIS-1,3-DICHLOROPROPENE	10	9.66	1		97	79 - 120%
108-10-1	4-METHYL-2-PENTANONE	40	34.8	10		87	50 - 150%
108-88-3	TOLUENE	10	9.88	1		99	83 - 121%
10061-02-6	TRANS-1,3-DICHLOROPROPENE	10	9.09	1		91	78 - 113%
79-00-5	1,1,2-TRICHLOROETHANE	10	9.05	1		90	82 - 122%
591-78-6	2-HEXANONE	40	33.5	10		84	50 - 150%
127-18-4	TETRACHLOROETHENE	10	11.2	1		112	79 - 136%
142-28-9	1,3-DICHLOROPROPANE	10	9.01	1		90	80 - 126%
124-48-1	DIBROMOCHLOROMETHANE	10	9.84	1		98	80 - 123%
106-93-4	1,2-DIBROMOETHANE	10	9.39	1		94	85 - 124%
544-10-5	1-CHLOROHEXANE	10	10	1		100	77 - 135%
108-90-7	CHLOROBENZENE	10	9.9	1		99	82 - 121%
630-20-6	1,1,2-TETRACHLOROETHANE	10	9.87	1		99	85 - 128%
100-41-4	ETHYLBENZENE	10	9.82	1		98	83 - 126%
136777-61-	M+P-XYLENE	20	20.2	1		101	82 - 129%
95-47-6	O-XYLENE	10	10.2	1		102	87 - 132%
100-42-5	STYRENE	10	10.1	1		101	82 - 123%
75-25-2	BROMOFORM	10	11.2	1		112	79 - 118%
98-82-8	ISOPROPYLBENZENE	10	10.4	1		104	75 - 132%
96-18-4	1,2,3-TRICHLOROPROPANE	10	8.97	1		90	77 - 128%
79-34-5	1,1,2,2-TETRACHLOROETHANE	10	8.38	1		84	74 - 130%
108-86-1	BROMOBENZENE	10	10	1		100	78 - 124%
103-65-1	N-PROPYLBENZENE	10	9.37	1		94	75 - 134%

Data Package ID: VL0812076-1

Date Printed: Thursday, December 11, 2008

ALS Paragon

LIMS Version: 6.215A

Page 2 of 6

# GC/MS Volatiles

## Method SW8260\_25B

### Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: Paragon Analytics

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: VL081210-4LCS

Sample Matrix: WATER

% Moisture: N/A

Prep Batch: VL081210-4

QCBatchID: VL081210-4-1

Sample Aliquot: 10 ml

Final Volume: 10 ml

Date Collected: N/A

Run ID: VL081210-4A

Result Units: UG/L

Date Extracted: 12/10/2008

Cleanup: NONE

Clean DF: 1

Date Analyzed: 12/10/2008

Basis: N/A

Prep Method: SW5030C

File Name: D24934

CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
95-49-8	2-CHLOROTOLUENE	10	9.28	1		93	77 - 128%
108-67-8	1,3,5-TRIMETHYLBENZENE	10	9.79	1		98	77 - 131%
106-43-4	4-CHLOROTOLUENE	10	9.25	1		93	79 - 128%
98-06-6	TERT-BUTYLBENZENE	10	9.97	1		100	76 - 134%
95-63-6	1,2,4-TRIMETHYLBENZENE	10	9.67	1		97	80 - 138%
135-98-8	SEC-BUTYLBENZENE	10	9.71	1		97	73 - 135%
541-73-1	1,3-DICHLOROBENZENE	10	9.92	1		99	79 - 126%
99-87-6	P-ISOPROPYLtoluene	10	10	1		100	72 - 132%
106-46-7	1,4-DICHLOROBENZENE	10	9.83	1		98	81 - 125%
104-51-8	N-BUTYLBENZENE	10	9.42	1		94	77 - 141%
95-50-1	1,2-DICHLOROBENZENE	10	9.76	1		98	82 - 128%
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	10	9.27	2		93	64 - 134%
120-82-1	1,2,4-TRICHLOROBENZENE	10	10.8	1		108	80 - 128%
87-68-3	HEXACHLOROBUTADIENE	10	12.4	1		124	70 - 136%
91-20-3	NAPHTHALENE	10	9.42	1		94	78 - 125%
87-61-6	1,2,3-TRICHLOROBENZENE	10	11.1	1		111	79 - 131%

Data Package ID: VL0812076-1

Date Printed: Thursday, December 11, 2008

ALS Paragon

LIMS Version: 6.215A

Page 3 of 6

# GC/MS Volatiles

## Method SW8260\_25B

### Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: Paragon Analytics

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: VL081210-4LCSD

Sample Matrix: WATER

% Moisture: N/A

Prep Batch: VL081210-4

Sample Aliquot: 10 ml

QCBatchID: VL081210-4-1

Final Volume: 10 ml

Date Collected: N/A

Run ID: VL081210-4A

Result Units: UG/L

Date Extracted: 12/10/2008

Cleanup: NONE

Clean DF: 1

Date Analyzed: 12/10/2008

Basis: N/A

Prep Method: SW5030C

File Name: D24935

CASNO	Target Analyte	Spike Added	LCSD Result	Reporting Limit	Result Qualifier	LCSD % Rec.	RPD Limit	RPD
75-71-8	DICHLORODIFLUOROMETHANE	10	10.3	1		103	20	7
74-87-3	CHLOROMETHANE	10	9.13	1		91	20	10
75-01-4	VINYL CHLORIDE	10	9.29	1		93	20	11
74-83-9	BROMOMETHANE	10	9.76	1		98	20	5
75-00-3	CHLOROETHANE	10	10.6	1		106	20	7
75-69-4	TRICHLOROFUOROMETHANE	10	10.4	1		104	20	6
75-35-4	1,1-DICHLOROETHENE	10	9.4	1		94	20	6
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHA	10	9.26	1		93	20	5
67-64-1	ACETONE	40	41.4	10		103	30	5
74-88-4	IODOMETHANE	10	10.8	1		108	20	4
75-15-0	CARBON DISULFIDE	10	9.62	1		96	20	8
75-09-2	METHYLENE CHLORIDE	10	9.59	1		96	20	3
156-60-5	TRANS-1,2-DICHLOROETHENE	10	9.83	1		98	20	4
1634-04-4	METHYL TERTIARY BUTYL ETHER	20	19.8	1		99	20	6
75-34-3	1,1-DICHLOROETHANE	10	9.25	1		92	20	5
108-05-4	VINYL ACETATE	10	10.4	2		104	20	8
156-59-2	CIS-1,2-DICHLOROETHENE	10	9.92	1		99	20	2
78-93-3	2-BUTANONE	40	38.5	10		96	30	13
74-97-5	BROMOCHLOROMETHANE	10	10.8	1		108	20	7
67-66-3	CHLOROFORM	10	9.63	1		96	20	2
71-55-6	1,1,1-TRICHLOROETHANE	10	9.74	1		97	20	4
594-20-7	2,2-DICHLOROPROPANE	10	9.33	1		93	20	6
56-23-5	CARBON TETRACHLORIDE	10	9.66	1		97	20	5
563-58-6	1,1-DICHLOROPROPENE	10	9.44	1		94	20	5
107-06-2	1,2-DICHLOROETHANE	10	9.65	1		96	20	3
71-43-2	BENZENE	10	9.54	1		95	20	2
79-01-6	TRICHLOROETHENE	10	9.84	1		98	20	3

Data Package ID: VL0812076-1

Date Printed: Thursday, December 11, 2008

ALS Paragon

LIMS Version: 6.215A

Page 4 of 6

# GC/MS Volatiles

## Method SW8260\_25B

### Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: Paragon Analytics

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: VL081210-4LCSD

Sample Matrix: WATER

% Moisture: N/A

Prep Batch: VL081210-4

QCBatchID: VL081210-4-1

Sample Aliquot: 10 ml

Final Volume: 10 ml

Date Collected: N/A

Run ID: VL081210-4A

Result Units: UG/L

Date Extracted: 12/10/2008

Cleanup: NONE

Clean DF: 1

Date Analyzed: 12/10/2008

Basis: N/A

Prep Method: SW5030C

File Name: D24935

CASNO	Target Analyte	Spike Added	LCSD Result	Reporting Limit	Result Qualifier	LCSD % Rec.	RPD Limit	RPD
78-87-5	1,2-DICHLOROPROPANE	10	9.51	1		95	20	1
74-95-3	DIBROMOMETHANE	10	10.2	1		102	20	8
75-27-4	BROMODICHLOROMETHANE	10	9.78	1		98	20	2
10061-01-5	CIS-1,3-DICHLOROPROPENE	10	9.98	1		100	20	3
108-10-1	4-METHYL-2-PENTANONE	40	39	10		98	30	11
108-88-3	TOLUENE	10	9.72	1		97	20	2
10061-02-6	TRANS-1,3-DICHLOROPROPENE	10	9.51	1		95	20	5
79-00-5	1,1,2-TRICHLOROETHANE	10	9.6	1		96	20	6
591-78-6	2-HEXANONE	40	37.3	10		93	30	11
127-18-4	TETRACHLOROETHENE	10	11.2	1		112	20	0
142-28-9	1,3-DICHLOROPROPANE	10	9.62	1		96	20	7
124-48-1	DIBROMOCHLOROMETHANE	10	10.5	1		105	20	7
106-93-4	1,2-DIBROMOETHANE	10	9.97	1		100	20	6
544-10-5	1-CHLOROHEXANE	10	9.73	1		97	20	3
108-90-7	CHLOROBENZENE	10	9.83	1		98	20	1
630-20-6	1,1,1,2-TETRACHLOROETHANE	10	9.95	1		100	20	1
100-41-4	ETHYLBENZENE	10	9.5	1		95	20	3
136777-61-	M+P-XYLENE	20	19.8	1		99	20	2
95-47-6	O-XYLENE	10	9.87	1		99	20	3
100-42-5	STYRENE	10	10.1	1		101	20	0
75-25-2	BROMOFORM	10	11.9	1	*	119	20	6
98-82-8	ISOPROPYLBENZENE	10	9.81	1		98	20	6
96-18-4	1,2,3-TRICHLOROPROPANE	10	9.87	1		99	20	10
79-34-5	1,1,2,2-TETRACHLOROETHANE	10	8.98	1		90	20	7
108-86-1	BROMOBENZENE	10	10.2	1		102	20	2
103-65-1	N-PROPYLBENZENE	10	9	1		90	20	4
95-49-8	2-CHLOROTOLUENE	10	8.94	1		89	20	4

Data Package ID: VL0812076-1

Date Printed: Thursday, December 11, 2008

ALS Paragon

LIMS Version: 6.215A

Page 5 of 6

# GC/MS Volatiles

## Method SW8260\_25B

### Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: Paragon Analytics

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: VL081210-4LCSD

Sample Matrix: WATER

% Moisture: N/A

Prep Batch: VL081210-4

QCBatchID: VL081210-4-1

Sample Aliquot: 10 ml

Final Volume: 10 ml

Date Collected: N/A

Run ID: VL081210-4A

Result Units: UG/L

Date Extracted: 12/10/2008

Cleanup: NONE

Clean DF: 1

Date Analyzed: 12/10/2008

Basis: N/A

Prep Method: SW5030C

File Name: D24935

CASNO	Target Analyte	Spike Added	LCSD Result	Reporting Limit	Result Qualifier	LCSD % Rec.	RPD Limit	RPD
108-67-8	1,3,5-TRIMETHYLBENZENE	10	9.37	1		94	20	4
106-43-4	4-CHLOROTOLUENE	10	9	1		90	20	3
98-06-6	TERT-BUTYLBENZENE	10	9.43	1		94	20	6
95-63-6	1,2,4-TRIMETHYLBENZENE	10	9.27	1		93	20	4
135-98-8	SEC-BUTYLBENZENE	10	9.21	1		92	20	5
541-73-1	1,3-DICHLOROBENZENE	10	9.79	1		98	20	1
99-87-6	P-ISOPROPYLtolUENE	10	9.46	1		95	20	6
106-46-7	1,4-DICHLOROBENZENE	10	9.73	1		97	20	1
104-51-8	N-BUTYLBENZENE	10	8.8	1		88	20	7
95-50-1	1,2-DICHLOROBENZENE	10	9.7	1		97	20	1
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	10	9.85	2		99	20	6
120-82-1	1,2,4-TRICHLOROBENZENE	10	10.7	1		107	20	1
87-68-3	HEXACHLOROBUTADIENE	10	11.6	1		116	20	7
91-20-3	NAPHTHALENE	10	9.71	1		97	20	3
87-61-6	1,2,3-TRICHLOROBENZENE	10	11.2	1		112	20	2

### Surrogate Recovery LCS/LCSD

CASNO	Target Analyte	Spike Added	LCS % Rec.	LCS Flag	LCSD % Rec.	LCSD Flag	Control Limits
460-00-4	4-BROMOFLUOROBENZENE	25	105		107		78 - 129
1868-53-7	DIBROMOFLUOROMETHANE	25	96		97		80 - 124
2037-26-5	TOLUENE-D8	25	96		96		81 - 119

Data Package ID: VL0812076-1

Date Printed: Thursday, December 11, 2008

ALS Paragon

LIMS Version: 6.215A

Page 6 of 6



# ALS Paragon



## Inorganics Case Narrative

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### Colorado Oil & Gas Conservation Commission

Complaint 200199690

Work Order Number: 0812076

1. This report consists of 1 water sample.
2. The sample was received cool and intact by ALS Paragon on 12/09/08.
3. The sample was prepared for analysis based on Methods for the Chemical Analysis of Waters and Wastes (MCAWW), May 1994 procedures and Environmental Monitoring Systems Laboratory (EMSL) Rev 2.1 procedures.
4. The sample was analyzed following MCAWW and EMSL procedures for the following methods:

Analyte	Method	SOP #
Alkalinity	310.1	1106 Rev 7
Bicarbonate	310.1	1106 Rev 7
Carbonate	310.1	1106 Rev 7
pH	150.1	1126 Rev 16
Specific conductance	120.1	1128 Rev 9
TDS	160.1	1101 Rev 10
Bromide	300.0	1113 Rev 11
Chloride	300.0	1113 Rev 11
Fluoride	300.0	1113 Rev 11
Nitrate as N	300.0	1113 Rev 11
Nitrite as N	300.0	1113 Rev 11
Sulfate	300.0	1113 Rev 11

5. All standards and solutions were used within their recommended shelf life.
6. The sample was prepared and analyzed within the established hold time for each analysis.

All in house quality control procedures were followed, as described below.

7. General quality control procedures.



- A preparation (method) blank and laboratory control sample (LCS) were prepared and analyzed with the samples in each applicable preparation batch. There were not more than 20 samples in each preparation batch.
  - The method blank associated with each applicable batch was below the reporting limit for the requested analytes. This indicates that no contaminants were introduced to the samples during preparation and analysis.
  - The LCS was within the acceptance limits for each applicable analysis.
  - All initial and continuing calibration blanks (ICB/CCB) associated with each applicable analytical batch were below the reporting limit for the requested analytes.
  - All initial and continuing calibration verifications (ICV/CCV) associated with each applicable analytical batch were within the acceptance criteria for the requested analytes. This indicates a valid calibration and stable instrument conditions.
8. Matrix specific quality control procedures.
- Per method requirements, matrix QC was performed for each analysis. Since a sample from this order number was not the selected quality control (QC) sample, matrix specific QC results are not included in this report.
9. Electrical conductivity screening indicated that the concentration of dissolved salts was high in the sample. Therefore, it was necessary to dilute the sample prior to injection into the ion chromatograph in order to minimize the amount of salts loaded into the analytical column.
- It was necessary to further dilute the sample in order to bring the chloride and sulfate concentrations into the analytical range of the ion chromatograph (IC).
- A reduced aliquot was taken of the sample for the TDS analysis. Reporting limits were elevated accordingly.
10. Manual integrations are performed when needed to provide consistent and defensible data following the guidelines in SOP 939 Revision 3.



The data contained in the following report have been reviewed and approved by the personnel listed below. In addition, ALS Paragon certifies that the analyses reported herein are true, complete and correct within the limits of the methods employed.

Megan Johnson  
Megan Johnson  
Inorganics Primary Data Reviewer

12/18/08  
Date

R.A. Cusick  
Inorganics Final Data Reviewer

12/18/08  
Date

# ALS Paragon

## Sample Number(s) Cross-Reference Table

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**Paragon OrderNum:** 0812076

**Client Name:** Colorado Oil & Gas Conservation Commission

**Client Project Name:** Complaint 200199690

**Client Project Number:**

**Client PO Number:** OE PHA 09000000004

---

Client Sample Number	Lab Sample Number	COC Number	Matrix	Date Collected	Time Collected
Angely WW	0812076-1		WATER	08-Dec-08	11:58
Trip Blank	0812076-2		WATER	08-Dec-08	



### **Inorganic Data Reporting Qualifiers**

The following qualifiers are used by the laboratory when reporting results of inorganic analyses.

- Concentration qualifier -- If the analyte was analyzed for but not detected a “U” is entered.
- QC qualifier -- Specified entries and their meanings are as follows:

N - Spiked sample recovery not within control limits.

\* - Duplicate analysis (relative percent difference) not within control limits.

Z - Calibration spike recovery not within control limits.



## CONDITION OF SAMPLE UPON RECEIPT FORM

Paragon Analytics

Client: COGCCWorkorder No: 0812076Project Manager: AWInitials: CDTDate: 12-9-08

1. Does this project require any <b>special handling</b> in addition to standard Paragon procedures?	YES	NO		
2. Are custody <b>seals on shipping containers</b> intact?	NONE	YES	NO	
3. Are Custody seals on <b>sample containers</b> intact?	(NONE)	YES	NO	
4. Is there a <b>COC (Chain-of-Custody) present</b> or other representative documents?	(YES)	YES	NO	
5. Are the <b>COC and bottle labels complete and legible</b> ?	(YES)	YES	NO	
6. Is the <b>COC in agreement</b> with samples received? (IDs, dates, times, no. of samples, no. of containers, matrix, requested analyses, etc.)	(YES)	YES	NO	
7. Were <b>airbills / shipping documents</b> present and/or removable?	DROP OFF	YES	NO	
8. Are all aqueous <b>samples requiring preservation</b> preserved correctly? (excluding volatiles)	N/A	YES	NO	
9. Are all aqueous <b>non-preserved samples pH 4-9</b> ?	N/A	YES	NO	
10. Is there <b>sufficient sample</b> for the requested analyses?	(YES)	YES	NO	
11. Were all samples placed in the <b>proper containers</b> for the requested analyses?	(YES)	YES	NO	
12. Are all samples within <b>holding times</b> for the requested analyses?	(YES)	YES	NO	
13. Were all sample containers received <b>intact</b> ? (not broken or leaking, etc.)	(YES)	YES	NO	
14. Are all samples requiring <b>no headspace</b> (VOC, GRO, RSK/MEE, Rx CN/S, radon) headspace free? Size of bubble: <u>&lt; green pea</u> ✓ <u>&gt; green pea</u>	N/A	YES	NO	
15. Do perchlorate LCMS-MS samples <b>have headspace</b> ? (at least 1/3 of container required)	(N/A)	YES	NO	
16. Were samples checked for and free from the presence of <b>residual chlorine</b> ? (Applicable when PM has indicated samples are from a chlorinated water source; note if field preservation with sodium thiosulfate was not observed.)	(N/A)	YES	NO	
17. Were the samples <b>shipped on ice</b> ?	(YES)	YES	NO	
18. Were cooler temperatures measured at 0.1-6.0°C?	IR gun used*: #2      #4	RAD ONLY	YES	NO
Cooler #: <u>1</u>				
Temperature (°C): <u>4.8</u>				
No. of custody seals on cooler: <u>2</u>				
DOT Survey/ Acceptance Information	External µR/hr reading:	<u>13</u>		
	Background µR/hr reading:	<u>10</u>		
Were external µR/hr readings ≤ two times background and within DOT acceptance criteria? <u>YES</u> NO / NA (If no. see Form 008.)				

Additional Information: PROVIDE DETAILS BELOW FOR A NO RESPONSE TO ANY QUESTION ABOVE, EXCEPT #1 AND #16.

HEADSPACE ALL &gt; green pea 0812076-1-1 0812076-2-1

↓      ↓      ↓      ↓

-2      -3      -2      -3

-3      -4      -3      -3

-4      -5      -4      -4

-5

If applicable, was the client contacted? YES NO / NA Contact: Peter ConstantasDate/Time: 12/9/08

e-mail

Project Manager Signature / Date: Connelly 12/9/08

# BICARBONATE AS CaCO<sub>3</sub>

## Method EPA310.1

### Sample Results

**Lab Name:** ALS Paragon

**Client Name:** Colorado Oil & Gas Conservation Commission

**Client Project ID:** Complaint 200199690

**Work Order Number:** 0812076

**Final Volume:** 100 ml

**Reporting Basis:** As Received

**Matrix:** WATER

**Prep Method:** NONE

**Result Units:** MG/L

---

Client Sample ID	Lab ID	Date Collected	Date Prepared	Date Analyzed	Percent Moisture	Dilution Factor	Result	Reporting Limit	Flag	Sample Aliquot
Angely WW	0812076-1	12/08/2008	12/16/2008	12/16/2008	N/A	1	57	5		100 ml

#### Comments:

---

1. ND or U = Not Detected at or above the client requested detection limit.

**Data Package ID:** ak0812076-1

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**Date Printed:** Wednesday, December 17, 2008

**ALS Paragon**

Page 1 of 3

LIMS Version: 6.215A

# CARBONATE AS CaCO<sub>3</sub>

## Method EPA310.1

### Sample Results

**Lab Name:** ALS Paragon

**Client Name:** Colorado Oil & Gas Conservation Commission

**Client Project ID:** Complaint 200199690

**Work Order Number:** 0812076

**Final Volume:** 100 ml

**Reporting Basis:** As Received

**Matrix:** WATER

**Prep Method:** NONE

**Result Units:** MG/L

---

Client Sample ID	Lab ID	Date Collected	Date Prepared	Date Analyzed	Percent Moisture	Dilution Factor	Result	Reporting Limit	Flag	Sample Aliquot
Angely WW	0812076-1	12/08/2008	12/16/2008	12/16/2008	N/A	1	5	5	U	100 ml

#### Comments:

---

1. ND or U = Not Detected at or above the client requested detection limit.

**Data Package ID:** ak0812076-1

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**Date Printed:** Wednesday, December 17, 2008

**ALS Paragon**

LIMS Version: 6.215A

Page 2 of 3

# TOTAL ALKALINITY AS CaCO<sub>3</sub>

## Method EPA310.1

### Sample Results

**Lab Name:** ALS Paragon

**Client Name:** Colorado Oil & Gas Conservation Commission

**Client Project ID:** Complaint 200199690

**Work Order Number:** 0812076

**Final Volume:** 100 ml

**Reporting Basis:** As Received

**Matrix:** WATER

**Prep Method:** NONE

**Result Units:** MG/L

---

Client Sample ID	Lab ID	Date Collected	Date Prepared	Date Analyzed	Percent Moisture	Dilution Factor	Result	Reporting Limit	Flag	Sample Aliquot
Angely WW	0812076-1	12/08/2008	12/16/2008	12/16/2008	N/A	1	57	5		100 ml

#### Comments:

---

1. ND or U = Not Detected at or above the client requested detection limit.

**Data Package ID:** ak0812076-1

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**Date Printed:** Wednesday, December 17, 2008

**ALS Paragon**

LIMS Version: 6.215A

Page 3 of 3

# pH

## Method EPA150.1

### Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:	Angely WW
Lab ID:	0812076-1

Sample Matrix: WATER

% Moisture: N/A

Date Collected: 08-Dec-08

Date Extracted: 12-Dec-08

Date Analyzed: 20-Dec-08

Prep Method: NONE

Prep Batch: PH081212-1

QCBatchID: PH081212-1-1

Run ID: ph081212-1a

Cleanup: NONE

Basis: As Received

File Name:

Sample Aliquot: 20 ml

Final Volume: 20 ml

Result Units: pH

Clean DF: 1

CASNO	Target Analyte	Dilution Factor	Result	Reporting Limit	Result Qualifier	EPA Qualifier
10-29-7	PH	1	7.67	0.1		

Data Package ID: *ph0812076-1*

Date Printed: Wednesday, December 17, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 1

# Specific Conductance in Water

## Method EPA120.1

### Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:	Angely WW
Lab ID:	0812076-1

Sample Matrix: WATER

% Moisture: N/A

Date Collected: 08-Dec-08

Date Extracted: 12-Dec-08

Date Analyzed: 12-Dec-08

Prep Method: NONE

Prep Batch: SC081212-1

QCBatchID: SC081212-1-1

Run ID: sc081212-1a

Cleanup: NONE

Basis: As Received

File Name:

Sample Aliquot: 45 ml

Final Volume: 45 ml

Result Units: umhos/cm

Clean DF: 1

CASNO	Target Analyte	Dilution Factor	Result	Reporting Limit	Result Qualifier	EPA Qualifier
10-34-4	SPECIFIC CONDUCTIVITY	1	2040	1		

Data Package ID: sc0812076-1

Date Printed: Wednesday, December 17, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 1

# Total Dissolved Solids

## Method EPA160.1 Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:	Angely WW
Lab ID:	0812076-1

Sample Matrix: WATER

% Moisture: N/A

Date Collected: 08-Dec-08

Date Extracted: 12-Dec-08

Date Analyzed: 12-Dec-08

Prep Method: NONE

Prep Batch: TD081212-1

QCBatchID: TD081212-1-1

Run ID: td081215-1a

Cleanup: NONE

Basis: As Received

File Name: Manual Entry

Sample Aliquot: 50 ml

Final Volume: 100 ml

Result Units: MG/L

Clean DF: 1

CASNO	Target Analyte	Dilution Factor	Result	Reporting Limit	Result Qualifier	EPA Qualifier
10-33-3	TOTAL DISSOLVED SOLIDS	1	1600	40		

Data Package ID: td0812076-1

Date Printed: Wednesday, December 17, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 1

# Ion Chromatography

## Method EPA300.0 Revision 2.1 Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID:	Angely WW
Lab ID:	0812076-1

Sample Matrix: WATER

% Moisture: N/A

Date Collected: 08-Dec-08

Date Extracted: 09-Dec-08

Date Analyzed: 09-Dec-08

Prep Method: NONE

Prep Batch: IC081209-1

QCBatchID: IC081209-1-1

Run ID: ic081209-1a

Cleanup: NONE

Basis: As Received

File Name: 81209\_032.DXD

Sample Aliquot: 5 ml

Final Volume: 5 ml

Result Units: MG/L

Clean DF: 1

CASNO	Target Analyte	Dilution Factor	Result	Reporting Limit	Result Qualifier	EPA Qualifier
16984-48-8	FLUORIDE	2	3.2	0.2		
16887-00-6	CHLORIDE	20	61	4		
14797-65-0	NITRITE AS N	2	0.2	0.2	U	
24959-67-9	BROMIDE	2	0.74	0.4		
14797-55-8	NITRATE AS N	2	0.4	0.4	U	
14808-79-8	SULFATE	20	1000	20		

Data Package ID: *ic0812076-1*

Date Printed: Wednesday, December 17, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 1

# BICARBONATE AS CaCO<sub>3</sub>

## Method EPA310.1

### Method Blank

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

---

Lab ID: AK081216-1MB

Sample Matrix: WATER  
% Moisture: N/A

Prep Batch: AK081216-1  
QCBatchID: AK081216-1-1  
Run ID: ak081216-1a  
Cleanup: NONE  
Basis: N/A

Sample Aliquot: 100 ml  
Final Volume: 100 ml  
Result Units: MG/L

---

Lab ID	Date Prepared	Date Analyzed	Percent Moisture	Dilution Factor	Result	Reporting Limit	Flag
AK081216-1MB	12/16/2008	12/16/2008	N/A	1	5	5	U

---

#### Comments:

- ND or U = Not Detected at or above the client requested detection limit.

---

Data Package ID: ak0812076-1

---

Date Printed: Wednesday, December 17, 2008

**ALS Paragon**

Page 1 of 3

LIMS Version: 6.215A

# CARBONATE AS CaCO<sub>3</sub>

## Method EPA310.1

### Method Blank

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

---

Lab ID: AK081216-1MB

Sample Matrix: WATER  
% Moisture: N/A

Prep Batch: AK081216-1  
QCBatchID: AK081216-1-1  
Run ID: ak081216-1a  
Cleanup: NONE  
Basis: N/A

Sample Aliquot: 100 ml  
Final Volume: 100 ml  
Result Units: MG/L

---

Lab ID	Date Prepared	Date Analyzed	Percent Moisture	Dilution Factor	Result	Reporting Limit	Flag
AK081216-1MB	12/16/2008	12/16/2008	N/A	1	5	5	U

---

#### Comments:

- ND or U = Not Detected at or above the client requested detection limit.

---

Data Package ID: ak0812076-1

---

Date Printed: Wednesday, December 17, 2008

**ALS Paragon**

Page 2 of 3

LIMS Version: 6.215A

# TOTAL ALKALINITY AS CaCO<sub>3</sub>

## Method EPA310.1

### Method Blank

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

---

Lab ID: AK081216-1MB

Sample Matrix: WATER  
% Moisture: N/A

Prep Batch: AK081216-1  
QCBatchID: AK081216-1-1  
Run ID: ak081216-1a  
Cleanup: NONE  
Basis: N/A

Sample Aliquot: 100 ml  
Final Volume: 100 ml  
Result Units: MG/L

---

Lab ID	Date Prepared	Date Analyzed	Percent Moisture	Dilution Factor	Result	Reporting Limit	Flag
AK081216-1MB	12/16/2008	12/16/2008	N/A	1	5	5	U

---

#### Comments:

- ND or U = Not Detected at or above the client requested detection limit.

---

Data Package ID: ak0812076-1

---

Date Printed: Wednesday, December 17, 2008

**ALS Paragon**

LIMS Version: 6.215A

Page 3 of 3

# TOTAL ALKALINITY AS CaCO<sub>3</sub>

Method EPA310.1

## Laboratory Control Sample

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

Client Project ID: Complaint 200199690

Lab ID: AK081216-1LCS

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 12/16/2008

Date Analyzed: 12/16/2008

Prep Batch: AK081216-1

QCBatchID: AK081216-1-1

Run ID: ak081216-1a

Cleanup: NONE

Basis: N/A

Sample Aliquot: 100 ml

Final Volume: 100 ml

Result Units: MG/L

CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
11-43-8	TOTAL ALKALINITY AS CaCO <sub>3</sub>	100	97.7	5		98	85 - 115

Data Package ID: ak0812076-1

Date Printed: Wednesday, December 17, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 1

# Total Dissolved Solids

## Method EPA160.1

### Method Blank

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: TD081212-1MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 12-Dec-08

Date Analyzed: 12-Dec-08

Prep Method: NONE

Prep Batch: TD081212-1

QCBatchID: TD081212-1-1

Run ID: td081215-1a

Cleanup: NONE

Basis: N/A

File Name: Manual Entry

Sample Aliquot: 100 ml

Final Volume: 100 ml

Result Units: MG/L

Clean DF: 1

CASNO	Target Analyte	DF	Result	Reporting Limit	Result Qualifier	EPA Qualifier
10-33-3	TOTAL DISSOLVED SOLIDS	1	20	20	U	

Data Package ID: td0812076-1

Date Printed: Wednesday, December 17, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 1

# Total Dissolved Solids

## Method EPA160.1 Laboratory Control Sample

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: TD081212-1LCS	Sample Matrix: WATER % Moisture: N/A Date Collected: N/A Date Extracted: 12/12/2008 Date Analyzed: 12/12/2008 Prep Method: NONE	Prep Batch: TD081212-1 QCBatchID: TD081212-1-1 Run ID: td081215-1a Cleanup: NONE Basis: N/A File Name: Manual Entry	Sample Aliquot: 100 ml Final Volume: 100 ml Result Units: MG/L Clean DF: 1				
<hr/>							
CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
10-33-3	TOTAL DISSOLVED SOLIDS	400	401	20		100	85 - 115%

Data Package ID: td0812076-1

Date Printed: Wednesday, December 17, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 1

# Ion Chromatography

## Method EPA300.0 Revision 2.1

### Method Blank

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: IC081209-1MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 09-Dec-08

Date Analyzed: 09-Dec-08

Prep Method: NONE

Prep Batch: IC081209-1

QCBatchID: IC081209-1-1

Run ID: ic081209-1a

Cleanup: NONE

Basis: N/A

File Name: 81209\_011.DXD

Sample Aliquot: 5 ml

Final Volume: 5 ml

Result Units: MG/L

Clean DF: 1

CASNO	Target Analyte	DF	Result	Reporting Limit	Result Qualifier	EPA Qualifier
16984-48-8	FLUORIDE	1	0.1	0.1	U	
16887-00-6	CHLORIDE	1	0.2	0.2	U	
14797-65-0	NITRITE AS N	1	0.1	0.1	U	
24959-67-9	BROMIDE	1	0.2	0.2	U	
14797-55-8	NITRATE AS N	1	0.2	0.2	U	
14808-79-8	SULFATE	1	1	1	U	

Data Package ID: ic0812076-1

Date Printed: Wednesday, December 17, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 1

# Ion Chromatography

Method EPA300.0 Revision 2.1

## Laboratory Control Sample

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: IC081209-1LCS	Sample Matrix: WATER % Moisture: N/A Date Collected: N/A Date Extracted: 12/09/2008 Date Analyzed: 12/09/2008 Prep Method: NONE	Prep Batch: IC081209-1 QCBatchID: IC081209-1-1 Run ID: ic081209-1a Cleanup: NONE Basis: N/A File Name: 81209_012.DXD	Sample Aliquot: 5 ml Final Volume: 5 ml Result Units: MG/L Clean DF: 1				
<hr/>							
CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
16984-48-8	FLUORIDE	2.5	2.49	0.1		100	90 - 110%
16887-00-6	CHLORIDE	5	5.39	0.2		108	90 - 110%
14797-65-0	NITRITE AS N	2	1.99	0.1		99	90 - 110%
24959-67-9	BROMIDE	5	5.26	0.2		105	90 - 110%
14797-55-8	NITRATE AS N	5	5.2	0.2		104	90 - 110%
14808-79-8	SULFATE	25	26.3	1		105	90 - 110%

Data Package ID: ic0812076-1

Date Printed: Wednesday, December 17, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 1



## A N A L Y S I S   R E P O R T

Lab #: 151580 Job #: 10746  
Sample Name/Number: Angely WW  
Company: Colorado Oil & Gas Conservation  
Date Sampled: 12/08/2008  
Container: Dissolved Gas Bottle  
Field/Site Name: Complaint 200199690  
Location: Huerfano County, CO  
Formation/Depth:  
Sampling Point:  
Date Received: 12/12/2008 Date Reported: 12/31/2008

Component	Chemical mol. %	Delta 13C per mil	Delta D per mil	Delta 18O per mil
Hydrogen Sulfide -----	nd			
Helium -----	nd			
Hydrogen -----	nd			
Argon -----	0.229			
Oxygen -----	2.19			
Nitrogen -----	10.47			
Carbon Dioxide -----	0.097			
Methane -----	86.96	-49.28	-235.4	
Ethane -----	0.0532			
Ethylene -----	na			
Propane -----	nd			
Iso-butane -----	nd			
N-butane -----	nd			
Iso-pentane -----	nd			
N-pentane -----	nd			
Hexanes + -----	nd			
Water -----		-104.3	-13.87	
Dissolved Inorganic Carbon -		-22.14		

Total BTU/cu.ft. dry @ 60deg F & 14.7psia, calculated: 882

Specific gravity, calculated: 0.612

Remarks: Analysis is of gas extracted from water by headspace equilibration. Analysis has been corrected for helium added to create headspace. Helium dilution factor = 0.46

nd = not detected. na = not analyzed. Isotopic composition of carbon is relative to VPDB. Isotopic composition of hydrogen and oxygen are relative to VSMOW. Calculations for BTU and specific gravity per D3588. Chemical compositions are normalized to 100%. Mol. % is approximately equal to vol. %. ASTM Chemical analysis based on standards accurate to within 2%



# ALS Paragon



## Metals Case Narrative

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### Colorado Oil & Gas Conservation Commission

Complaint 200199690

Work Order Number: 0812076

1. This report consists of 1 water sample.
2. The sample was received cool and intact by ALS Paragon on 12/09/08.
3. The sample was to be analyzed for dissolved metals. The sample was filtered through a 0.45 micron filter and preserved with nitric acid to a pH less than two prior to analysis.
4. The sample was prepared for analysis based on Methods for the Determination of Metals in Environmental Samples – Supplement 1 procedures.

Prior to analysis by Trace ICP, an ionization buffer was added to the sample and associated QC to improve the sodium and potassium quantitation.

For analysis by Trace ICP and ICP-MS, the sample was digested following method 200.2 and SOP 806 Rev. 13.

The sample was prepared for ICP-MS analysis of arsenic and selenium by passing the digested sample and associated QC through a cation exchange column. The cation exchange column removes cations from the matrix and eliminates the CaCl<sub>2</sub> (mass 75) interferences on arsenic.

5. The sample was analyzed following Methods for the Determination of Metals in Environmental Samples – Supplement 1 procedures.

Analysis by Trace ICP followed method 200.7 and SOP 807 Rev. 11.

The relationship between intensity and concentration for each element is established using at least four standards, one of which is a blank solution.

During sample analysis concentrations are computed by the software and the results are printed in mg/L. The instrument software does not provide a printout which gives both intensity and concentration. The validity of the calibration equation is tested by analyzing the following



solutions: a blank, a low level check solution with concentrations near the reporting limit, an Initial Calibration Verification (ICV) standard from a 2<sup>nd</sup> source standard solution with concentrations near the middle of the analytical range, a Continuing Calibration Verification (CCV) standard with concentrations at two times those in the ICV, and a readback of the highest calibration standard.

These solutions provide verification that the calibration equations are functioning properly throughout the analytical range of the instrument. During sample analysis dilutions are made for analytes found at concentrations above the highest calibration standard. No results are taken from extrapolations beyond the highest standard.

Analysis by ICP-MS followed method 200.8 and SOP 827 Rev. 6.

The relationship between intensity and concentration for each element is established using at least four standards, one of which is a blank solution. A calibration equation relating instrument response to concentration is developed by the instrument software. The equation is a higher order polynomial. This type of equation is used to improve quantitation accuracy at lower concentrations where the relationship between concentration and instrument response is non-linear.

During sample analysis concentrations are computed by the software and the results are printed in ug/L. The validity of the calibration equation is tested by analyzing the following solutions: a blank, a low level check solution with concentrations near the reporting limit, an Initial Calibration Verification (ICV) standard from a 2<sup>nd</sup> source standard solution with concentrations near the middle of the analytical range, a Continuing Calibration Verification (CCV) standard with concentrations near the middle of the analytical range but different than those in the ICV, and a readback of the highest calibration standard.

These solutions provide verification that the calibration equations are functioning properly throughout the analytical range of the instrument. During sample analysis dilutions are made for analytes found at concentrations above the highest calibration standard. No results are taken from extrapolations beyond the highest standard.

6. All standards and solutions are NIST traceable and were used within their recommended shelf life.
7. The sample was prepared and analyzed within the established hold times.

All in house quality control procedures were followed, as described below.

8. General quality control procedures.
  - A filter (method) blank and laboratory control sample were filtered, preserved, and digested at the same time as the samples. There were not more than 20 samples associated with each filtered blank and laboratory control sample.
  - The filter (method) blank associated with each digestion batch was below the practical quantitation limit for each requested analyte.



- The laboratory control sample associated with each digestion batch was within the acceptance limits. This indicates complete digestion according to the method.
- All initial and continuing calibration blanks associated with each analytical batch were below the practical quantitation limits for the requested analytes, with the exception of CCB6 for strontium. The method blank and laboratory control sample were the only samples associated with this order number that were bracketed by this CCB. Strontium was not detected in the method blank, and was within control limits in the laboratory control sample.
- All initial and continuing calibration verifications associated with each analytical batch were within the acceptance criteria for the requested analytes, with the exception of CCV6 for strontium. The method blank and laboratory control sample were the only samples associated with this order number that were bracketed by this CCV. Strontium was not detected in the method blank, and was within control limits in the laboratory control sample.
- The interference check samples associated with Method 200.8 were analyzed.
- The interference check samples associated with Method 200.7 were within acceptance criteria.

9. Matrix specific quality control procedures.

Per method requirements, matrix QC was performed for each analysis. Since a sample from this order number was not the selected quality control (QC) sample, matrix specific QC results are not included in this report.

10. The sample required a dilution to bring sodium into the analytical range of the Trace ICP.

It is a standard ALS Paragon practice that samples for ICP-MS are analyzed at a dilution.

11. Sodium Adsorption Ration (SAR) was determined by calculation based on a reference from the client. Calcium, magnesium, and sodium concentrations were determined by ICP, Method 200.7.

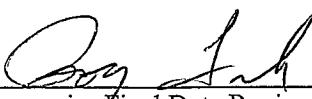
$$\text{SAR} = \text{Na}/(((\text{Ca} + \text{Mg})/2)^{1/2})$$

The analyte results are the me/L concentrations based on conversions from their mg/L concentrations.

The data contained in the following report have been reviewed and approved by the personnel listed below. In addition, ALS Paragon certifies that the analyses reported herein are true, complete and correct within the limits of the methods employed.

  
Emily Knodel  
Inorganics Primary Data Reviewer

12-22-08  
Date

  
Bob Link  
Inorganics Final Data Reviewer

12/22/08  
Date



## **Inorganic Data Reporting Qualifiers**

The following qualifiers are used by the laboratory when reporting results of inorganic analyses.

- Result qualifier -- If the analyte was analyzed for but not detected a "U" is entered.
- QC qualifier -- Specified entries and their meanings are as follows:
  - E - The reported value is estimated because of the presence of interference. An explanatory note may be included in the narrative.
  - M - Duplicate injection precision was not met.
  - N - Spiked sample recovery not within control limits. A post spike is analyzed for all ICP analyses when the matrix spike and or spike duplicate fail and the native sample concentration is less than four times the spike added concentration.
  - Z - Spiked recovery not within control limits. An explanatory note may be included in the narrative.
  - \* - Duplicate analysis (relative percent difference) not within control limits.
  - S - SAR value is estimated as one or more analytes used in the calculation were not detected above the detection limit.

# ALS Paragon

## Sample Number(s) Cross-Reference Table

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**Paragon OrderNum:** 0812076

**Client Name:** Colorado Oil & Gas Conservation Commission

**Client Project Name:** Complaint 200199690

**Client Project Number:**

**Client PO Number:** OE PHA 09000000004

---

Client Sample Number	Lab Sample Number	COC Number	Matrix	Date Collected	Time Collected
Angely WW	0812076-1		WATER	08-Dec-08	11:58
Trip Blank	0812076-2		WATER	08-Dec-08	



## CONDITION OF SAMPLE UPON RECEIPT FORM

Paragon Analytics

Client: COGCCWorkorder No: 0812076Project Manager: AWInitials: CDTDate: 12-9-08

1. Does this project require any <b>special handling</b> in addition to standard Paragon procedures?	YES	NO		
2. Are custody <b>seals on shipping containers</b> intact?	NONE	YES	NO	
3. Are Custody seals on <b>sample containers</b> intact?	(NONE)	YES	NO	
4. Is there a <b>COC (Chain-of-Custody) present</b> or other representative documents?	(YES)	YES	NO	
5. Are the <b>COC and bottle labels complete and legible</b> ?	(YES)	YES	NO	
6. Is the <b>COC in agreement</b> with samples received? (IDs, dates, times, no. of samples, no. of containers, matrix, requested analyses, etc.)	(YES)	YES	NO	
7. Were <b>airbills / shipping documents</b> present and/or removable?	DROP OFF	YES	NO	
8. Are all aqueous <b>samples requiring preservation</b> preserved correctly? (excluding volatiles)	N/A	YES	NO	
9. Are all aqueous <b>non-preserved samples pH 4-9</b> ?	N/A	YES	NO	
10. Is there <b>sufficient sample</b> for the requested analyses?	(YES)	YES	NO	
11. Were all samples placed in the <b>proper containers</b> for the requested analyses?	(YES)	YES	NO	
12. Are all samples within <b>holding times</b> for the requested analyses?	(YES)	YES	NO	
13. Were all sample containers received <b>intact</b> ? (not broken or leaking, etc.)	(YES)	YES	NO	
14. Are all samples requiring <b>no headspace</b> (VOC, GRO, RSK/MEE, Rx CN/S, radon) headspace free? Size of bubble: <u>&lt; green pea</u> ✓ <u>&gt; green pea</u>	N/A	YES	NO	
15. Do perchlorate LCMS-MS samples <b>have headspace</b> ? (at least 1/3 of container required)	(N/A)	YES	NO	
16. Were samples checked for and free from the presence of <b>residual chlorine</b> ? (Applicable when PM has indicated samples are from a chlorinated water source; note if field preservation with sodium thiosulfate was not observed.)	(N/A)	YES	NO	
17. Were the samples <b>shipped on ice</b> ?	(YES)	YES	NO	
18. Were cooler temperatures measured at 0.1-6.0°C?	IR gun used*: #2      #4	RAD ONLY	YES	NO
Cooler #: <u>1</u>				
Temperature (°C): <u>4.8</u>				
No. of custody seals on cooler: <u>2</u>				
DOT Survey/ Acceptance Information	External µR/hr reading:	<u>13</u>		
	Background µR/hr reading:	<u>10</u>		
Were external µR/hr readings ≤ two times background and within DOT acceptance criteria? <u>YES</u> NO / NA (If no. see Form 008.)				

Additional Information: PROVIDE DETAILS BELOW FOR A NO RESPONSE TO ANY QUESTION ABOVE, EXCEPT #1 AND #16.

HEADSPACE ALL &gt; green pea 0812076-1-1 0812076-2-1

↓      ↓      ↓      ↓

-2      -3      -2      -3

-3      -4      -3      -3

-4      -5      -4      -4

-5

If applicable, was the client contacted? YES NO / NA Contact: Peter ConstantasDate/Time: 12/9/08

e-mail

Project Manager Signature / Date: Connelly 12/9/08

# Dissolved Metals by 200.7

## Method EPA200.7 Revision 4.4 Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID: Angely WW	Sample Matrix: WATER	Prep Batch: IP081211-4	Sample Aliquot: 50 g
Lab ID: 0812076-1	% Moisture: N/A	QCBatchID: IP081211-4-1	Final Volume: 50 g
	Date Collected: 08-Dec-08	Run ID: IT081216-1A7	Result Units: mg/l
	Date Extracted: 11-Dec-08	Cleanup: NONE	Clean DF: 1
	Date Analyzed: 16-Dec-08	Basis: As Received	
	Prep Method: EPA200.2 Rev 2.8	File Name: T81216A.	

CASNO	Target Analyte	Dilution Factor	Result	Reporting Limit	Result Qualifier	EPA Qualifier
7440-39-3	BARIUM	1	0.1	0.1	U	
7440-41-7	BERYLLIUM	1	0.002	0.002	U	
7440-42-8	BORON	1	0.45	0.1		
7440-70-2	CALCIUM	1	190	1		
7440-47-3	CHROMIUM	1	0.01	0.01	U	
7440-48-4	COBALT	1	0.01	0.01	U	
7440-50-8	COPPER	1	0.01	0.01	U	
7439-89-6	IRON	1	0.1	0.1	U	
7439-93-2	LITHIUM	1	0.029	0.01		
7439-95-4	MAGNESIUM	1	2.5	1		
7439-96-5	MANGANESE	1	0.19	0.01		
7440-02-0	NICKEL	1	0.02	0.02	U	
7440-09-7	POTASSIUM	1	2.2	1		
7440-23-5	SODIUM	5	260	5		
	SODIUM ADSORPTION RATIO	5	5.2	0.85		
7440-24-6	STRONTIUM	1	3.8	0.01		
7440-66-6	ZINC	1	0.02	0.02	U	

Data Package ID: IT0812076-1

Date Printed: Monday, December 22, 2008

ALS Paragon

LIMS Version: 6.217A

Page 1 of 1

# Dissolved Metals by 200.8

## Method EPA200.8 Revision 5.4 Sample Results

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Field ID: Angely WW	Sample Matrix: WATER	Prep Batch: IP081211-4	Sample Aliquot: 50 g
Lab ID: 0812076-1	% Moisture: N/A	QCBatchID: IP081211-4-2	Final Volume: 50 g
	Date Collected: 08-Dec-08	Run ID: IM081212-1A5	Result Units: UG/L
	Date Extracted: 11-Dec-08	Cleanup: NONE	Clean DF: 1
	Date Analyzed: 12-Dec-08	Basis: As Received	
	Prep Method: EPA200.2 Rev 2.8	File Name: 12DEC08A	

CASNO	Target Analyte	Dilution Factor	Result	Reporting Limit	Result Qualifier	EPA Qualifier
7440-36-0	ANTIMONY	10	0.45	0.3		
7440-38-2	ARSENIC	10	2	2	U	
7440-43-9	CADMIUM	10	0.3	0.3	U	
7439-92-1	LEAD	10	0.5	0.5	U	
7439-98-7	MOLYBDENUM	10	2	1		
7782-49-2	SELENIUM	10	1	1	U	
7440-22-4	SILVER	10	0.1	0.1	U	
7440-28-0	THALLIUM	10	0.2	0.2	U	
7440-61-1	URANIUM	10	0.1	0.1	U	

Data Package ID: IM0812076-1

Date Printed: Monday, December 22, 2008

ALS Paragon

LIMS Version: 6.217A

Page 1 of 1

# Metals by 200.7

## Method EPA200.7 Revision 4.4 Method Blank

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: F081210-1MB

Sample Matrix: WATER  
% Moisture: N/A  
Date Collected: N/A  
Date Extracted: 11-Dec-08  
Date Analyzed: 16-Dec-08  
Prep Method: EPA200.2 Rev 2.8

Prep Batch: IP081211-4  
QCBatchID: IP081211-4-1  
Run ID: IT081216-1A7  
Cleanup: NONE  
Basis: N/A  
File Name: T81216A.

Sample Aliquot: 50 g  
Final Volume: 50 g  
Result Units: mg/l  
Clean DF: 1

CASNO	Target Analyte	DF	Result	Reporting Limit	Result Qualifier	EPA Qualifier
7440-39-3	BARIUM	1	0.1	0.1	U	
7440-41-7	BERYLLIUM	1	0.002	0.002	U	
7440-42-8	BORON	1	0.1	0.1	U	
7440-70-2	CALCIUM	1	1	1	U	
7440-47-3	CHROMIUM	1	0.01	0.01	U	
7440-48-4	COBALT	1	0.01	0.01	U	
7440-50-8	COPPER	1	0.01	0.01	U	
7439-89-6	IRON	1	0.1	0.1	U	
7439-93-2	LITHIUM	1	0.01	0.01	U	
7439-95-4	MAGNESIUM	1	1	1	U	
7439-96-5	MANGANESE	1	0.01	0.01	U	
7440-02-0	NICKEL	1	0.02	0.02	U	
7440-09-7	POTASSIUM	1	1	1	U	
7440-23-5	SODIUM	1	1	1	U	
7440-24-6	STRONTIUM	1	0.01	0.01	U	
7440-66-6	ZINC	1	0.02	0.02	U	

Data Package ID: IT0812076-1

Date Printed: Monday, December 22, 2008

ALS Paragon

LIMS Version: 6.217A

Page 1 of 1

# Metals by 200.7

## Method EPA200.7 Revision 4.4 Laboratory Control Sample

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: F081210-1LCS	Sample Matrix: WATER % Moisture: N/A Date Collected: N/A Date Extracted: 12/11/2008 Date Analyzed: 12/16/2008 Prep Method: EPA200.22.8	Prep Batch: IP081211-4 QCBatchID: IP081211-4-1 Run ID: IT081216-1A7 Cleanup: NONE Basis: N/A File Name: T81216A.	Sample Aliquot: 50 g Final Volume: 50 g Result Units: mg/l Clean DF: 1
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CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
7440-39-3	BARIUM	2	2.02	0.1		101	85 - 115%
7440-41-7	BERYLLIUM	0.05	0.0449	0.002		90	85 - 115%
7440-42-8	BORON	1	1.01	0.1		101	85 - 115%
7440-70-2	CALCIUM	40	39.9	1		100	85 - 115%
7440-47-3	CHROMIUM	0.2	0.199	0.01		99	85 - 115%
7440-48-4	COBALT	0.5	0.479	0.01		96	85 - 115%
7440-50-8	COPPER	0.25	0.237	0.01		95	85 - 115%
7439-89-6	IRON	1	1.01	0.1		101	85 - 115%
7439-93-2	LITHIUM	0.5	0.518	0.01		104	85 - 115%
7439-95-4	MAGNESIUM	40	40.5	1		101	85 - 115%
7439-96-5	MANGANESE	0.5	0.49	0.01		98	85 - 115%
7440-02-0	NICKEL	0.5	0.5	0.02		100	85 - 115%
7440-09-7	POTASSIUM	40	43.2	1		108	85 - 115%
7440-23-5	SODIUM	40	40	1		100	85 - 115%
7440-24-6	STRONTIUM	0.5	0.484	0.01		97	85 - 115%
7440-66-6	ZINC	0.5	0.509	0.02		102	85 - 115%

Data Package ID: IT0812076-1

Date Printed: Monday, December 22, 2008

ALS Paragon

LIMS Version: 6.217A

Page 1 of 1

# Metals by 200.8

## Method EPA200.8 Revision 5.4 Method Blank

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: F081210-1MB

Sample Matrix: WATER  
% Moisture: N/A  
Date Collected: N/A  
Date Extracted: 11-Dec-08  
Date Analyzed: 12-Dec-08  
Prep Method: EPA200.2 Rev 2.8

Prep Batch: IP081211-4  
QCBatchID: IP081211-4-2  
Run ID: IM081212-1A5  
Cleanup: NONE  
Basis: N/A  
File Name: 12DEC08A

Sample Aliquot: 50 g  
Final Volume: 50 g  
Result Units: UG/L  
Clean DF: 1

CASNO	Target Analyte	DF	Result	Reporting Limit	Result Qualifier	EPA Qualifier
7440-36-0	ANTIMONY	10	0.3	0.3	U	
7440-38-2	ARSENIC	10	2	2	U	
7440-43-9	CADMIUM	10	0.3	0.3	U	
7439-92-1	LEAD	10	0.5	0.5	U	
7439-98-7	MOLYBDENUM	10	1	1	U	
7782-49-2	SELENIUM	10	1	1	U	
7440-22-4	SILVER	10	0.1	0.1	U	
7440-28-0	THALLIUM	10	0.2	0.2	U	
7440-61-1	URANIUM	10	0.1	0.1	U	

Data Package ID: IM0812076-1

Date Printed: Monday, December 22, 2008

ALS Paragon

LIMS Version: 6.217A

Page 1 of 1

# Metals by 200.8

## Method EPA200.8 Revision 5.4 Laboratory Control Sample

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: FM81210-1LCS

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 12/11/2008

Date Analyzed: 12/12/2008

Prep Method: EPA200.22.8

Prep Batch: IP081211-4

QCBatchID: IP081211-4-2

Run ID: IM081212-1A5

Cleanup: NONE

Basis: N/A

File Name: 12DEC08A

Sample Aliquot: 50 g

Final Volume: 50 g

Result Units: UG/L

Clean DF: 1

CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
7440-36-0	ANTIMONY	20	19.7	0.3		99	85 - 115%
7440-38-2	ARSENIC	40	41.9	2		105	85 - 115%
7440-43-9	CADMIUM	20	19.6	0.3		98	85 - 115%
7439-92-1	LEAD	100	100	0.5		100	85 - 115%
7439-98-7	MOLYBDENUM	20	19.7	1		99	85 - 115%
7782-49-2	SELENIUM	40	40.9	1		102	85 - 115%
7440-22-4	SILVER	20	20.5	0.1		103	85 - 115%
7440-28-0	THALLIUM	1	0.848	0.2		85	85 - 115%
7440-61-1	URANIUM	20	20.6	0.1		103	85 - 115%

Data Package ID: IM0812076-1

Date Printed: Monday, December 22, 2008

ALS Paragon

LIMS Version: 6.217A

Page 1 of 1



# ALS Paragon



## Total Organic Carbon Case Narrative

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### Colorado Oil & Gas Conservation Commission

Complaint 200199690

**Order Number - 0812076**

1. This report consists of 1 water sample.
2. The sample was received cool and intact by ALS Paragon on 12/09/08.
3. The sample had been correctly preserved for the requested analysis.
4. The sample was prepared for analysis based on Methods for the Chemical Analysis of Waters and Wastes (MCAWW), May 1994 procedures.
5. The sample was analyzed following MCAWW procedures for the following method:

<u>Analyte</u>	<u>Method</u>	<u>SOP #</u>
TOC (Total Organic Carbon)	415.1	670 Rev 12
6. All standards and solutions were used within their recommended shelf life.
7. The sample was prepared and analyzed within the established hold time for TOC analysis.

All in house quality control procedures were followed, as described below.

8. General quality control procedures.
  - n A preparation (method) blank, laboratory control sample (LCS) and laboratory control sample duplicate (LCSD) were prepared and analyzed with the samples in this preparation batch. There were not more than 20 samples in this preparation batch.
  - n The method blank associated with this batch was below the reporting limit for the requested analyte. This indicates that no contaminants were introduced to the samples during preparation and analysis.
  - n The LCS and LCSD were within the acceptance limits for TOC analysis.



- All continuing calibration verifications (CCV) associated with this batch were within the acceptance criteria for the requested analyte. This indicates a valid calibration and stable instrument conditions.

9. Matrix specific quality control procedures.

Since a sample from this Order Number was not the selected quality control (QC) sample, matrix specific QC results are not included in this report.

The data contained in the following report have been reviewed and approved by the personnel listed below. In addition, ALS Paragon certifies that the analyses reported herein are true, complete and correct within the limits of the methods employed.

SLJ

Sharon L. Jobes  
Organics Primary Data Reviewer

12-16-08

Date

B. prastit

Organics Final Data Reviewer

12-16-08

Date

# ALS Paragon

## Sample Number(s) Cross-Reference Table

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**Paragon OrderNum:** 0812076

**Client Name:** Colorado Oil & Gas Conservation Commission

**Client Project Name:** Complaint 200199690

**Client Project Number:**

**Client PO Number:** OE PHA 09000000004

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Client Sample Number	Lab Sample Number	COC Number	Matrix	Date Collected	Time Collected
Angely WW	0812076-1		WATER	08-Dec-08	11:58
Trip Blank	0812076-2		WATER	08-Dec-08	



## CONDITION OF SAMPLE UPON RECEIPT FORM

Paragon Analytics

Client: COGCCWorkorder No: 0812076Project Manager: AWInitials: CDTDate: 12-9-08

1. Does this project require any <b>special handling</b> in addition to standard Paragon procedures?	YES	NO	
2. Are custody <b>seals on shipping containers</b> intact?	NONE	YES	NO
3. Are Custody seals on <b>sample containers</b> intact?	(NONE)	YES	NO
4. Is there a <b>COC (Chain-of-Custody) present</b> or other representative documents?	(YES)	YES	NO
5. Are the <b>COC and bottle labels complete and legible</b> ?	(YES)	YES	NO
6. Is the <b>COC in agreement</b> with samples received? (IDs, dates, times, no. of samples, no. of containers, matrix, requested analyses, etc.)	(YES)	YES	NO
7. Were <b>airbills / shipping documents</b> present and/or removable?	DROP OFF	YES	NO
8. Are all aqueous <b>samples requiring preservation</b> preserved correctly? (excluding volatiles)	N/A	YES	NO
9. Are all aqueous <b>non-preserved samples pH 4-9</b> ?	N/A	YES	NO
10. Is there <b>sufficient sample</b> for the requested analyses?	(YES)	YES	NO
11. Were all samples placed in the <b>proper containers</b> for the requested analyses?	(YES)	YES	NO
12. Are all samples within <b>holding times</b> for the requested analyses?	(YES)	YES	NO
13. Were all sample containers received <b>intact</b> ? (not broken or leaking, etc.)	(YES)	YES	NO
14. Are all samples requiring <b>no headspace</b> (VOC, GRO, RSK/MEE, Rx CN/S, radon) headspace free? Size of bubble: <u>&lt; green pea</u> ✓ <u>&gt; green pea</u>	N/A	YES	NO
15. Do perchlorate LCMS-MS samples <b>have headspace</b> ? (at least 1/3 of container required)	(N/A)	YES	NO
16. Were samples checked for and free from the presence of <b>residual chlorine</b> ? (Applicable when PM has indicated samples are from a chlorinated water source; note if field preservation with sodium thiosulfate was not observed.)	(N/A)	YES	NO
17. Were the samples <b>shipped on ice</b> ?	(YES)	YES	NO
18. Were cooler temperatures measured at 0.1-6.0°C?	IR gun used*: #2      #4 RAD ONLY	YES	NO
Cooler #:	<u>1</u>		
Temperature (°C):	<u>4.8</u>		
No. of custody seals on cooler:	<u>2</u>		
External µR/hr reading:	<u>13</u>		
Background µR/hr reading:	<u>10</u>		
Were external µR/hr readings ≤ two times background and within DOT acceptance criteria? <u>YES</u> NO / NA (If no. see Form 008.)			

Additional Information: PROVIDE DETAILS BELOW FOR A NO RESPONSE TO ANY QUESTION ABOVE, EXCEPT #1 AND #16.

HEADSPACE ALL > green pea    0812076-1-1    0812076-2-1

↓      -2  
 ↓      -3  
 ↓      -4  
 ↓      -5

If applicable, was the client contacted? YES NO / NA Contact: Peter ConstantasDate/Time: 12/9/08

e-mail

Project Manager Signature / Date: Connelly 12/9/08



### **Inorganic Data Reporting Qualifiers**

The following qualifiers are used by the laboratory when reporting results of inorganic analyses.

- Concentration qualifier -- If the analyte was analyzed for but not detected a “U” is entered.
- QC qualifier -- Specified entries and their meanings are as follows:
  - N - Spiked sample recovery not within control limits.
  - \* - Duplicate analysis (relative percent difference) not within control limits.
  - B - The method blank for the analysis contained the analyte of interest above the reporting limit.

# TOTAL ORGANIC CARBON

Method EPA415.1

## Sample Results

Lab Name: ALS Paragon

Client Name: Colorado Oil & Gas Conservation Commission

Client Project ID: Complaint 200199690

Work Order Number: 0812076

Final Volume: 40 ml

Reporting Basis: As Received

Matrix: WATER

Prep Method: NONE

Result Units: MG/L

---

Client Sample ID	Lab ID	Date Collected	Date Prepared	Date Analyzed	Percent Moisture	Dilution Factor	Result	Reporting Limit	Flag	Sample Aliquot
Angely WW	0812076-1	12/08/2008	12/12/2008	12/12/2008	N/A	1	1.6	1		40 ml

### Comments:

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1. ND or U = Not Detected at or above the client requested detection limit.

Data Package ID: MO0812076-1

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Date Printed: Monday, December 15, 2008

ALS Paragon

Page 1 of 1

LIMS Version: 6.215A

# TOTAL ORGANIC CARBON

## Method EPA415.1

### Method Blank

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

---

Lab ID: MO081212-1MB

Sample Matrix: WATER  
% Moisture: N/A

Prep Batch: MO081212-1  
QCBatchID: MO081212-1-1  
Run ID: MO081212-1A  
Cleanup: NONE  
Basis: N/A

Sample Aliquot: 40 ml  
Final Volume: 40 ml  
Result Units: MG/L

---

Lab ID	Date Prepared	Date Analyzed	Percent Moisture	Dilution Factor	Result	Reporting Limit	Flag
MO081212-1MB	12/12/2008	12/12/2008	N/A	1	1	1	U

#### Comments:

---

- ND or U = Not Detected at or above the client requested detection limit.

Data Package ID: MO0812076-1

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Date Printed: Monday, December 15, 2008

**ALS Paragon**

Page 1 of 1

LIMS Version: 6.215A

# Organic Carbon

## Method EPA415.1

### Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: ALS Paragon

Work Order Number: 0812076

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200199690

Lab ID: MO081212-1LCS	Sample Matrix: WATER % Moisture: N/A Date Collected: N/A Date Extracted: 12/12/2008 Date Analyzed: 12/12/2008 Prep Method: NONE	Prep Batch: MO081212-1 QCBatchID: MO081212-1-1 Run ID: MO081212-1A Cleanup: NONE Basis: N/A File Name: 12121027	Sample Aliquot: 40 ml Final Volume: 40 ml Result Units: MG/L Clean DF: 1
-----------------------	--	--	---

CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
10-35-5	TOTAL ORGANIC CARBON	15	15.5	1		103	85 - 115%

Lab ID: MO081212-1LCSD	Sample Matrix: WATER % Moisture: N/A Date Collected: N/A Date Extracted: 12/12/2008 Date Analyzed: 12/12/2008 Prep Method: NONE	Prep Batch: MO081212-1 QCBatchID: MO081212-1-1 Run ID: MO081212-1A Cleanup: NONE Basis: N/A File Name: 12121027	Sample Aliquot: 40 ml Final Volume: 40 ml Result Units: MG/L Clean DF: 1
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CASNO	Target Analyte	Spike Added	LCSD Result	Reporting Limit	Result Qualifier	LCSD % Rec.	RPD Limit	RPD
10-35-5	TOTAL ORGANIC CARBON	15	15.4	1		103	20	1

Data Package ID: MO0812076-1

Date Printed: Monday, December 15, 2008

ALS Paragon

LIMS Version: 6.215A

Page 1 of 1