

State of Colorado  
**Oil and Gas Conservation Commission**

1120 Lincoln Street, Suite 801, Denver, Colorado 80203 (303)894-2100 Fax:(303)894-2109



FOR OGCC USE ONLY

**SITE INVESTIGATION AND REMEDIATION WORKPLAN**

This form shall be submitted to the Director for approval prior to the initiation of site investigation and remediation activities. Form 27 is intended to be used whenever possible. Additional documentation will be required when large volumes of soil and groundwater have been impacted or involve large facilities with multiple source areas. See Rule 910. Attach as many pages as needed to fully describe the proposed work.

OGCC Employee:

☐ Spill ☐ Complaint  
☐ Inspection ☐ NOAV

Tracking No:

**CAUSE OF CONDITION BEING INVESTIGATED AND REMEDIATED**

☐ Spill or Release ☐ Plug & Abandon ☐ Central Facility Closure ☐ Site/Facility Closure ☒ Other (describe): Pit Closure

OGCC Operator Number: 96850

Name of Operator: Williams Production RMT Company

Address: 1058 County Road 215

City: Parachute State: CO Zip: 81635

Contact Name and Telephone:

Karolina Blaney

No: 970-683-2295

Fax: 970-285-9573

API Number: N/A

County: Garfield

Facility Name: Chevron TR 21-20-597

Facility Number: 284697

Well Name: Chevron TR 21-20-597

Well Number: N/A

Location: (QtrQtr, Sec, Twp, Rng, Meridian): SENW, SEC 20, T5S, R97W, 6th PM Latitude: 39.601501 Longitude: -108.302601

**TECHNICAL CONDITIONS**

Type of Waste Causing Impact (crude oil, condensate, produced water, etc): Produced Water

**Site Conditions:** Is location within a sensitive area (according to Rule 901e)? ☒ Y ☐ N If yes, attach evaluation.

Adjacent land use (cultivated, irrigated, dry land farming, industrial, residential, etc.): Rangeland, Non Crop Land

Soil type, if not previously identified on Form 2A or Federal Surface Use Plan: Parachute Irigul-Complex, 5-30% slopes

Potential receptors (water wells within 1/4 mi, surface waters, etc.): Wiess Creek lies approximately 1769 feet to the west.

**Description of Impact** (if previously provided, refer to that form or document):

Impacted Media (check):

- ☒ Soils  
☐ Vegetation  
☐ Groundwater  
☐ Surface Water

Extent of Impact:

Please See Attached Notice of Completion Report  
for Remediation # 5949

How Determined:

Visual observation, field screening equipment, and analytical analysis

**REMEDIATION WORKPLAN**

**Describe initial action taken** (if previously provided, refer to that form or document):

Please See Attached Notice of Completion Report for Remediation # 5949

**Describe how source is to be removed:**

Please See Attached Notice of Completion Report for Remediation # 5949

**Describe how remediation of existing impacts is to be accomplished, including removal and disposal at an injection well or licensed facility, land treatment on site, removal of impacted groundwater, insitu bioremediation, burning of oily vegetation, etc.:**

Please See Attached Notice of Completion Report for Remediation # 5949



Tracking Number: REM # 5949  
Name of Operator: WILLIAMS  
OGCC Operator No: \_\_\_\_\_  
Received Date: \_\_\_\_\_  
Well Name & No: \_\_\_\_\_  
Facility Name & No: CHEVRON TR 21-20-597

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**REMEDIATION WORKPLAN (Cont.)**

OGCC Employee: \_\_\_\_\_

**If groundwater has been impacted, describe proposed monitoring plan** (# of wells or sample points, sampling schedule, analytical methods, etc.):

Please See Attached Notice of Completion Report for Remediation # 5949

**Describe reclamation plan.** Discuss existing and new grade recontouring; method and testing of compaction alleviation; and reseeding program, including location of new seed, seed mix and noxious weed prevention. Attach diagram or drawing. Use additional sheet for description if required.

Please See Attached Notice of Completion Report for Remediation # 5949

**Attach samples and analytical results taken to verify remediation of impacts. Show locations of samples on an onsite schematic or drawing.**

**Is further site investigation required?** ☒ Y ☐ N If yes, describe:

Please See Attached Notice of Completion Report for Remediation # 5949

**Final disposition of E&P waste** (landtreated and disposed onsite, name of licensed disposal facility, recycling, reuse, etc.):

Please See Attached Notice of Completion Report for Remediation # 5949

**IMPLEMENTATION SCHEDULE**

Date Site Investigation Began: <u>August 14, 2011</u>	Date Site Investigation Completed: <u>August 14, 2011</u>	Date Remediation Plan Submitted: <u>July 12, 2011</u>
Remediation Start Date: <u>August 16, 2011</u>	Anticipated Completion Date: <u>November 4, 2011</u>	Actual Completion Date: <u>November 4, 2011</u>

I hereby certify that the statements made in this form are, to the best of my knowledge, true, correct, and complete.

Print Name: Karolina Blaney

Signed: \_\_\_\_\_

Title: Environmental Specialist

Date: 12/15/2011

OGCC Approved: \_\_\_\_\_

Title: \_\_\_\_\_

Date: 01/26/2012

Please notify OGCC when  
pit backfill and reclamation  
are completed, to issue NFA letter and close the project.

FOR Chris Canfield  
EPS NW Region

al

## Sensitive Area Determination Checklist

<b>Williams Production RMT Company – Highlands</b>		
<b>Person(s) Conducting Field Inspection</b>	Ashlee Lane	9/9/10
	<i>Biologist</i>	
<b>Site Information</b>		
Location:	TR 22-20-597	Time: 1230
Type of Facility:	Existing Well Pad	
<b>Environmental Conditions</b>	Clear and extremely windy; soil conditions are dry.	
Temperature (°F)	70°	

Has the proposed, new or existing location been designated as a sensitive area?

☒ Yes      ☐ No

### **SURFACE WATER**

1. Are there any surface water features or SWSAs adjacent to or within ¼ mile of the proposed/new or existing facility?

☐ Yes      ☒ No

If yes, list type of surface water feature(s), i.e. rivers, creeks, streams, seeps, springs, wetlands: There were three springs identified that are outside of the ¼ mile buffer zone that will be elaborated on in the ground water comments section.

If yes, describe location relative to facility:

2. Could a potential release from the facility reach surface water features?

☐ Yes      ☒ No

If yes, describe the pathway a release from the facility would likely follow to determine if the potential to impact surface water is high or low.

3. Is the potential to impact surface water from a facility release high or low?

☐ High      ☒ Low

## GROUNDWATER

1. Will the proposed/new or existing facility have any pits which will contain hydrocarbons and chlorides or other E&P wastes?

☒ Yes      ☐ No

If yes, List the pit type(s): Multi-well pit.

2. Is the site of the proposed facility underlain by an unconfined aquifer or recharge zone?

☒ Yes      ☐ No

3. Is the hydraulic conductivity of the underlying soil or geologic material  $\leq 1.0 \times 10^{-7}$  cm/sec?

☐ Yes      ☒ No

4. Is the proposed facility located within 1/8 mile of a domestic water well or 1/4 mile of a public water supply well which would use the same aquifer?

☐ Yes      ☒ No

5. Is the proposed facility located within a 100 year floodplain?

☐ Yes (*Sensitive Area*)      ☒ No (*If no, proceed to question #6.*)

6. Is the depth to groundwater known?

☐ Yes (*If yes, follow instructions provided in 6(a) of this section.*)

☒ No (*If no, follow instructions provided in 6(b) of this section.*)

- (a) If yes, could a potential release from the proposed facility reach groundwater?

☐ Yes      ☐ No

If yes, explain:

- (b) If no:

(i) Evaluate surrounding soils, topography, and vegetation which may suggest the presence of shallow groundwater.

(ii) Gather information from surrounding well data in order to determine a depth to groundwater, i.e. State Engineers Office.

7. Is the potential to impact ground water from the facility in the event of a release high or low?

☐ High      ☒ Low

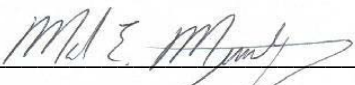
### **Additional Comments:**

As stated in the surface water section of this sensitive area determination, there are no surface water features identified within the ¼ mile buffer zone of the existing facility. There are several large valleys located to the northwest, southwest and southeast of the facility; however no defined channels have been identified on the USGS topographic maps in addition to the site investigation confirming this finding. The facility as it is currently constructed, limits the flow directions of a potential release to the western, southern, and eastern edges of the facility. If a release were to migrate off the facility it would run down the hillsides towards the above mentioned valleys. However the potential for fluids to reach any of the identified intermittent channels in the valleys would be very low due to the thick vegetative cover consisting of service berry, oak brush, sage brush, grasses, the moderate to high infiltration rates of the underlying soils, and the distance a potential release would have to travel in order to reach any identified intermittent stream channel. There are currently adequate Best Management Practices (BMPs) installed in the form of a perimeter berm on the western, southern, and eastern edges of the facility. These BMPs should be monitored and maintained to further ensure site containment in the event of a release. With the current BMPs in place, the topographical setting of the facility, the thick vegetation surrounding the facility, the moderate to high infiltration rates of the underlying soil, and the distance to any identified intermittent channels the potential to impact surface water features outside the quarter mile buffer zone would be deemed very low.

The State Engineer's Office and USGS records were reviewed and no records were revealed that would provide additional information pertaining to the depth to groundwater. The vegetative cover in the immediate vicinity of the facility, service berry, oak brush, and, sage brush does not suggest the presence of shallow groundwater. However, there were two springs identified on the USGS topographic map southwest of the existing facility approximately 2,562 feet (SWNW Sec 20) and southeast of the facility approximately 1,746 feet (NWSE Sec 20). In addition one additional spring was identified during the site investigation approximately 2,070 feet northwest of the facility (NWNW Sec 20). The facility resides in the Uintah formation, which like the Green River Formation, tends to be fractured both vertically and horizontally allowing fluids to migrate in the subsurface over large distances. Based on the topographical setting of the facility, it is not anticipated that an overland release would impact groundwater and thus potentially the spring to the northwest due to the short duration time involved and the fact it would spread out over a large area. The greatest potential for impact to groundwater would be from a release that occurred over a longer period of time such as a leaking pit, due to the close proximity of the subject pit to the springs and the likelihood of fractured bedrock. Note however that due to the topographical setting of the existing facility, the greatest potential for any impacts to the above mentioned springs would be to the springs to the northwest and southeast of the facility. Therefore it would be highly recommended that the pit be lined in accordance to COGCC criteria and tested prior to placement of any materials into it.



Based on the information collected during the site investigation and desktop review, the potential to impact surface water has been deemed very low. The greatest potential for impacts from the facility would be to groundwater due to the geologic conditions in the area and the relatively close proximity of the springs to the northwest and southeast of the existing facility. With this potential to impact groundwater, the facility should be designated as being in a sensitive area

Inspector Signature(s):  Date: 9/15/2010  
Mark E. Mumby, *Project Manager/RPG*  
HRL Compliance Solutions, Inc.

 Date: 9/9/2010  
Ashlee Lane, *Biologist*  
HRL Compliance Solutions, Inc.

***WILLIAMS PRODUCTION RMT COMPANY  
TRAIL RIDGE FIELD  
CHEVRON TR 21-20-597  
NOTICE OF COMPLETION REPORT FOR  
REMEDATION # 5949***

***August 2011***

Prepared For:



1058 County Road 215  
P.O. Box 370  
Parachute, Colorado 81635

Prepared By:



744 Horizon Court, Suite 140  
Grand Junction, CO 81506  
Phone: 970-243-3271  
Fax: 970-243-3280

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Appendix 4: Treatment Cell Confirmation Raw Analytical Data

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## Form 27 Attachment

### **Introduction**

The purpose of this Notice of Completion report – for the closure of the Williams TR 21-20-597 production pit (COGCC Facility ID # 284697; hereinafter also referred to as TR 22-20-597) – is to provide detailed information and findings analysis for the previously submitted and approved (remediation number 5949) Colorado Oil and Gas Conservation Commission (COGCC) Site Investigation and Remediation Workplan, Form 27. This report will provide the documentation necessary to demonstrate a comprehensive and diligent investigation of the pit and adjacent environment which was obtained as described and in accordance with all appropriate county, state and federal rules and regulations.

The subject Form 27 was delivered via electronic email on July 12, 2011. Preliminary approval to proceed with closure of the subject pit was issued by the COGCC and obtained by Williams Production RMT Company (Williams) on July 26, 2011; at which time the aforementioned remediation number was issued. Closure activities began in August 15, 2011 and were concluded on November 8, 2011. Information in this report includes, but is not limited to: field screening results; laboratory analytical; subliner soil remediation; liner recycling; and bioremediation of the excavated impacted soils.

### **Evacuation of Pit Contents**

Remaining pit contents were removed from the pit using hydro-vac trucks and placed in a lined bermed containment to have free liquids removed via filter press.

The filter press sludge was placed into the aforementioned lined bermed containment cell, profiled for disposal/characterization purposes, and transported to ECDC Environmental for disposal on September 5, 2011.

### **Background Sampling**

Three samples were collected from the up-gradient undisturbed hillsides surrounding the pad. All background samples were analyzed for arsenic as well as additional analysis at one location which included inorganic parameters of COGCC Table 910-1(i.e. SAR, EC, pH). Refer to Table 4 and Appendix 3 for background sampling results.

### **Pit Liner Investigation and Integrity Assessment**

The pit liner system – containing two of layers of poly synthetic material/liner and one layer of felt was present within the pit. A pit liner investigation was unable to be conducted prior to removal due to scheduling and miscommunication of crews.

## **Pit Liner Removal**

Removal of the pit liners consisted of a crew cutting the liner along the crest of the pit at an elevation adjacent to the surface of the well pad. A trackhoe bucket was utilized to grab sections of the liner for extraction and place them in a lined earthen bermed containment cell for subsequent management. Sections of liner that contained residual or trace amounts of sludge were pulled, placed into the containment cell, and allowed to dry. Liners were stored in a lined bermed containment until being banded to pallets to be recycled. During the liner removal, the bottom section of the liner was accessible for inspection and revealed no signs of tears or holes.

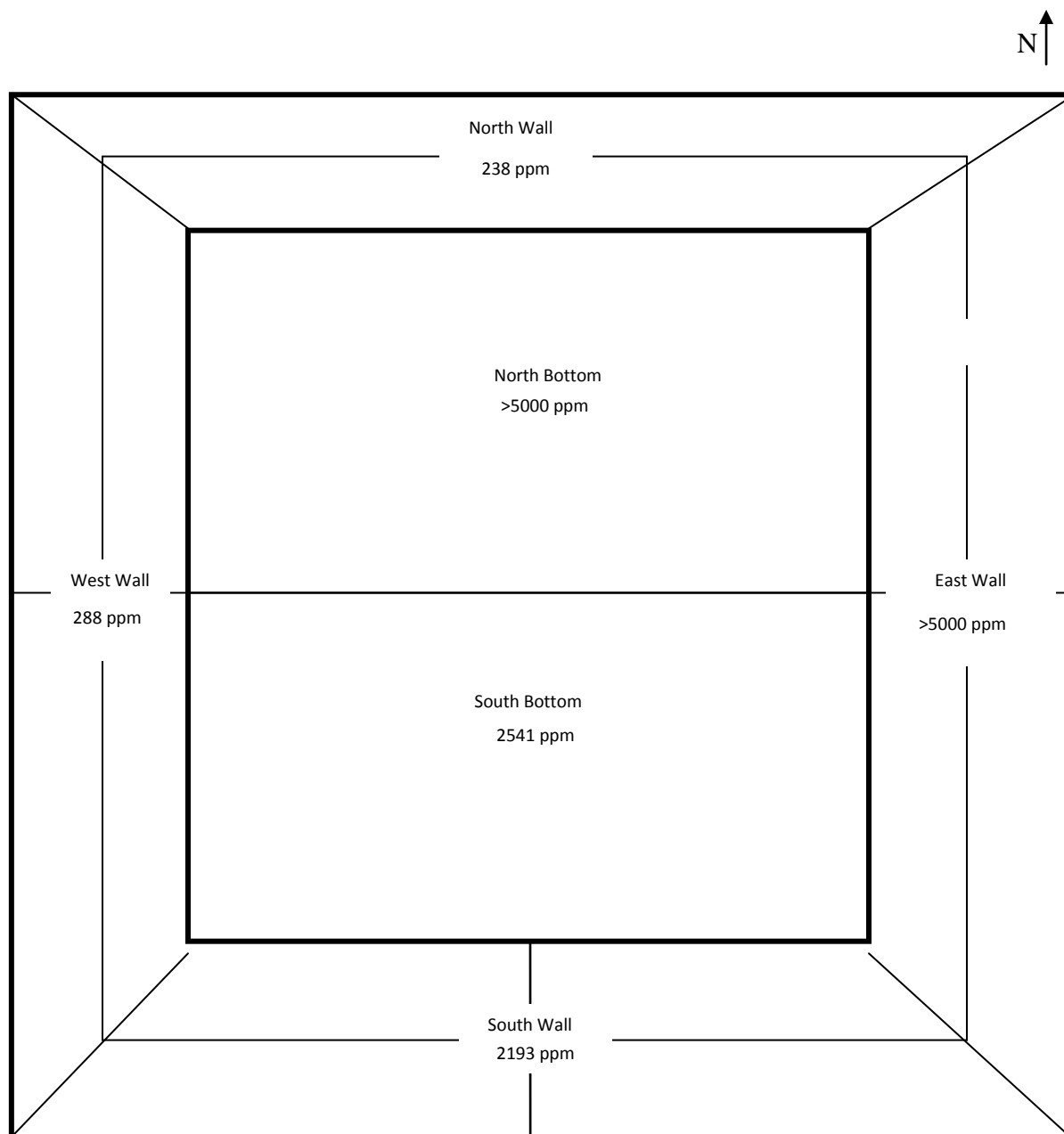
## **Subliner Soil Investigation and Activities**

Subliner soils, examined below the pit lining, were inspected visually and through the use of specialized field screening equipment (identified below) to identify areas which may exceed standards set forth in Table 910-1 of the COGCC 900-Series Rule for hydrocarbons within the soil. Soils below the second lining system on the pit bottom, southern and eastern walls were stained black and contained a moderate hydrocarbon odor. The western wall did not contain any staining nor did the soil exhibit any hydrocarbon odor.

Field screening of the pit footprint and walls was performed along the entire pit in a grid pattern of sections. The pit bottom was separated into two sections, west and east, and a five point composite sample was collected from each of the half sections, with a depth of 0-6 inches below the surface and analyzed utilizing a PetroFlag (PetroFlag<sup>®</sup>) hydrocarbon detector. In addition to the bottom, a five point composite sample was collected from each of the pit walls and field screened for hydrocarbons.

Figure 1 outlines the pit sampling nomenclature and field screening results using the PetroFlag unit. Figure 2 is a GIS map of the pit outlining sample locations within the pit as well as background sample locations from the nearby uphill undisturbed soil.

Figure 1  
PetroFlag Results and Pit Sampling ID Layout



Facility Name: Chevron TR 21-20-597  
Remediation # 5949  
Facility ID: 284697

Name of Operator: Williams Production RMT Company  
Latitude: 39.601501 Longitude -108.302601  
Location (QtrQtr, Sec, Twp, Rng, Meridian): SENW, Sec 20, T5S, R97W, 6th PM

COGCC Operator # 96850  
County: Garfield

Table 1: PetroFlag Hydrocarbon Initial Field Screening Results

Sample ID	Results mg/kg
North Wall	238
East Wall	>5000
South Wall	2193
West Wall	288
North Bottom	>5000
South Bottom	2541

Note: All results are in mg/kg

Highlighted numbers indicate areas that warranted additional inspection and analysis

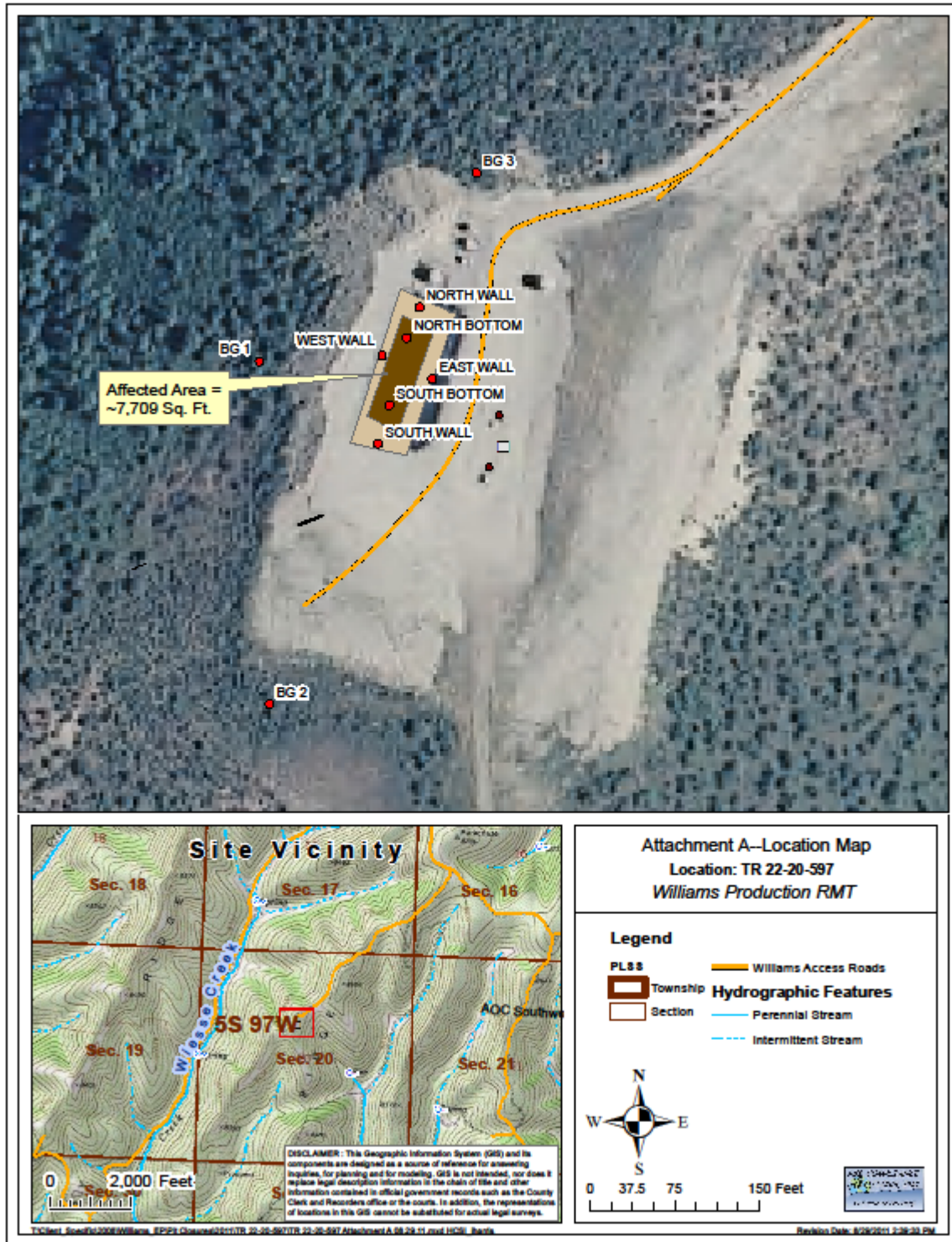
Facility Name: Chevron TR 21-20-597  
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Name of Operator: Williams Production RMT Company  
Latitude: 39.601501 Longitude -108.302601  
Location (QtrQtr, Sec, Twp, Rng, Meridian): SENW, Sec 20, T5S, R97W, 6th PM

COGCC Operator # 96850  
County: Garfield

Figure 2

GIS Map of Sampling Locations



Field screening results are provided in Table 1 and indicate that remediation is required due to TPH concentrations being above COGCC Table 910-1 standards.

### **Remediation Activities**

Soil exhibiting dark stains and hydrocarbon odors were located on the pit bottom as well as the east and south pit walls, indicating the potential presence of hydrocarbon concentrations exceeding 500 ppm and thus required remediation. The pit bottom, east wall, and south wall was excavated to a depth of approximately 2 feet in areas containing a potential hydrocarbon concentration above 500 ppm. Field screening at that depth indicated that soils located on the east wall and northern pit bottom still exceeded COGCC Table 910-1 concentrations and required additional excavation. The north pit bottom and eastern wall was excavated to a total of 6 feet, where discoloration within soil was no longer present at the excavated depth and field screening results indicated that hydrocarbon concentrations were below 500 ppm. Confirmation samples were collected and analyzed for COGCC Table 910-1.

- Confirmation samples, in accordance with Rule 905.b.(4), were collected from the sides walls at a position that was centered vertically and horizontally. These samples were collected for confirmation of compliance with COGCC Rule 910 and Table 910-1; as well as verification of field screening analysis. Two (2) additional grab samples were collected from the base of the pit, dividing the bottom of the pit into halves, which included the low point of the base, to demonstrate compliance in accordance with Rule 905.b.(1).
- A Trimble Geo XT 2008 was used to satisfy requirements outlined in COGCC Rule 215 for collecting GPS locations of each confirmation sample location from the pit walls and pit footprint.
- Visual inspection of the pit bottoms, field screening techniques, and sampling procedures were followed in accordance with Williams Highlands Pit Closure Plan (COGCC document #01175818).

Confirmation samples indicated that the pit bottom, as well as the southern pit wall still exceeded COGCC Table 910-1 for hydrocarbon concentrations exceeding 500 ppm in the DRO range. An additional 3 feet was excavated from the pit bottom and re-sampled for DRO. Confirmation samples collected at 8 feet below the original pit bottom indicated that the DRO concentrations were below COGCC Table 910-1, no additional remediation was required.

Analytical data presented in Table 2 provides results for the confirmation sampling performed after the initial 2 feet of soil was excavated from the pit footprint (raw analytical results are available for review in Appendix 1 of this report) at various depths and Table 3 provides confirmation sampling analysis of additional excavation performed on the pit bottom and southern wall (raw analytical results are available for review in Appendix 2)

## **Sample Analysis**

See attached Table 2 (additional detail provided in Appendix 1) for summary of pit bottom and side wall raw analytical results, and Table 3 (additional detail provided in Appendix 2) which provides raw analytical results for the additional excavation on the north and south pit bottom, as well as the southern pit wall analytical results, and Table 4 (additional detail provided in Appendix 3) for background analytical results. See attached Table 5 for confirmation analysis of the bioremediated soil located within the treatment cell (additional detail provided in Appendix 4).

## **Management of Stockpiled Material**

The pit liner was segregated according to material type and placed in a bermed containment. Plastic lining material was placed in the south end of the containment and felt liners were placed on the north end. High Plains Services compressed and collected the liners and bound them to pallets for transportation to be recycled.

Excavated soils from within the pit was placed in treatment cells, no thicker than 18 inches and treated on site with bioremediation product. See Table 5 (additional detail is provided in Appendix 4) for the treatment cell soil samples collected from the treated soil, providing confirmation that soil meets COGCC Table 910-1 allowable concentrations.

## **Backfill Material**

The backfill material utilized was from the stockpiled soil present on the east side of the pad from the initial construction of the pit.

- The soil was placed in lifts and was not compacted beyond the point of making an impenetrable layer but sufficient to suppose subsequent operations and prevent subsidence.
- The pit was reclaimed in accordance with the COGCC 1000 Series Rule in addition to all SUA/COA's per the land owner.

## **Exceptions to COGCC Table 910-1**

The only exceedances with COGCC Table 910-1 are within the confines of constituents listed for inorganics and metals (i.e. arsenic). Refer to Appendix 5 for the Sundry Notice for consideration of background arsenic concentrations in the immediate area of the subject facility.

## **Analytical Data Management**

See Appendix 1 for post excavated pit bottom and wall raw analytical data, Appendix 2 for additional excavation performed on the north and south pit bottom as well as the southern pit wall confirmation analytical data, and Appendix 3 for background analytical data. Appendix 4 provides confirmation analytical data for the successfully treated soil within the treatment cell.



## Figures

**Figure 3**



**Visual representation of the pit facing north after Excavation**

## **Summary Tables**

Table 2: Post Excavation Pit Bottom Analytical Results

Pit Bottom and Side Walls Post Excavation of 2 feet.	SAMPLE LOCATIONS					
	N. Bottom	N. Wall	S. Bottom	S. Wall	W. Wall	E. Wall
TEPH (DRO) (mg/kg)	1100	37	1400	910	420	13
TVPH (GRO) (mg/kg)	ND	ND	ND	ND	ND	ND
BENZENE (mg/kg)	ND	ND	ND	ND	ND	ND
TOLUENE (mg/kg)	ND	ND	ND	ND	ND	ND
ETHYLBENZENE (mg/kg)	ND	ND	ND	ND	ND	ND
XYLENE TOTAL (mg/kg)	ND	ND	ND	ND	ND	ND
ACENAPHTHENE (mg/kg)	ND	ND	ND	ND	ND	ND
ACENAPHTHYLENE (mg/kg)	NS	NS	NS	NS	NS	NS
ANTHRACENE (mg/kg)	ND	ND	ND	ND	ND	ND
BENZO(A)ANTHRACENE (mg/kg)	ND	ND	ND	ND	ND	ND
BENZO(A)PYRENE (mg/kg)	ND	ND	ND	ND	ND	ND
BENZO(B)FLUORANTHENE (mg/kg)	ND	ND	ND	ND	ND	ND
BENZO(G,H,I)PERYLENE (mg/kg)	ND	ND	ND	ND	ND	ND
BENZO(K)FLUORANTHENE (mg/kg)	ND	ND	ND	ND	ND	ND
CHRYSENE (mg/kg)	ND	ND	ND	ND	ND	ND
DIBENZO(A,H)ANTHRACENE (mg/kg)	ND	ND	ND	ND	ND	ND
FLUORANTHENE (mg/kg)	ND	ND	ND	ND	0.06	ND
FLUORENE (mg/kg)	ND	ND	ND	ND	0.042	ND
INDENO(1,2,3-CD)PYRENE (mg/kg)	ND	ND	ND	ND	ND	ND
1-METHYLNAPHTHALENE (mg/kg)	NS	NS	NS	NS	NS	NS
2-METHYLNAPHTHALENE (mg/kg)	NS	NS	NS	NS	NS	NS
NAPHTHALENE (mg/kg)	ND	ND	ND	ND	0.14	ND
PHENANTHRENE (mg/kg)	NS	NS	NS	NS	NS	NS
PYRENE (mg/kg)	ND	ND	0.045	ND	0.063	ND
ARSENIC (mg/kg)	2	1.7	2.2	2	4.1	2.8
BARIUM (mg/kg)	620	700	290	550	1000	380
CADMIUM (mg/kg)	ND	ND	0.78	ND	ND	ND
CHROMIUM (mg/kg)	17	43	16	22	30	38
CHROMIUM (III) (mg/kg)	17	43	16	22	31	37
CHROMIUM (IV) (mg/kg)	ND	ND	ND	ND	ND	ND
COPPER (mg/kg)	6.9	22	7.9	12	19	13
LEAD (mg/kg)	12	19	13	18	14	17
MERCURY (mg/kg)	ND	0.027	0.024	0.02	0.043	0.023
NICKEL (mg/kg)	10	24	10	14	18	18
SELENIUM (mg/kg)	ND	ND	ND	ND	ND	ND
SILVER (mg/kg)	ND	ND	ND	ND	ND	ND
ZINC (mg/kg)	70	72	48	67	78	64
ELECTRICAL CONDUCTIVITY (EC) (mmho/cm)	0.36	0.81	4.21	3.63	3.33	0.83
pH	8.35	8.14	8.21	8.58	8.65	7.75
SODIUM ABSORPTION RATIO (SAR)	0.7	10	71.2	62.8	42.7	2.8
CALCIUM (ppm)	69	113	281	254	503	158
MAGNESIUM (ppm)	15	21	35	34	46	30
SODIUM (ppm)	24	441	4811	4024	3740	144

Note: all results are in, mg/kg = milligram per kilogram, unless noted  
Exceedances are highlighted in yellow.

Table 3: Pit Bottom and Southern Wall – Additional Excavation

	North Bottom	South Bottom	South Wall
<b>Post Excavation Pit Bottom @ 6 ft</b>			
TEPH (DRO)	31	71	110

Note: All results are in, mg/L = milligrams per liter, unless noted otherwise

Table 4: Background Analytical Data

	Arsenic	Sodium Absorption Ratio (unitless)	Electro Conductivity (mmhos/cm)	pH (unitless)
BKGD 1	5.4	46.8	3.01	7.02
BKGD 2	5.2			
BKGD 3	5.7			

All results are in, mg/kg = milligram per kilogram, unless noted otherwise

Table 5: Treatment Cell Confirmation Analytical Data

Treatment Cell Confirmation	SAMPLE LOCATIONS	
	North Side	South South
TEPH (DRO) (mg/kg)	96	47
TVPH (GRO) (mg/kg)	ND	ND
BENZENE (mg/kg)	ND	ND
TOLUENE (mg/kg)	ND	ND
ETHYLBENZENE (mg/kg)	ND	ND
XYLENE TOTAL (mg/kg)	ND	ND
ACENAPHTHENE (mg/kg)	ND	ND
ACENAPHTHYLENE (mg/kg)	NS	NS
ANTHRACENE (mg/kg)	ND	ND
BENZO(A)ANTHRACENE (mg/kg)	ND	ND
BENZO(A)PYRENE (mg/kg)	ND	ND
BENZO(B)FLUORANTHENE (mg/kg)	ND	ND
BENZO(G,H,I)PERYLENE (mg/kg)	ND	ND
BENZO(K)FLUORANTHENE (mg/kg)	ND	ND
CHRYSENE (mg/kg)	ND	ND
DIBENZO(A,H)ANTHRACENE (mg/kg)	ND	ND
FLUORANTHENE (mg/kg)	ND	ND
FLUORENE (mg/kg)	ND	ND
INDENO(1,2,3-CD)PYRENE (mg/kg)	ND	ND
1-METHYLNAPHTHALENE (mg/kg)	NS	NS
2-METHYLNAPHTHALENE (mg/kg)	NS	NS
NAPHTHALENE (mg/kg)	ND	ND
PHENANTHRENE (mg/kg)	NS	NS
PYRENE (mg/kg)	ND	ND
ARSENIC (mg/kg)	2.7	3.9
BARIUM (mg/kg)	500	540
CADMIUM (mg/kg)	ND	ND
CHROMIUM (mg/kg)	17	16
CHROMIUM (III) (mg/kg)	17	16
CHROMIUM (IV) (mg/kg)	ND	ND
COPPER (mg/kg)	10	10
LEAD (mg/kg)	22	20
MERCURY (mg/kg)	ND	ND
NICKEL (mg/kg)	12	11
SELENIUM (mg/kg)	ND	ND
SILVER (mg/kg)	ND	ND
ZINC (mg/kg)	48	44
ELECTRICAL CONDUCTIVITY (EC) (mmho/cm)	3.03	2.22
pH	8.95	8.93
SODIUM ABSORPTION RATIO (SAR)	30.3	26.7

Note: All results are reported in mg/kg, unless otherwise noted.

## **Appendix 1: Pit Bottom and Walls Confirmation after 2 ft Raw Analytical Data**



08-Sep-2011

Kris Rowe  
HRL Compliance Solutions  
744 Horizon Ct. Suite 140  
Grand Junction, CO 81506

Re: **Williams TR 22-20-597 Pad LOE 8/24/11**

Work Order: **1108912**

Dear Kris,

ALS Environmental received 6 samples on 30-Aug-2011 10:30 AM for the analyses presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested.

QC sample results for this data met laboratory specifications. Any exceptions are noted in the Case Narrative, or noted with qualifiers in the report or QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained from ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

The total number of pages in this report is 49.

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

Sincerely,

A handwritten signature in cursive script that reads "Ann Preston".

Electronically approved by: Ann Preston

Ann Preston  
Project Manager



Certificate No: IL100452

ADDRESS 3352 128th Avenue Holland, Michigan 49424-9263 | PHONE (616) 399-6070 | FAX (616) 399-6185

ALS GROUP USA, CORP Part of the ALS Laboratory Group A Campbell Brothers Limited Company

Environmental The ALS logo, which is a stylized blue triangle with a yellow flame inside.

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RIGHT SOLUTIONS RIGHT PARTNER



**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11  
**Work Order:** 1108912

## Work Order Sample Summary

<u>Lab Samp ID</u>	<u>Client Sample ID</u>	<u>Matrix</u>	<u>Tag Number</u>	<u>Collection Date</u>	<u>Date Received</u>	<u>Hold</u>
1108912-01	TR 22-20-597 N. Bottom	Soil		8/24/2011 14:20	8/30/2011 10:30	<input type="checkbox"/>
1108912-02	TR 22-20-597 N. Wall	Soil		8/24/2011 13:45	8/30/2011 10:30	<input type="checkbox"/>
1108912-03	TR 22-20-597 S Bottom	Soil		8/24/2011 14:15	8/30/2011 10:30	<input type="checkbox"/>
1108912-04	TR 22-20-597 S. Wall	Soil		8/24/2011 14:00	8/30/2011 10:30	<input type="checkbox"/>
1108912-05	TR 22-20-597 W. Wall	Soil		8/24/2011 13:50	8/30/2011 10:30	<input type="checkbox"/>
1108912-06	TR 22-20-597 E. Wall	Soil		8/24/2011 14:10	8/30/2011 10:30	<input type="checkbox"/>

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**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11  
**Work Order:** 1108912

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**Case Narrative**

Batch 35228 sample TR-22-20-597 S Bottom MS/MSD recoveries for Barium, Lead and Zinc were above control limits, The results for these elements in the parent sample may be biased high due to matrix interference. The MS/MSD recoveries for Silver were below control limits. The results for Silver in the parent sample may be biased low due to matrix interference. The MS recoveries for Chromium and Copper were below control limits, but both the MSD recoveries and RPDs met quality control criteria.

Batch 35235 sample 1108912-04B PAH surrogate recoveries were above control limits, but all PAH compounds were ND. Samples 1108912-01 and 1108912-03 had a few PAH compounds that were reported with a dilution and raised reporting limits due to matrix interference.

The samples for pH were received after the hold time had expired.

**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11  
**WorkOrder:** 1108912

## **QUALIFIERS, ACRONYMS, UNITS**

<b><u>Qualifier</u></b>	<b><u>Description</u></b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL

<b><u>Acronym</u></b>	<b><u>Description</u></b>
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
TDL	Target Detection Limit

<b><u>Units Reported</u></b>	<b><u>Description</u></b>
% of sample	Percent of Sample
µg/Kg-dry as noted	Micrograms per Kilogram Dry Weight
mg/Kg-dry	Milligrams per Kilogram Dry Weight
s.u.	Standard Units

# ALS Group USA, Corp

Date: 08-Sep-11

**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11  
**Sample ID:** TR 22-20-597 N. Bottom  
**Collection Date:** 8/24/2011 02:20 PM

**Work Order:** 1108912  
**Lab ID:** 1108912-01  
**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<b>DIESEL RANGE ORGANICS BY GC-FID</b>						
			<b>SW8015M</b>		Prep Date: <b>8/30/2011</b>	Analyst: <b>RM</b>
<b>DRO (C10-C28)</b>	<b>1,100</b>		<b>4.7</b>	<b>mg/Kg-dry</b>	1	8/31/2011 05:35 PM
Surr: 4-Terphenyl-d14	60.7		39-115	%REC	1	8/31/2011 05:35 PM
<b>GASOLINE RANGE ORGANICS BY GC-FID</b>						
			<b>SW8015</b>			Analyst: <b>RM</b>
GRO (C6-C10)	ND		5.8	mg/Kg-dry	100	8/31/2011 09:51 PM
Surr: Toluene-d8	105		50-150	%REC	100	8/31/2011 09:51 PM
<b>MERCURY BY CVAA</b>						
			<b>SW7471</b>		Prep Date: <b>8/31/2011</b>	Analyst: <b>LR</b>
Mercury	ND		0.020	mg/Kg-dry	1	9/1/2011 02:28 PM
<b>METALS BY ICP-MS</b>						
			<b>SW6020A</b>		Prep Date: <b>8/30/2011</b>	Analyst: <b>CES</b>
Arsenic	2.0		1.8	mg/Kg-dry	4	8/31/2011 07:09 PM
Barium	620		1.8	mg/Kg-dry	4	8/31/2011 07:09 PM
Cadmium	ND		0.70	mg/Kg-dry	4	8/31/2011 07:09 PM
Chromium	17		1.8	mg/Kg-dry	4	8/31/2011 07:09 PM
Copper	6.9		1.8	mg/Kg-dry	4	8/31/2011 07:09 PM
Lead	12		1.8	mg/Kg-dry	4	9/1/2011 12:42 PM
Nickel	10		1.8	mg/Kg-dry	4	8/31/2011 07:09 PM
Selenium	ND		1.8	mg/Kg-dry	4	8/31/2011 07:09 PM
Silver	ND		1.8	mg/Kg-dry	4	8/31/2011 07:09 PM
Zinc	70		3.5	mg/Kg-dry	4	8/31/2011 07:09 PM
<b>SUBCONTRACTED ANALYSES</b>						
Subcontracted Analyses	Rcvd 9/6/11		<b>SUBCONTRACT</b>			Analyst: <b>A&amp;LGL</b>
			as noted		1	9/6/2011
<b>SEMI-VOLATILE ORGANIC COMPOUNDS</b>						
			<b>SW8270</b>		Prep Date: <b>8/30/2011</b>	Analyst: <b>HL</b>
Acenaphthene	ND		680	µg/Kg-dry	20	9/1/2011 11:02 PM
Anthracene	ND		34	µg/Kg-dry	1	8/31/2011 07:16 PM
Benzo(a)anthracene	ND		34	µg/Kg-dry	1	8/31/2011 07:16 PM
Benzo(a)pyrene	ND		34	µg/Kg-dry	1	8/31/2011 07:16 PM
Benzo(b)fluoranthene	ND		34	µg/Kg-dry	1	8/31/2011 07:16 PM
Benzo(g,h,i)perylene	ND		34	µg/Kg-dry	1	8/31/2011 07:16 PM
Benzo(k)fluoranthene	ND		34	µg/Kg-dry	1	8/31/2011 07:16 PM
Chrysene	ND		34	µg/Kg-dry	1	8/31/2011 07:16 PM
Dibenzo(a,h)anthracene	ND		34	µg/Kg-dry	1	8/31/2011 07:16 PM
Fluoranthene	ND		34	µg/Kg-dry	1	8/31/2011 07:16 PM
Fluorene	ND		680	µg/Kg-dry	20	9/1/2011 11:02 PM
Indeno(1,2,3-cd)pyrene	ND		34	µg/Kg-dry	1	8/31/2011 07:16 PM
Naphthalene	ND		34	µg/Kg-dry	1	8/31/2011 07:16 PM
Pyrene	ND		34	µg/Kg-dry	1	8/31/2011 07:16 PM
Surr: 2,4,6-Tribromophenol	139		34-140	%REC	1	8/31/2011 07:16 PM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

# ALS Group USA, Corp

Date: 08-Sep-11

**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11  
**Sample ID:** TR 22-20-597 N. Bottom  
**Collection Date:** 8/24/2011 02:20 PM

**Work Order:** 1108912  
**Lab ID:** 1108912-01  
**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<i>Surr: 2-Fluorobiphenyl</i>	86.4		12-100	%REC	20	9/1/2011 11:02 PM
<i>Surr: 2-Fluorophenol</i>	101		33-117	%REC	1	8/31/2011 07:16 PM
<i>Surr: 4-Terphenyl-d14</i>	106		25-137	%REC	1	8/31/2011 07:16 PM
<i>Surr: Nitrobenzene-d5</i>	66.8		37-107	%REC	20	9/1/2011 11:02 PM
<i>Surr: Phenol-d6</i>	103		40-106	%REC	1	8/31/2011 07:16 PM
<b>VOLATILE ORGANIC COMPOUNDS</b>			<b>SW8260</b>			Analyst: <b>BG</b>
Benzene	ND		120	µg/Kg-dry	100	9/1/2011 08:28 AM
Ethylbenzene	ND		120	µg/Kg-dry	100	9/1/2011 08:28 AM
m,p-Xylene	ND		120	µg/Kg-dry	100	9/1/2011 08:28 AM
o-Xylene	ND		120	µg/Kg-dry	100	9/1/2011 08:28 AM
Toluene	ND		120	µg/Kg-dry	100	9/1/2011 08:28 AM
Xylenes, Total	ND		350	µg/Kg-dry	100	9/1/2011 08:28 AM
<i>Surr: 1,2-Dichloroethane-d4</i>	103		70-120	%REC	100	9/1/2011 08:28 AM
<i>Surr: 4-Bromofluorobenzene</i>	102		75-120	%REC	100	9/1/2011 08:28 AM
<i>Surr: Dibromofluoromethane</i>	94.6		85-115	%REC	100	9/1/2011 08:28 AM
<i>Surr: Toluene-d8</i>	97.4		85-115	%REC	100	9/1/2011 08:28 AM
<b>CHROMIUM, TRIVALENT</b>			<b>CALCULATION</b>			Analyst: <b>JJG</b>
Chromium, Trivalent	17			mg/kg-dry	1	9/2/2011 02:35 PM
<b>CHROMIUM, HEXAVALENT</b>			<b>SW7196A</b>		Prep Date: <b>9/1/2011</b>	Analyst: <b>MB</b>
Chromium, Hexavalent	ND		0.58	mg/Kg-dry	1	9/2/2011 01:00 PM
<b>MOISTURE</b>			<b>A2540 G</b>			Analyst: <b>CG</b>
Moisture	14		0.050	% of sample	1	8/30/2011 02:49 PM
<b>PH</b>			<b>SW9045D</b>			Analyst: <b>JJG</b>
pH	8.35	H		s.u.	1	8/30/2011 10:15 AM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

# ALS Group USA, Corp

Date: 08-Sep-11

Client: HRL Compliance Solutions

Project: Williams TR 22-20-597 Pad LOE 8/24/11

Work Order: 1108912

Sample ID: TR 22-20-597 N. Wall

Lab ID: 1108912-02

Collection Date: 8/24/2011 01:45 PM

Matrix: SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<b>DIESEL RANGE ORGANICS BY GC-FID</b>						
<b>DRO (C10-C28)</b>	<b>37</b>		<b>SW8015M</b>		Prep Date: <b>8/30/2011</b>	Analyst: <b>RM</b>
			<b>4.9</b>	<b>mg/Kg-dry</b>	<b>1</b>	8/31/2011 05:35 PM
Surr: 4-Terphenyl-d14	89.7		39-115	%REC	1	8/31/2011 05:35 PM
<b>GASOLINE RANGE ORGANICS BY GC-FID</b>						
<b>GRO (C6-C10)</b>	<b>ND</b>		<b>SW8015</b>			Analyst: <b>RM</b>
			<b>5.9</b>	<b>mg/Kg-dry</b>	<b>100</b>	8/31/2011 10:17 PM
Surr: Toluene-d8	106		50-150	%REC	100	8/31/2011 10:17 PM
<b>MERCURY BY CVAA</b>						
<b>Mercury</b>	<b>0.027</b>		<b>SW7471</b>		Prep Date: <b>8/31/2011</b>	Analyst: <b>LR</b>
			<b>0.018</b>	<b>mg/Kg-dry</b>	<b>1</b>	9/1/2011 02:30 PM
<b>METALS BY ICP-MS</b>						
<b>Arsenic</b>	<b>1.8</b>		<b>SW6020A</b>		Prep Date: <b>8/30/2011</b>	Analyst: <b>CES</b>
			<b>1.6</b>	<b>mg/Kg-dry</b>	<b>4</b>	8/31/2011 07:14 PM
<b>Barium</b>	<b>700</b>		<b>16</b>	<b>mg/Kg-dry</b>	<b>40</b>	9/1/2011 12:47 PM
Cadmium	ND		0.62	mg/Kg-dry	4	8/31/2011 07:14 PM
<b>Chromium</b>	<b>43</b>		<b>1.6</b>	<b>mg/Kg-dry</b>	<b>4</b>	8/31/2011 07:14 PM
<b>Copper</b>	<b>22</b>		<b>1.6</b>	<b>mg/Kg-dry</b>	<b>4</b>	8/31/2011 07:14 PM
<b>Lead</b>	<b>19</b>		<b>1.6</b>	<b>mg/Kg-dry</b>	<b>4</b>	9/1/2011 12:53 PM
<b>Nickel</b>	<b>24</b>		<b>1.6</b>	<b>mg/Kg-dry</b>	<b>4</b>	8/31/2011 07:14 PM
Selenium	ND		1.6	mg/Kg-dry	4	8/31/2011 07:14 PM
Silver	ND		1.6	mg/Kg-dry	4	8/31/2011 07:14 PM
<b>Zinc</b>	<b>72</b>		<b>3.1</b>	<b>mg/Kg-dry</b>	<b>4</b>	8/31/2011 07:14 PM
<b>SUBCONTRACTED ANALYSES</b>						
<b>Subcontracted Analyses</b>		<b>Rcvd 9/6/11</b>	<b>SUBCONTRACT</b>			Analyst: <b>A&amp;LGL</b>
			<b>as noted</b>		<b>1</b>	9/6/2011
<b>SEMI-VOLATILE ORGANIC COMPOUNDS</b>						
<b>Acenaphthene</b>	<b>ND</b>		<b>SW8270</b>		Prep Date: <b>8/30/2011</b>	Analyst: <b>HL</b>
			<b>35</b>	<b>µg/Kg-dry</b>	<b>1</b>	8/31/2011 06:45 PM
<b>Anthracene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	<b>1</b>	8/31/2011 06:45 PM
<b>Benzo(a)anthracene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	<b>1</b>	8/31/2011 06:45 PM
<b>Benzo(a)pyrene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	<b>1</b>	8/31/2011 06:45 PM
<b>Benzo(b)fluoranthene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	<b>1</b>	8/31/2011 06:45 PM
<b>Benzo(g,h,i)perylene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	<b>1</b>	8/31/2011 06:45 PM
<b>Benzo(k)fluoranthene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	<b>1</b>	8/31/2011 06:45 PM
<b>Chrysene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	<b>1</b>	8/31/2011 06:45 PM
<b>Dibenzo(a,h)anthracene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	<b>1</b>	8/31/2011 06:45 PM
<b>Fluoranthene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	<b>1</b>	8/31/2011 06:45 PM
<b>Fluorene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	<b>1</b>	8/31/2011 06:45 PM
<b>Indeno(1,2,3-cd)pyrene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	<b>1</b>	8/31/2011 06:45 PM
<b>Naphthalene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	<b>1</b>	8/31/2011 06:45 PM
<b>Pyrene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	<b>1</b>	8/31/2011 06:45 PM
Surr: 2,4,6-Tribromophenol	111		34-140	%REC	1	8/31/2011 06:45 PM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

# ALS Group USA, Corp

Date: 08-Sep-11

**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11  
**Sample ID:** TR 22-20-597 N. Wall  
**Collection Date:** 8/24/2011 01:45 PM

**Work Order:** 1108912  
**Lab ID:** 1108912-02  
**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<i>Surr: 2-Fluorobiphenyl</i>	69.8		12-100	%REC	1	8/31/2011 06:45 PM
<i>Surr: 2-Fluorophenol</i>	77.3		33-117	%REC	1	8/31/2011 06:45 PM
<i>Surr: 4-Terphenyl-d14</i>	98.3		25-137	%REC	1	8/31/2011 06:45 PM
<i>Surr: Nitrobenzene-d5</i>	67.1		37-107	%REC	1	8/31/2011 06:45 PM
<i>Surr: Phenol-d6</i>	75.9		40-106	%REC	1	8/31/2011 06:45 PM
<b>VOLATILE ORGANIC COMPOUNDS</b>			<b>SW8260</b>			Analyst: <b>BG</b>
Benzene	ND		120	µg/Kg-dry	100	9/1/2011 08:53 AM
Ethylbenzene	ND		120	µg/Kg-dry	100	9/1/2011 08:53 AM
m,p-Xylene	ND		120	µg/Kg-dry	100	9/1/2011 08:53 AM
o-Xylene	ND		120	µg/Kg-dry	100	9/1/2011 08:53 AM
Toluene	ND		120	µg/Kg-dry	100	9/1/2011 08:53 AM
Xylenes, Total	ND		360	µg/Kg-dry	100	9/1/2011 08:53 AM
<i>Surr: 1,2-Dichloroethane-d4</i>	100		70-120	%REC	100	9/1/2011 08:53 AM
<i>Surr: 4-Bromofluorobenzene</i>	99.8		75-120	%REC	100	9/1/2011 08:53 AM
<i>Surr: Dibromofluoromethane</i>	93.2		85-115	%REC	100	9/1/2011 08:53 AM
<i>Surr: Toluene-d8</i>	97.5		85-115	%REC	100	9/1/2011 08:53 AM
<b>CHROMIUM, TRIVALENT</b>			<b>CALCULATION</b>			Analyst: <b>JJG</b>
Chromium, Trivalent	43			mg/kg-dry	1	9/2/2011 02:35 PM
<b>CHROMIUM, HEXAVALENT</b>			<b>SW7196A</b>		Prep Date: <b>9/1/2011</b>	Analyst: <b>MB</b>
Chromium, Hexavalent	ND		0.59	mg/Kg-dry	1	9/2/2011 01:00 PM
<b>MOISTURE</b>			<b>A2540 G</b>			Analyst: <b>CG</b>
Moisture	16		0.050	% of sample	1	8/30/2011 02:49 PM
<b>PH</b>			<b>SW9045D</b>			Analyst: <b>JJG</b>
pH	8.14	H		s.u.	1	8/30/2011 10:15 AM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

# ALS Group USA, Corp

Date: 08-Sep-11

Client: HRL Compliance Solutions

Project: Williams TR 22-20-597 Pad LOE 8/24/11

Work Order: 1108912

Sample ID: TR 22-20-597 S Bottom

Lab ID: 1108912-03

Collection Date: 8/24/2011 02:15 PM

Matrix: SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
DIESEL RANGE ORGANICS BY GC-FID			SW8015M		Prep Date: 8/30/2011	Analyst: RM
DRO (C10-C28)	1,400		4.6	mg/Kg-dry	1	8/31/2011 05:58 PM
Surr: 4-Terphenyl-d14	80.9		39-115	%REC	1	8/31/2011 05:58 PM
GASOLINE RANGE ORGANICS BY GC-FID			SW8015			Analyst: RM
GRO (C6-C10)	ND		5.7	mg/Kg-dry	100	8/31/2011 10:43 PM
Surr: Toluene-d8	103		50-150	%REC	100	8/31/2011 10:43 PM
MERCURY BY CVAA			SW7471		Prep Date: 8/31/2011	Analyst: LR
Mercury	0.024		0.021	mg/Kg-dry	1	9/1/2011 02:32 PM
METALS BY ICP-MS			SW6020A		Prep Date: 8/30/2011	Analyst: CES
Arsenic	2.2		1.7	mg/Kg-dry	4	8/31/2011 07:19 PM
Barium	290		1.7	mg/Kg-dry	4	8/31/2011 07:19 PM
Cadmium	0.78		0.70	mg/Kg-dry	4	8/31/2011 07:19 PM
Chromium	16		1.7	mg/Kg-dry	4	8/31/2011 07:19 PM
Copper	7.9		1.7	mg/Kg-dry	4	8/31/2011 07:19 PM
Lead	13		1.7	mg/Kg-dry	4	9/1/2011 12:58 PM
Nickel	10		1.7	mg/Kg-dry	4	8/31/2011 07:19 PM
Selenium	ND		1.7	mg/Kg-dry	4	8/31/2011 07:19 PM
Silver	ND		1.7	mg/Kg-dry	4	8/31/2011 07:19 PM
Zinc	48		3.5	mg/Kg-dry	4	8/31/2011 07:19 PM
SUBCONTRACTED ANALYSES			SUBCONTRACT			Analyst: A&LGL
Subcontracted Analyses	Rcvd 9/6/11		as noted		1	9/6/2011
SEMI-VOLATILE ORGANIC COMPOUNDS			SW8270		Prep Date: 8/30/2011	Analyst: HL
Acenaphthene	ND		670	µg/Kg-dry	20	9/1/2011 11:33 PM
Anthracene	ND		670	µg/Kg-dry	20	9/1/2011 11:33 PM
Benzo(a)anthracene	ND		33	µg/Kg-dry	1	8/31/2011 06:14 PM
Benzo(a)pyrene	ND		33	µg/Kg-dry	1	8/31/2011 06:14 PM
Benzo(b)fluoranthene	ND		33	µg/Kg-dry	1	8/31/2011 06:14 PM
Benzo(g,h,i)perylene	ND		33	µg/Kg-dry	1	8/31/2011 06:14 PM
Benzo(k)fluoranthene	ND		33	µg/Kg-dry	1	8/31/2011 06:14 PM
Chrysene	ND		33	µg/Kg-dry	1	8/31/2011 06:14 PM
Dibenzo(a,h)anthracene	ND		33	µg/Kg-dry	1	8/31/2011 06:14 PM
Fluoranthene	ND		670	µg/Kg-dry	20	9/1/2011 11:33 PM
Fluorene	ND		670	µg/Kg-dry	20	9/1/2011 11:33 PM
Indeno(1,2,3-cd)pyrene	ND		33	µg/Kg-dry	1	8/31/2011 06:14 PM
Naphthalene	ND		33	µg/Kg-dry	1	8/31/2011 06:14 PM
Pyrene	45		33	µg/Kg-dry	1	8/31/2011 06:14 PM
Surr: 2,4,6-Tribromophenol	85.2		34-140	%REC	20	9/1/2011 11:33 PM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.



# ALS Group USA, Corp

Date: 08-Sep-11

**Client:** HRL Compliance Solutions

**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

**Work Order:** 1108912

**Sample ID:** TR 22-20-597 S Bottom

**Lab ID:** 1108912-03

**Collection Date:** 8/24/2011 02:15 PM

**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<i>Surr: 2-Fluorobiphenyl</i>	84.8		12-100	%REC	20	9/1/2011 11:33 PM
<i>Surr: 2-Fluorophenol</i>	80.5		33-117	%REC	1	8/31/2011 06:14 PM
<i>Surr: 4-Terphenyl-d14</i>	83.5		25-137	%REC	1	8/31/2011 06:14 PM
<i>Surr: Nitrobenzene-d5</i>	71.2		37-107	%REC	20	9/1/2011 11:33 PM
<i>Surr: Phenol-d6</i>	88.1		40-106	%REC	1	8/31/2011 06:14 PM
<b>VOLATILE ORGANIC COMPOUNDS</b>			<b>SW8260</b>			Analyst: <b>BG</b>
Benzene	ND		110	µg/Kg-dry	100	9/1/2011 09:19 AM
Ethylbenzene	ND		110	µg/Kg-dry	100	9/1/2011 09:19 AM
m,p-Xylene	ND		110	µg/Kg-dry	100	9/1/2011 09:19 AM
o-Xylene	ND		110	µg/Kg-dry	100	9/1/2011 09:19 AM
Toluene	ND		110	µg/Kg-dry	100	9/1/2011 09:19 AM
Xylenes, Total	ND		340	µg/Kg-dry	100	9/1/2011 09:19 AM
<i>Surr: 1,2-Dichloroethane-d4</i>	101		70-120	%REC	100	9/1/2011 09:19 AM
<i>Surr: 4-Bromofluorobenzene</i>	100		75-120	%REC	100	9/1/2011 09:19 AM
<i>Surr: Dibromofluoromethane</i>	94.7		85-115	%REC	100	9/1/2011 09:19 AM
<i>Surr: Toluene-d8</i>	97.2		85-115	%REC	100	9/1/2011 09:19 AM
<b>CHROMIUM, TRIVALENT</b>			<b>CALCULATION</b>			Analyst: <b>JJG</b>
Chromium, Trivalent	16			mg/kg-dry	1	9/2/2011 02:35 PM
<b>CHROMIUM, HEXAVALENT</b>			<b>SW7196A</b>		Prep Date: <b>9/1/2011</b>	Analyst: <b>MB</b>
Chromium, Hexavalent	ND		0.56	mg/Kg-dry	1	9/2/2011 01:00 PM
<b>MOISTURE</b>			<b>A2540 G</b>			Analyst: <b>CG</b>
Moisture	12		0.050	% of sample	1	8/30/2011 02:49 PM
<b>PH</b>			<b>SW9045D</b>			Analyst: <b>JJG</b>
pH	8.21	H		s.u.	1	8/30/2011 10:15 AM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

# ALS Group USA, Corp

Date: 08-Sep-11

**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11  
**Sample ID:** TR 22-20-597 S. Wall  
**Collection Date:** 8/24/2011 02:00 PM

**Work Order:** 1108912  
**Lab ID:** 1108912-04  
**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<b>DIESEL RANGE ORGANICS BY GC-FID</b>						
<b>DRO (C10-C28)</b>	<b>910</b>		<b>SW8015M</b>		Prep Date: <b>8/31/2011</b>	Analyst: <b>RM</b>
			<b>4.8</b>	<b>mg/Kg-dry</b>	1	9/1/2011 03:18 PM
Surr: 4-Terphenyl-d14	51.2		39-115	%REC	1	9/1/2011 03:18 PM
<b>GASOLINE RANGE ORGANICS BY GC-FID</b>						
<b>GRO (C6-C10)</b>	<b>ND</b>		<b>SW8015</b>			Analyst: <b>RM</b>
			<b>5.9</b>	<b>mg/Kg-dry</b>	100	8/31/2011 11:09 PM
Surr: Toluene-d8	106		50-150	%REC	100	8/31/2011 11:09 PM
<b>MERCURY BY CVAA</b>						
<b>Mercury</b>	<b>0.020</b>		<b>SW7471</b>		Prep Date: <b>8/31/2011</b>	Analyst: <b>LR</b>
			<b>0.019</b>	<b>mg/Kg-dry</b>	1	9/1/2011 02:34 PM
<b>METALS BY ICP-MS</b>						
<b>Arsenic</b>	<b>2.0</b>		<b>SW6020A</b>		Prep Date: <b>8/30/2011</b>	Analyst: <b>CES</b>
			<b>1.6</b>	<b>mg/Kg-dry</b>	4	8/31/2011 07:45 PM
<b>Barium</b>	<b>550</b>		<b>1.6</b>	<b>mg/Kg-dry</b>	4	8/31/2011 07:45 PM
Cadmium	ND		0.66	mg/Kg-dry	4	8/31/2011 07:45 PM
<b>Chromium</b>	<b>22</b>		<b>1.6</b>	<b>mg/Kg-dry</b>	4	8/31/2011 07:45 PM
<b>Copper</b>	<b>12</b>		<b>1.6</b>	<b>mg/Kg-dry</b>	4	8/31/2011 07:45 PM
<b>Lead</b>	<b>18</b>		<b>1.6</b>	<b>mg/Kg-dry</b>	4	9/1/2011 01:24 PM
<b>Nickel</b>	<b>14</b>		<b>1.6</b>	<b>mg/Kg-dry</b>	4	8/31/2011 07:45 PM
Selenium	ND		1.6	mg/Kg-dry	4	8/31/2011 07:45 PM
Silver	ND		1.6	mg/Kg-dry	4	8/31/2011 07:45 PM
<b>Zinc</b>	<b>67</b>		<b>3.3</b>	<b>mg/Kg-dry</b>	4	8/31/2011 07:45 PM
<b>SUBCONTRACTED ANALYSES</b>						
<b>Subcontracted Analyses</b>		<b>Rcvd 9/6/11</b>	<b>SUBCONTRACT</b>			Analyst: <b>A&amp;LGL</b>
			<b>as noted</b>		1	9/6/2011
<b>SEMI-VOLATILE ORGANIC COMPOUNDS</b>						
<b>Acenaphthene</b>	<b>ND</b>		<b>SW8270</b>		Prep Date: <b>8/31/2011</b>	Analyst: <b>HL</b>
			<b>350</b>	<b>µg/Kg-dry</b>	10	9/8/2011 03:59 PM
<b>Anthracene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	1	9/2/2011 01:07 AM
<b>Benzo(a)anthracene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	1	9/2/2011 01:07 AM
<b>Benzo(a)pyrene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	1	9/2/2011 01:07 AM
<b>Benzo(b)fluoranthene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	1	9/2/2011 01:07 AM
<b>Benzo(g,h,i)perylene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	1	9/2/2011 01:07 AM
<b>Benzo(k)fluoranthene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	1	9/2/2011 01:07 AM
<b>Chrysene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	1	9/2/2011 01:07 AM
<b>Dibenzo(a,h)anthracene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	1	9/2/2011 01:07 AM
<b>Fluoranthene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	1	9/2/2011 01:07 AM
<b>Fluorene</b>	<b>ND</b>		<b>350</b>	<b>µg/Kg-dry</b>	10	9/8/2011 03:59 PM
<b>Indeno(1,2,3-cd)pyrene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	1	9/2/2011 01:07 AM
<b>Naphthalene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	1	9/2/2011 01:07 AM
<b>Pyrene</b>	<b>ND</b>		<b>35</b>	<b>µg/Kg-dry</b>	1	9/2/2011 01:07 AM
Surr: 2,4,6-Tribromophenol	134		34-140	%REC	1	9/2/2011 01:07 AM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

# ALS Group USA, Corp

Date: 08-Sep-11

**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11  
**Sample ID:** TR 22-20-597 S. Wall  
**Collection Date:** 8/24/2011 02:00 PM

**Work Order:** 1108912  
**Lab ID:** 1108912-04  
**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<i>Surr: 2-Fluorobiphenyl</i>	118	S	12-100	%REC	10	9/8/2011 03:59 PM
<i>Surr: 2-Fluorophenol</i>	94.3		33-117	%REC	1	9/2/2011 01:07 AM
<i>Surr: 4-Terphenyl-d14</i>	99.6		25-137	%REC	1	9/2/2011 01:07 AM
<i>Surr: Nitrobenzene-d5</i>	123	S	37-107	%REC	10	9/8/2011 03:59 PM
<i>Surr: Phenol-d6</i>	90.9		40-106	%REC	1	9/2/2011 01:07 AM
<b>VOLATILE ORGANIC COMPOUNDS</b>			<b>SW8260</b>			Analyst: <b>BG</b>
Benzene	ND		120	µg/Kg-dry	100	9/1/2011 09:44 AM
Ethylbenzene	ND		120	µg/Kg-dry	100	9/1/2011 09:44 AM
m,p-Xylene	ND		120	µg/Kg-dry	100	9/1/2011 09:44 AM
o-Xylene	ND		120	µg/Kg-dry	100	9/1/2011 09:44 AM
Toluene	ND		120	µg/Kg-dry	100	9/1/2011 09:44 AM
Xylenes, Total	ND		350	µg/Kg-dry	100	9/1/2011 09:44 AM
<i>Surr: 1,2-Dichloroethane-d4</i>	101		70-120	%REC	100	9/1/2011 09:44 AM
<i>Surr: 4-Bromofluorobenzene</i>	98.0		75-120	%REC	100	9/1/2011 09:44 AM
<i>Surr: Dibromofluoromethane</i>	94.4		85-115	%REC	100	9/1/2011 09:44 AM
<i>Surr: Toluene-d8</i>	94.7		85-115	%REC	100	9/1/2011 09:44 AM
<b>CHROMIUM, TRIVALENT</b>			<b>CALCULATION</b>			Analyst: <b>JJG</b>
Chromium, Trivalent	22			mg/kg-dry	1	9/2/2011 02:35 PM
<b>CHROMIUM, HEXAVALENT</b>			<b>SW7196A</b>		Prep Date: <b>9/1/2011</b>	Analyst: <b>MB</b>
Chromium, Hexavalent	ND		0.57	mg/Kg-dry	1	9/2/2011 01:00 PM
<b>MOISTURE</b>			<b>A2540 G</b>			Analyst: <b>CG</b>
Moisture	15		0.050	% of sample	1	8/30/2011 02:49 PM
<b>PH</b>			<b>SW9045D</b>			Analyst: <b>JJG</b>
pH	8.58	H		s.u.	1	8/30/2011 10:15 AM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

# ALS Group USA, Corp

Date: 08-Sep-11

**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11  
**Sample ID:** TR 22-20-597 W. Wall  
**Collection Date:** 8/24/2011 01:50 PM

**Work Order:** 1108912  
**Lab ID:** 1108912-05  
**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<b>DIESEL RANGE ORGANICS BY GC-FID</b>						
<b>DRO (C10-C28)</b>	<b>420</b>		<b>SW8015M</b>		Prep Date: <b>8/31/2011</b>	Analyst: <b>RM</b>
<i>Surr: 4-Terphenyl-d14</i>	<i>71.8</i>		<i>39-115</i>	<i>%REC</i>	<i>1</i>	<i>9/1/2011 03:18 PM</i>
<b>GASOLINE RANGE ORGANICS BY GC-FID</b>						
<b>GRO (C6-C10)</b>	<b>ND</b>		<b>SW8015</b>			Analyst: <b>RM</b>
<i>Surr: Toluene-d8</i>	<i>104</i>		<i>50-150</i>	<i>%REC</i>	<i>100</i>	<i>8/31/2011 11:35 PM</i>
<b>MERCURY BY CVAA</b>						
<b>Mercury</b>	<b>0.043</b>		<b>SW7471</b>		Prep Date: <b>8/31/2011</b>	Analyst: <b>LR</b>
			<b>0.021</b>	<b>mg/Kg-dry</b>	<b>1</b>	<b>9/1/2011 02:41 PM</b>
<b>METALS BY ICP-MS</b>						
<b>Arsenic</b>	<b>4.1</b>		<b>SW6020A</b>		Prep Date: <b>8/30/2011</b>	Analyst: <b>CES</b>
<b>Barium</b>	<b>1,000</b>		<b>1.7</b>	<b>mg/Kg-dry</b>	<b>4</b>	<b>8/31/2011 09:37 PM</b>
Cadmium	ND		<b>17</b>	<b>mg/Kg-dry</b>	<b>40</b>	<b>9/1/2011 01:29 PM</b>
<b>Chromium</b>	<b>30</b>		<b>0.66</b>	<b>mg/Kg-dry</b>	<b>4</b>	<b>8/31/2011 09:37 PM</b>
<b>Copper</b>	<b>19</b>		<b>1.7</b>	<b>mg/Kg-dry</b>	<b>4</b>	<b>8/31/2011 09:37 PM</b>
<b>Lead</b>	<b>14</b>		<b>1.7</b>	<b>mg/Kg-dry</b>	<b>4</b>	<b>8/31/2011 09:37 PM</b>
<b>Nickel</b>	<b>18</b>		<b>1.7</b>	<b>mg/Kg-dry</b>	<b>4</b>	<b>8/31/2011 09:37 PM</b>
Selenium	ND		<b>1.7</b>	<b>mg/Kg-dry</b>	<b>4</b>	<b>8/31/2011 09:37 PM</b>
Silver	ND		<b>1.7</b>	<b>mg/Kg-dry</b>	<b>4</b>	<b>8/31/2011 09:37 PM</b>
<b>Zinc</b>	<b>78</b>		<b>3.3</b>	<b>mg/Kg-dry</b>	<b>4</b>	<b>8/31/2011 09:37 PM</b>
<b>SUBCONTRACTED ANALYSES</b>						
<b>Subcontracted Analyses</b>	<b>Rcvd 9/6/11</b>		<b>SUBCONTRACT</b>			Analyst: <b>A&amp;LGL</b>
			<b>as noted</b>		<b>1</b>	<b>9/6/2011</b>
<b>SEMI-VOLATILE ORGANIC COMPOUNDS</b>						
<b>Acenaphthene</b>	<b>ND</b>		<b>SW8270</b>		Prep Date: <b>8/31/2011</b>	Analyst: <b>HL</b>
<b>Anthracene</b>	<b>ND</b>		<b>33</b>	<b>µg/Kg-dry</b>	<b>1</b>	<b>9/2/2011 02:10 AM</b>
<b>Benzo(a)anthracene</b>	<b>ND</b>		<b>33</b>	<b>µg/Kg-dry</b>	<b>1</b>	<b>9/2/2011 02:10 AM</b>
<b>Benzo(a)pyrene</b>	<b>ND</b>		<b>33</b>	<b>µg/Kg-dry</b>	<b>1</b>	<b>9/2/2011 02:10 AM</b>
<b>Benzo(b)fluoranthene</b>	<b>ND</b>		<b>33</b>	<b>µg/Kg-dry</b>	<b>1</b>	<b>9/2/2011 02:10 AM</b>
<b>Benzo(g,h,i)perylene</b>	<b>ND</b>		<b>33</b>	<b>µg/Kg-dry</b>	<b>1</b>	<b>9/2/2011 02:10 AM</b>
<b>Benzo(k)fluoranthene</b>	<b>ND</b>		<b>33</b>	<b>µg/Kg-dry</b>	<b>1</b>	<b>9/2/2011 02:10 AM</b>
<b>Chrysene</b>	<b>ND</b>		<b>33</b>	<b>µg/Kg-dry</b>	<b>1</b>	<b>9/2/2011 02:10 AM</b>
<b>Dibenzo(a,h)anthracene</b>	<b>ND</b>		<b>33</b>	<b>µg/Kg-dry</b>	<b>1</b>	<b>9/2/2011 02:10 AM</b>
<b>Fluoranthene</b>	<b>60</b>		<b>33</b>	<b>µg/Kg-dry</b>	<b>1</b>	<b>9/2/2011 02:10 AM</b>
<b>Fluorene</b>	<b>42</b>		<b>33</b>	<b>µg/Kg-dry</b>	<b>1</b>	<b>9/2/2011 02:10 AM</b>
<b>Indeno(1,2,3-cd)pyrene</b>	<b>ND</b>		<b>33</b>	<b>µg/Kg-dry</b>	<b>1</b>	<b>9/2/2011 02:10 AM</b>
<b>Naphthalene</b>	<b>140</b>		<b>33</b>	<b>µg/Kg-dry</b>	<b>1</b>	<b>9/2/2011 02:10 AM</b>
<b>Pyrene</b>	<b>63</b>		<b>33</b>	<b>µg/Kg-dry</b>	<b>1</b>	<b>9/2/2011 02:10 AM</b>
<i>Surr: 2,4,6-Tribromophenol</i>	<i>105</i>		<i>34-140</i>	<i>%REC</i>	<i>1</i>	<i>9/2/2011 02:10 AM</i>

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

# ALS Group USA, Corp

Date: 08-Sep-11

**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11  
**Sample ID:** TR 22-20-597 W. Wall  
**Collection Date:** 8/24/2011 01:50 PM

**Work Order:** 1108912  
**Lab ID:** 1108912-05  
**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<i>Surr: 2-Fluorobiphenyl</i>	92.0		12-100	%REC	1	9/2/2011 02:10 AM
<i>Surr: 2-Fluorophenol</i>	84.4		33-117	%REC	1	9/2/2011 02:10 AM
<i>Surr: 4-Terphenyl-d14</i>	103		25-137	%REC	1	9/2/2011 02:10 AM
<i>Surr: Nitrobenzene-d5</i>	79.7		37-107	%REC	1	9/2/2011 02:10 AM
<i>Surr: Phenol-d6</i>	84.8		40-106	%REC	1	9/2/2011 02:10 AM
<b>VOLATILE ORGANIC COMPOUNDS</b>			<b>SW8260</b>			Analyst: <b>BG</b>
Benzene	ND		110	µg/Kg-dry	100	9/1/2011 10:09 AM
Ethylbenzene	ND		110	µg/Kg-dry	100	9/1/2011 10:09 AM
<b>m,p-Xylene</b>	<b>120</b>		<b>110</b>	<b>µg/Kg-dry</b>	100	9/1/2011 10:09 AM
o-Xylene	ND		110	µg/Kg-dry	100	9/1/2011 10:09 AM
Toluene	ND		110	µg/Kg-dry	100	9/1/2011 10:09 AM
Xylenes, Total	ND		340	µg/Kg-dry	100	9/1/2011 10:09 AM
<i>Surr: 1,2-Dichloroethane-d4</i>	100		70-120	%REC	100	9/1/2011 10:09 AM
<i>Surr: 4-Bromofluorobenzene</i>	100		75-120	%REC	100	9/1/2011 10:09 AM
<i>Surr: Dibromofluoromethane</i>	96.2		85-115	%REC	100	9/1/2011 10:09 AM
<i>Surr: Toluene-d8</i>	99.3		85-115	%REC	100	9/1/2011 10:09 AM
<b>CHROMIUM, TRIVALENT</b>			<b>CALCULATION</b>			Analyst: <b>JJG</b>
<b>Chromium, Trivalent</b>	<b>31</b>			<b>mg/kg-dry</b>	1	9/2/2011 02:35 PM
<b>CHROMIUM, HEXAVALENT</b>			<b>SW7196A</b>		Prep Date: <b>9/1/2011</b>	Analyst: <b>MB</b>
Chromium, Hexavalent	ND		0.55	mg/Kg-dry	1	9/2/2011 01:00 PM
<b>MOISTURE</b>			<b>A2540 G</b>			Analyst: <b>CG</b>
<b>Moisture</b>	<b>12</b>		<b>0.050</b>	<b>% of sample</b>	1	8/30/2011 02:49 PM
<b>PH</b>			<b>SW9045D</b>			Analyst: <b>JJG</b>
<b>pH</b>	<b>8.65</b>	H		<b>s.u.</b>	1	8/30/2011 10:15 AM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

# ALS Group USA, Corp

Date: 08-Sep-11

Client: HRL Compliance Solutions

Project: Williams TR 22-20-597 Pad LOE 8/24/11

Work Order: 1108912

Sample ID: TR 22-20-597 E. Wall

Lab ID: 1108912-06

Collection Date: 8/24/2011 02:10 PM

Matrix: SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<b>DIESEL RANGE ORGANICS BY GC-FID</b>						
<b>DRO (C10-C28)</b>	<b>13</b>		<b>SW8015M</b>		Prep Date: <b>8/31/2011</b>	Analyst: <b>RM</b>
			<b>4.8</b>	<b>mg/Kg-dry</b>	<b>1</b>	9/1/2011 03:42 PM
Surr: 4-Terphenyl-d14	45.5		39-115	%REC	1	9/1/2011 03:42 PM
<b>GASOLINE RANGE ORGANICS BY GC-FID</b>						
<b>GRO (C6-C10)</b>	<b>ND</b>		<b>SW8015</b>			Analyst: <b>RM</b>
			<b>5.8</b>	<b>mg/Kg-dry</b>	<b>100</b>	9/1/2011 12:01 PM
Surr: Toluene-d8	103		50-150	%REC	100	9/1/2011 12:01 PM
<b>MERCURY BY CVAA</b>						
<b>Mercury</b>	<b>0.023</b>		<b>SW7471</b>		Prep Date: <b>8/31/2011</b>	Analyst: <b>LR</b>
			<b>0.021</b>	<b>mg/Kg-dry</b>	<b>1</b>	9/1/2011 02:43 PM
<b>METALS BY ICP-MS</b>						
<b>Arsenic</b>	<b>2.8</b>		<b>SW6020A</b>		Prep Date: <b>8/30/2011</b>	Analyst: <b>CES</b>
			<b>1.8</b>	<b>mg/Kg-dry</b>	<b>4</b>	8/31/2011 10:04 PM
<b>Barium</b>	<b>380</b>		<b>1.8</b>	<b>mg/Kg-dry</b>	<b>4</b>	8/31/2011 10:04 PM
Cadmium	ND		0.73	mg/Kg-dry	4	8/31/2011 10:04 PM
<b>Chromium</b>	<b>38</b>		<b>1.8</b>	<b>mg/Kg-dry</b>	<b>4</b>	8/31/2011 10:04 PM
<b>Copper</b>	<b>13</b>		<b>1.8</b>	<b>mg/Kg-dry</b>	<b>4</b>	8/31/2011 10:04 PM
<b>Lead</b>	<b>17</b>		<b>1.8</b>	<b>mg/Kg-dry</b>	<b>4</b>	8/31/2011 10:04 PM
<b>Nickel</b>	<b>18</b>		<b>1.8</b>	<b>mg/Kg-dry</b>	<b>4</b>	8/31/2011 10:04 PM
Selenium	ND		1.8	mg/Kg-dry	4	8/31/2011 10:04 PM
Silver	ND		1.8	mg/Kg-dry	4	8/31/2011 10:04 PM
<b>Zinc</b>	<b>64</b>		<b>3.7</b>	<b>mg/Kg-dry</b>	<b>4</b>	8/31/2011 10:04 PM
<b>SUBCONTRACTED ANALYSES</b>						
<b>Subcontracted Analyses</b>	<b>Rcvd 9/6/11</b>		<b>SUBCONTRACT</b>			Analyst: <b>A&amp;LGL</b>
			<b>as noted</b>		<b>1</b>	9/6/2011
<b>SEMI-VOLATILE ORGANIC COMPOUNDS</b>						
<b>Acenaphthene</b>	<b>ND</b>		<b>SW8270</b>		Prep Date: <b>9/6/2011</b>	Analyst: <b>HL</b>
			<b>34</b>	<b>µg/Kg-dry</b>	<b>1</b>	9/8/2011 02:36 AM
<b>Anthracene</b>	<b>ND</b>		<b>34</b>	<b>µg/Kg-dry</b>	<b>1</b>	9/8/2011 02:36 AM
<b>Benzo(a)anthracene</b>	<b>ND</b>		<b>34</b>	<b>µg/Kg-dry</b>	<b>1</b>	9/8/2011 02:36 AM
<b>Benzo(a)pyrene</b>	<b>ND</b>		<b>34</b>	<b>µg/Kg-dry</b>	<b>1</b>	9/8/2011 02:36 AM
<b>Benzo(b)fluoranthene</b>	<b>ND</b>		<b>34</b>	<b>µg/Kg-dry</b>	<b>1</b>	9/8/2011 02:36 AM
<b>Benzo(g,h,i)perylene</b>	<b>ND</b>		<b>34</b>	<b>µg/Kg-dry</b>	<b>1</b>	9/8/2011 02:36 AM
<b>Benzo(k)fluoranthene</b>	<b>ND</b>		<b>34</b>	<b>µg/Kg-dry</b>	<b>1</b>	9/8/2011 02:36 AM
<b>Chrysene</b>	<b>ND</b>		<b>34</b>	<b>µg/Kg-dry</b>	<b>1</b>	9/8/2011 02:36 AM
<b>Dibenzo(a,h)anthracene</b>	<b>ND</b>		<b>34</b>	<b>µg/Kg-dry</b>	<b>1</b>	9/8/2011 02:36 AM
<b>Fluoranthene</b>	<b>ND</b>		<b>34</b>	<b>µg/Kg-dry</b>	<b>1</b>	9/8/2011 02:36 AM
<b>Fluorene</b>	<b>ND</b>		<b>34</b>	<b>µg/Kg-dry</b>	<b>1</b>	9/8/2011 02:36 AM
<b>Indeno(1,2,3-cd)pyrene</b>	<b>ND</b>		<b>34</b>	<b>µg/Kg-dry</b>	<b>1</b>	9/8/2011 02:36 AM
<b>Naphthalene</b>	<b>ND</b>		<b>34</b>	<b>µg/Kg-dry</b>	<b>1</b>	9/8/2011 02:36 AM
<b>Pyrene</b>	<b>ND</b>		<b>34</b>	<b>µg/Kg-dry</b>	<b>1</b>	9/8/2011 02:36 AM
Surr: 2,4,6-Tribromophenol	99.8		34-140	%REC	1	9/8/2011 02:36 AM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

# ALS Group USA, Corp

Date: 08-Sep-11

**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11  
**Sample ID:** TR 22-20-597 E. Wall  
**Collection Date:** 8/24/2011 02:10 PM

**Work Order:** 1108912  
**Lab ID:** 1108912-06  
**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<i>Surr: 2-Fluorobiphenyl</i>	66.8		12-100	%REC	1	9/8/2011 02:36 AM
<i>Surr: 2-Fluorophenol</i>	83.0		33-117	%REC	1	9/8/2011 02:36 AM
<i>Surr: 4-Terphenyl-d14</i>	96.1		25-137	%REC	1	9/8/2011 02:36 AM
<i>Surr: Nitrobenzene-d5</i>	70.2		37-107	%REC	1	9/8/2011 02:36 AM
<i>Surr: Phenol-d6</i>	78.4		40-106	%REC	1	9/8/2011 02:36 AM
<b>VOLATILE ORGANIC COMPOUNDS</b>			<b>SW8260</b>			Analyst: <b>BG</b>
Benzene	ND		120	µg/Kg-dry	100	9/1/2011 10:34 AM
Ethylbenzene	ND		120	µg/Kg-dry	100	9/1/2011 10:34 AM
<b>m,p-Xylene</b>	<b>130</b>		<b>120</b>	<b>µg/Kg-dry</b>	100	9/1/2011 10:34 AM
o-Xylene	ND		120	µg/Kg-dry	100	9/1/2011 10:34 AM
Toluene	ND		120	µg/Kg-dry	100	9/1/2011 10:34 AM
Xylenes, Total	ND		350	µg/Kg-dry	100	9/1/2011 10:34 AM
<i>Surr: 1,2-Dichloroethane-d4</i>	98.0		70-120	%REC	100	9/1/2011 10:34 AM
<i>Surr: 4-Bromofluorobenzene</i>	99.1		75-120	%REC	100	9/1/2011 10:34 AM
<i>Surr: Dibromofluoromethane</i>	94.3		85-115	%REC	100	9/1/2011 10:34 AM
<i>Surr: Toluene-d8</i>	99.0		85-115	%REC	100	9/1/2011 10:34 AM
<b>CHROMIUM, TRIVALENT</b>			<b>CALCULATION</b>			Analyst: <b>JJG</b>
<b>Chromium, Trivalent</b>	<b>37</b>			<b>mg/kg-dry</b>	1	9/2/2011 02:35 PM
<b>CHROMIUM, HEXAVALENT</b>			<b>SW7196A</b>		Prep Date: <b>9/1/2011</b>	Analyst: <b>MB</b>
Chromium, Hexavalent	ND		0.58	mg/Kg-dry	1	9/2/2011 01:00 PM
<b>MOISTURE</b>			<b>A2540 G</b>			Analyst: <b>CG</b>
<b>Moisture</b>	<b>15</b>		<b>0.050</b>	<b>% of sample</b>	1	8/30/2011 02:49 PM
<b>PH</b>			<b>SW9045D</b>			Analyst: <b>JJG</b>
<b>pH</b>	<b>7.75</b>	H		<b>s.u.</b>	1	8/30/2011 10:15 AM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

Report Number: F11244-0176

Account Number: 91000

# A & L GREAT LAKES LABORATORIES, INC.

3505 Conestoga Drive • Fort Wayne, Indiana 46808-4413 • Phone 260-483-4759 • Fax 260-483-5274

www.algreatlakes.com • lab@algreatlakes.com



**QUALITY ANALYSES FOR INFORMED DECISIONS**

TO: ALS LABORATORY GROUP  
3352 128TH AVE  
HOLLAND, MI 49424-9263

RE: 1108912

DATE RECEIVED: 09/01/2011

DATE REPORTED: 09/06/2011

PAGE: 1

P.O. NUMBER: 20-122010668

ATTN: ANN PRESTON

## REPORT OF ANALYSIS

LAB NO.	SAMPLE ID	ANALYSIS	RESULT	UNIT	METHOD
82829	01C	Sat'd Paste Extraction with DIW	1		USDA Handbook 60
		Conductivity (ECe)	0.36	mmho/cm	USDA Handbook 60
		Calcium (Sat'd Paste)	69	ppm	USDA Handbook 60
		Magnesium (Sat'd Paste)	15	ppm	USDA Handbook 60
		Sodium (Sat'd Paste)	24	ppm	USDA Handbook 60
		Sodium Adsorption Ratio	0.7	-	USDA Handbook 60
82830	02C	Sat'd Paste Extraction with DIW	1		USDA Handbook 60
		Conductivity (ECe)	0.81	mmho/cm	USDA Handbook 60
		Calcium (Sat'd Paste)	113	ppm	USDA Handbook 60
		Magnesium (Sat'd Paste)	21	ppm	USDA Handbook 60
		Sodium (Sat'd Paste)	441	ppm	USDA Handbook 60
		Sodium Adsorption Ratio	10.0	-	USDA Handbook 60
82831	03C	Sat'd Paste Extraction with DIW	1		USDA Handbook 60
		Conductivity (ECe)	4.21	mmho/cm	USDA Handbook 60
		Calcium (Sat'd Paste)	281	ppm	USDA Handbook 60
		Magnesium (Sat'd Paste)	35	ppm	USDA Handbook 60
		Sodium (Sat'd Paste)	4811	ppm	USDA Handbook 60
		Sodium Adsorption Ratio	71.2	-	USDA Handbook 60
82832	04C	Sat'd Paste Extraction with DIW	1		USDA Handbook 60
		Conductivity (ECe)	3.63	mmho/cm	USDA Handbook 60
		Calcium (Sat'd Paste)	254	ppm	USDA Handbook 60
		Magnesium (Sat'd Paste)	34	ppm	USDA Handbook 60
		Sodium (Sat'd Paste)	4024	ppm	USDA Handbook 60
		Sodium Adsorption Ratio	62.8	-	USDA Handbook 60



Report Number: F11244-0176

Account Number: 91000

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**QUALITY ANALYSES FOR INFORMED DECISIONS**

TO: ALS LABORATORY GROUP  
3352 128TH AVE  
HOLLAND, MI 49424-9263

RE: 1108912

DATE RECEIVED: 09/01/2011

DATE REPORTED: 09/06/2011

PAGE: 2

P.O. NUMBER: 20-122010668

ATTN: ANN PRESTON

## REPORT OF ANALYSIS

LAB NO.	SAMPLE ID	ANALYSIS	RESULT	UNIT	METHOD
82833	05C	Sat'd Paste Extraction with DIW	1		USDA Handbook 60
		Conductivity (ECe)	3.33	mmho/cm	USDA Handbook 60
		Calcium (Sat'd Paste)	503	ppm	USDA Handbook 60
		Magnesium (Sat'd Paste)	46	ppm	USDA Handbook 60
		Sodium (Sat'd Paste)	3740	ppm	USDA Handbook 60
		Sodium Adsorption Ratio	42.7	-	USDA Handbook 60
82834	06C	Sat'd Paste Extraction with DIW	1		USDA Handbook 60
		Conductivity (ECe)	0.83	mmho/cm	USDA Handbook 60
		Calcium (Sat'd Paste)	158	ppm	USDA Handbook 60
		Magnesium (Sat'd Paste)	30	ppm	USDA Handbook 60
		Sodium (Sat'd Paste)	144	ppm	USDA Handbook 60
		Sodium Adsorption Ratio	2.8	-	USDA Handbook 60

# ALS Group USA, Corp

Date: 08-Sep-11

**Client:** HRL Compliance Solutions

## QC BATCH REPORT

**Work Order:** 1108912

**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

Batch ID: **35214**

Instrument ID **GC8**

Method: **SW8015M**

<b>MBLK</b>	Sample ID: <b>DBLKS1-35214-35214</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>8/31/2011 12:52 PM</b>			
Client ID:	Run ID: <b>GC8_110830A</b>				SeqNo: <b>1722245</b>		Prep Date: <b>8/30/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	ND	4.2								
<i>Surr: 4-Terphenyl-d14</i>	1.312	0	1.667	0	78.7	39-115	0			

<b>LCS</b>	Sample ID: <b>DLCSS1-35214-35214</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>8/31/2011 02:01 PM</b>			
Client ID:	Run ID: <b>GC8_110830A</b>				SeqNo: <b>1723379</b>		Prep Date: <b>8/30/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	150.4	4.2	166.7	0	90.2	60-130	0			
<i>Surr: 4-Terphenyl-d14</i>	1.083	0	1.667	0	65	39-115	0			

<b>LCSD</b>	Sample ID: <b>DLCSDS1-35214-35214</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>8/31/2011 02:01 PM</b>			
Client ID:	Run ID: <b>GC8_110830A</b>				SeqNo: <b>1723419</b>		Prep Date: <b>8/30/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	175.6	4.2	166.7	0	105	60-130	150.4	15.5	30	
<i>Surr: 4-Terphenyl-d14</i>	1.288	0	1.667	0	77.3	39-115	1.083	17.3	30	

<b>MS</b>	Sample ID: <b>1108889-20B MS</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>8/31/2011 02:24 PM</b>			
Client ID:	Run ID: <b>GC8_110830A</b>				SeqNo: <b>1723380</b>		Prep Date: <b>8/30/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	276.9	7.9	316.4	4.124	86.2	60-130	0			
<i>Surr: 4-Terphenyl-d14</i>	1.939	0	3.164	0	61.3	39-115	0			

<b>MSD</b>	Sample ID: <b>1108889-20B MSD</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>8/31/2011 02:24 PM</b>			
Client ID:	Run ID: <b>GC8_110830A</b>				SeqNo: <b>1723420</b>		Prep Date: <b>8/30/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	226.2	8.1	322.4	4.124	68.9	60-130	276.9	20.2	30	
<i>Surr: 4-Terphenyl-d14</i>	1.819	0	3.224	0	56.4	39-115	1.939	6.35	30	

The following samples were analyzed in this batch:

1108912-01B	1108912-02B	1108912-03B
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**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35236**      Instrument ID **GC8**      Method: **SW8015M**

<b>MBLK</b>	Sample ID: <b>DBLKS1-35236-35236</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/1/2011 02:55 PM</b>			
Client ID:	Run ID: <b>GC8_110901A</b>				SeqNo: <b>1723709</b>		Prep Date: <b>8/31/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	ND	4.2								
<i>Surr: 4-Terphenyl-d14</i>	1.36	0	1.667	0	81.6	39-115	0			

<b>LCS</b>	Sample ID: <b>DLCSS1-35236-35236</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/1/2011 12:59 PM</b>			
Client ID:	Run ID: <b>GC8_110901A</b>				SeqNo: <b>1724697</b>		Prep Date: <b>8/31/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	158.1	4.2	166.7	0	94.9	60-130	0			
<i>Surr: 4-Terphenyl-d14</i>	1.182	0	1.667	0	70.9	39-115	0			

<b>LCSD</b>	Sample ID: <b>DLCSDS1-35236-35236</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/1/2011 01:23 PM</b>			
Client ID:	Run ID: <b>GC8_110901A</b>				SeqNo: <b>1724676</b>		Prep Date: <b>8/31/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	168	4.2	166.7	0	101	60-130	158.1	6.05	30	
<i>Surr: 4-Terphenyl-d14</i>	1.06	0	1.667	0	63.6	39-115	1.182	10.8	30	

<b>MS</b>	Sample ID: <b>1108926-27B MS</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/1/2011 01:23 PM</b>			
Client ID:	Run ID: <b>GC8_110901A</b>				SeqNo: <b>1724698</b>		Prep Date: <b>8/31/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	287.3	8.1	324.7	3.173	87.5	60-130	0			
<i>Surr: 4-Terphenyl-d14</i>	2.023	0	3.247	0	62.3	39-115	0			

<b>MSD</b>	Sample ID: <b>1108926-27B MSD</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/1/2011 01:46 PM</b>			
Client ID:	Run ID: <b>GC8_110901A</b>				SeqNo: <b>1724677</b>		Prep Date: <b>8/31/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	356.8	8.3	331.9	3.173	107	60-130	287.3	21.6	30	
<i>Surr: 4-Terphenyl-d14</i>	2.177	0	3.319	0	65.6	39-115	2.023	7.32	30	

The following samples were analyzed in this batch:      1108912-04B      1108912-05B      1108912-06B

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **R94173**      Instrument ID **GC9**      Method: **SW8015**

<b>MBLK</b>	Sample ID: <b>MBLK-R94173-R94173</b>				Units: <b>µg/L</b>		Analysis Date: <b>8/31/2011 03:19 PM</b>			
Client ID:	Run ID: <b>GC9_110831A</b>				SeqNo: <b>1723625</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
GRO (C6-C10)	ND	200								
<i>Surr: Toluene-d8</i>	<i>110.1</i>	<i>0</i>	<i>100</i>	<i>0</i>	<i>110</i>	<i>70-130</i>	<i>0</i>			

<b>LCS</b>	Sample ID: <b>LCS-R94173-R94173</b>				Units: <b>µg/L</b>		Analysis Date: <b>8/31/2011 02:01 PM</b>			
Client ID:	Run ID: <b>GC9_110831A</b>				SeqNo: <b>1723623</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
GRO (C6-C10)	26600	200	25000	0	106	70-130	0			
<i>Surr: Toluene-d8</i>	<i>111.9</i>	<i>0</i>	<i>100</i>	<i>0</i>	<i>112</i>	<i>70-130</i>	<i>0</i>			

<b>LCSD</b>	Sample ID: <b>LCSD-R94173-R94173</b>				Units: <b>µg/L</b>		Analysis Date: <b>8/31/2011 02:27 PM</b>			
Client ID:	Run ID: <b>GC9_110831A</b>				SeqNo: <b>1723624</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
GRO (C6-C10)	25570	200	25000	0	102	70-130	26600	3.94	30	
<i>Surr: Toluene-d8</i>	<i>105.8</i>	<i>0</i>	<i>100</i>	<i>0</i>	<i>106</i>	<i>70-130</i>	<i>111.9</i>	<i>5.56</i>	<i>30</i>	

<b>MS</b>	Sample ID: <b>1108889-20A MS</b>				Units: <b>µg/Kg</b>		Analysis Date: <b>9/1/2011 01:18 AM</b>			
Client ID:	Run ID: <b>GC9_110831A</b>				SeqNo: <b>1723645</b>		Prep Date:		DF: <b>50</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
GRO (C6-C10)	1300000	2,500	1250000	0	104	70-130	0			
<i>Surr: Toluene-d8</i>	<i>5097</i>	<i>0</i>	<i>5000</i>	<i>0</i>	<i>102</i>	<i>50-150</i>	<i>0</i>			

<b>MS</b>	Sample ID: <b>1108889-06A MS</b>				Units: <b>µg/Kg</b>		Analysis Date: <b>9/1/2011 12:27 PM</b>			
Client ID:	Run ID: <b>GC9_110831A</b>				SeqNo: <b>1723648</b>		Prep Date:		DF: <b>50</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
GRO (C6-C10)	1367000	2,500	1250000	0	109	70-130	0			
<i>Surr: Toluene-d8</i>	<i>5460</i>	<i>0</i>	<i>5000</i>	<i>0</i>	<i>109</i>	<i>50-150</i>	<i>0</i>			

<b>MSD</b>	Sample ID: <b>1108889-20A MSD</b>				Units: <b>µg/Kg</b>		Analysis Date: <b>9/1/2011 01:44 AM</b>			
Client ID:	Run ID: <b>GC9_110831A</b>				SeqNo: <b>1723646</b>		Prep Date:		DF: <b>50</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
GRO (C6-C10)	1259000	2,500	1250000	0	101	70-130	1300000	3.16	30	
<i>Surr: Toluene-d8</i>	<i>5130</i>	<i>0</i>	<i>5000</i>	<i>0</i>	<i>103</i>	<i>50-150</i>	<i>5097</i>	<i>0.645</i>	<i>30</i>	

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **R94173** Instrument ID **GC9** Method: **SW8015**

**MSD** Sample ID: **1108889-06A MSD** Units: **µg/Kg** Analysis Date: **9/1/2011 12:53 PM**

Client ID: Run ID: **GC9\_110831A** SeqNo: **1723649** Prep Date: DF: **50**

Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
GRO (C6-C10)	1321000	2,500	1250000	0	106	70-130	1367000	3.38	30	
<i>Surr: Toluene-d8</i>	<i>5160</i>	<i>0</i>	<i>5000</i>	<i>0</i>	<i>103</i>	<i>50-150</i>	<i>5460</i>	<i>5.64</i>	<i>30</i>	

The following samples were analyzed in this batch:

1108912-01A	1108912-02A	1108912-03A
1108912-04A	1108912-05A	1108912-06A

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35252** Instrument ID **HG1** Method: **SW7471**

<b>MBLK</b>	Sample ID: <b>MBLK-35252-35252</b>				Units: <b>mg/Kg</b>			Analysis Date: <b>9/1/2011 01:49 PM</b>		
Client ID:	Run ID: <b>HG1_110901A</b>				SeqNo: <b>1723335</b>			Prep Date: <b>8/31/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Mercury	ND	0.020								

<b>LCS</b>	Sample ID: <b>LCS-35252-35252</b>				Units: <b>mg/Kg</b>			Analysis Date: <b>9/1/2011 01:51 PM</b>		
Client ID:	Run ID: <b>HG1_110901A</b>				SeqNo: <b>1723336</b>			Prep Date: <b>8/31/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Mercury	0.1721	0.020	0.1665		0	103	80-120	0		

<b>LCSD</b>	Sample ID: <b>LCSD-35252-35252</b>				Units: <b>mg/Kg</b>			Analysis Date: <b>9/1/2011 01:53 PM</b>		
Client ID:	Run ID: <b>HG1_110901A</b>				SeqNo: <b>1723337</b>			Prep Date: <b>8/31/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Mercury	0.1744	0.020	0.1665		0	105	80-120	0.1721	1.35	20

<b>MS</b>	Sample ID: <b>1108867-03A MS</b>				Units: <b>mg/Kg</b>			Analysis Date: <b>9/1/2011 02:00 PM</b>		
Client ID:	Run ID: <b>HG1_110901A</b>				SeqNo: <b>1723340</b>			Prep Date: <b>8/31/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Mercury	0.1475	0.017	0.1415	0.0005556	104	75-125		0		

<b>MSD</b>	Sample ID: <b>1108867-03A MSD</b>				Units: <b>mg/Kg</b>			Analysis Date: <b>9/1/2011 02:02 PM</b>		
Client ID:	Run ID: <b>HG1_110901A</b>				SeqNo: <b>1723341</b>			Prep Date: <b>8/31/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Mercury	0.1574	0.018	0.1507	0.0005556	104	75-125	0.1475	6.52	35	

The following samples were analyzed in this batch:

1108912-01B	1108912-02B	1108912-03B
1108912-04B	1108912-05B	1108912-06B

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35228** Instrument ID **ICPMS1** Method: **SW6020A**

<b>MBLK</b>	Sample ID: <b>MBLK-35228-35228</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>8/31/2011 06:11 PM</b>			
Client ID:	Run ID: <b>ICPMS1_110831A</b>				SeqNo: <b>1722949</b>		Prep Date: <b>8/30/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	ND	0.25								
Barium	ND	0.25								
Cadmium	0.001396	0.10								J
Chromium	0.004901	0.25								J
Copper	ND	0.25								
Lead	ND	0.25								
Nickel	ND	0.25								
Selenium	ND	0.25								
Silver	ND	0.25								
Zinc	0.04614	0.50								J

<b>LCS</b>	Sample ID: <b>LCS-35228-35228</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>8/31/2011 06:16 PM</b>			
Client ID:	Run ID: <b>ICPMS1_110831A</b>				SeqNo: <b>1722950</b>		Prep Date: <b>8/30/2011</b>		DF: <b>2</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	4.609	0.50	5	0	92.2	80-120	0			
Barium	4.557	0.50	5	0	91.1	80-120	0			
Cadmium	4.603	0.20	5	0	92.1	80-120	0			
Chromium	4.679	0.50	5	0	93.6	80-120	0			
Copper	4.727	0.50	5	0	94.5	80-120	0			
Lead	4.696	0.50	5	0	93.9	80-120	0			
Nickel	4.736	0.50	5	0	94.7	80-120	0			
Selenium	4.543	0.50	5	0	90.9	80-120	0			
Silver	4.487	0.50	5	0	89.7	80-120	0			
Zinc	4.629	1.0	5	0	92.6	80-120	0			

<b>LCSD</b>	Sample ID: <b>LCSD-35228-35228</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>8/31/2011 06:21 PM</b>			
Client ID:	Run ID: <b>ICPMS1_110831A</b>				SeqNo: <b>1722951</b>		Prep Date: <b>8/30/2011</b>		DF: <b>2</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	4.403	0.50	5	0	88.1	80-120	4.609	4.57	20	
Barium	4.427	0.50	5	0	88.5	80-120	4.557	2.89	20	
Cadmium	4.452	0.20	5	0	89	80-120	4.603	3.34	20	
Chromium	4.506	0.50	5	0	90.1	80-120	4.679	3.77	20	
Copper	4.482	0.50	5	0	89.6	80-120	4.727	5.32	20	
Lead	4.481	0.50	5	0	89.6	80-120	4.696	4.69	20	
Nickel	4.513	0.50	5	0	90.3	80-120	4.736	4.82	20	
Selenium	4.486	0.50	5	0	89.7	80-120	4.543	1.26	20	
Silver	4.313	0.50	5	0	86.3	80-120	4.487	3.95	20	
Zinc	4.459	1.0	5	0	89.2	80-120	4.629	3.74	20	

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35228**      Instrument ID **ICPMS1**      Method: **SW6020A**

MS				Sample ID: <b>1108912-03BMS</b>			Units: <b>mg/Kg</b>		Analysis Date: <b>8/31/2011 07:30 PM</b>	
Client ID: <b>TR 22-20-597 S Bottom</b>				Run ID: <b>ICPMS1_110831A</b>			SeqNo: <b>1722962</b>		Prep Date: <b>8/30/2011</b>	
									DF: <b>4</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	8.025	1.4	6.84	1.905	89.5	80-120	0			
Barium	302.3	1.4	6.84	252.2	734	80-120	0			SO
Cadmium	7.111	0.55	6.84	0.6827	94	80-120	0			
Chromium	19.5	1.4	6.84	14.36	75.2	80-120	0			S
Copper	12.23	1.4	6.84	6.889	78.1	80-120	0			S
Nickel	14.9	1.4	6.84	8.84	88.6	80-120	0			
Selenium	6.865	1.4	6.84	0.8821	87.5	80-120	0			
Silver	5.261	1.4	6.84	0.04208	76.3	80-120	0			S
Zinc	51.6	2.7	6.84	41.83	143	80-120	0			SO

MS				Sample ID: <b>1108912-03BMS</b>			Units: <b>mg/Kg</b>		Analysis Date: <b>9/1/2011 01:08 PM</b>	
Client ID: <b>TR 22-20-597 S Bottom</b>				Run ID: <b>ICPMS1_110901A</b>			SeqNo: <b>1723489</b>		Prep Date: <b>8/30/2011</b>	
									DF: <b>4</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Lead	21.47	1.4	6.84	11.44	147	80-120	0			S

MSD				Sample ID: <b>1108912-03BMSD</b>			Units: <b>mg/Kg</b>		Analysis Date: <b>8/31/2011 07:35 PM</b>	
Client ID: <b>TR 22-20-597 S Bottom</b>				Run ID: <b>ICPMS1_110831A</b>			SeqNo: <b>1722963</b>		Prep Date: <b>8/30/2011</b>	
									DF: <b>4</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	8.815	1.5	7.692	1.905	89.8	80-120	8.025	9.39	25	
Barium	312.6	1.5	7.692	252.2	786	80-120	302.3	3.35	25	SO
Cadmium	7.6	0.62	7.692	0.6827	89.9	80-120	7.111	6.65	25	
Chromium	22.05	1.5	7.692	14.36	100	80-120	19.5	12.3	25	
Copper	13.9	1.5	7.692	6.889	91.2	80-120	12.23	12.8	25	
Nickel	16.44	1.5	7.692	8.84	98.8	80-120	14.9	9.79	25	
Selenium	7.225	1.5	7.692	0.8821	82.5	80-120	6.865	5.11	25	
Silver	6.138	1.5	7.692	0.04208	79.3	80-120	5.261	15.4	25	S
Zinc	57.14	3.1	7.692	41.83	199	80-120	51.6	10.2	25	SO

MSD				Sample ID: <b>1108912-03BMSD</b>			Units: <b>mg/Kg</b>		Analysis Date: <b>9/1/2011 01:14 PM</b>	
Client ID: <b>TR 22-20-597 S Bottom</b>				Run ID: <b>ICPMS1_110901A</b>			SeqNo: <b>1723491</b>		Prep Date: <b>8/30/2011</b>	
									DF: <b>4</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Lead	22.34	1.5	7.692	11.44	142	80-120	21.47	3.96	25	S

The following samples were analyzed in this batch:

1108912-01B	1108912-02B	1108912-03B
1108912-04B	1108912-05B	1108912-06B

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.



**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35213**      Instrument ID **SVMS4**      Method: **SW8270**

MBLK		Sample ID: <b>SBLKS1-35213-35213</b>				Units: <b>µg/Kg</b>		Analysis Date: <b>8/31/2011 10:59 AM</b>		
Client ID:		Run ID: <b>SVMS4_110831A</b>				SeqNo: <b>1721969</b>		Prep Date: <b>8/30/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Acenaphthene	ND	30								
Anthracene	ND	30								
Benzo(a)anthracene	ND	30								
Benzo(a)pyrene	ND	30								
Benzo(b)fluoranthene	ND	30								
Benzo(g,h,i)perylene	ND	30								
Benzo(k)fluoranthene	ND	30								
Chrysene	ND	30								
Dibenzo(a,h)anthracene	ND	30								
Fluoranthene	ND	30								
Fluorene	ND	30								
Indeno(1,2,3-cd)pyrene	ND	30								
Naphthalene	ND	30								
Pyrene	ND	30								
<i>Surr: 2,4,6-Tribromophenol</i>	1260	0	1667	0	75.6	34-140		0		
<i>Surr: 2-Fluorobiphenyl</i>	1153	0	1667	0	69.2	12-100		0		
<i>Surr: 2-Fluorophenol</i>	1222	0	1667	0	73.3	33-117		0		
<i>Surr: 4-Terphenyl-d14</i>	1431	0	1667	0	85.9	25-137		0		
<i>Surr: Nitrobenzene-d5</i>	1131	0	1667	0	67.8	37-107		0		
<i>Surr: Phenol-d6</i>	1236	0	1667	0	74.1	40-106		0		

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35213**      Instrument ID **SVMS4**      Method: **SW8270**

LCS		Sample ID: <b>SLCSS1-35213-35213</b>				Units: <b>µg/Kg</b>		Analysis Date: <b>8/31/2011 09:57 AM</b>		
Client ID:		Run ID: <b>SVMS4_110831A</b>				SeqNo: <b>1722118</b>		Prep Date: <b>8/30/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Acenaphthene	1153	30	1333	0	86.5	45-110	0			
Anthracene	1336	30	1333	0	100	55-105	0			
Benzo(a)anthracene	1278	30	1333	0	95.9	50-110	0			
Benzo(a)pyrene	1194	30	1333	0	89.6	50-110	0			
Benzo(b)fluoranthene	1253	30	1333	0	94	45-115	0			
Benzo(g,h,i)perylene	1199	30	1333	0	89.9	40-125	0			
Benzo(k)fluoranthene	1264	30	1333	0	94.8	45-115	0			
Chrysene	1368	30	1333	0	103	55-110	0			
Dibenzo(a,h)anthracene	1226	30	1333	0	92	40-125	0			
Fluoranthene	1239	30	1333	0	92.9	55-115	0			
Fluorene	1256	30	1333	0	94.2	50-110	0			
Indeno(1,2,3-cd)pyrene	1213	30	1333	0	91	40-120	0			
Naphthalene	924.3	30	1333	0	69.3	40-105	0			
Pyrene	1201	30	1333	0	90.1	45-125	0			
<i>Surr: 2,4,6-Tribromophenol</i>	1522	0	1667	0	91.3	34-140	0			
<i>Surr: 2-Fluorobiphenyl</i>	1238	0	1667	0	74.3	12-100	0			
<i>Surr: 2-Fluorophenol</i>	1112	0	1667	0	66.7	33-117	0			
<i>Surr: 4-Terphenyl-d14</i>	1590	0	1667	0	95.4	25-137	0			
<i>Surr: Nitrobenzene-d5</i>	1163	0	1667	0	69.8	37-107	0			
<i>Surr: Phenol-d6</i>	1197	0	1667	0	71.8	40-106	0			

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35213**      Instrument ID **SVMS4**      Method: **SW8270**

LCSD		Sample ID: <b>SLCSDS1-35213-35213</b>				Units: <b>µg/Kg</b>		Analysis Date: <b>8/31/2011 10:28 AM</b>		
Client ID:		Run ID: <b>SVMS4_110831A</b>				SeqNo: <b>1722119</b>		Prep Date: <b>8/30/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Acenaphthene	1259	30	1333	0	94.5	45-110	1153	8.84	25	
Anthracene	1367	30	1333	0	103	55-105	1336	2.32	25	
Benzo(a)anthracene	1307	30	1333	0	98	50-110	1278	2.24	25	
Benzo(a)pyrene	1216	30	1333	0	91.2	50-110	1194	1.83	25	
Benzo(b)fluoranthene	1174	30	1333	0	88.1	45-115	1253	6.51	25	
Benzo(g,h,i)perylene	1224	30	1333	0	91.8	40-125	1199	2.06	25	
Benzo(k)fluoranthene	1240	30	1333	0	93	45-115	1264	1.89	25	
Chrysene	1383	30	1333	0	104	55-110	1368	1.11	25	
Dibenzo(a,h)anthracene	1266	30	1333	0	95	40-125	1226	3.18	25	
Fluoranthene	1274	30	1333	0	95.5	55-115	1239	2.79	25	
Fluorene	1302	30	1333	0	97.7	50-110	1256	3.62	25	
Indeno(1,2,3-cd)pyrene	1253	30	1333	0	94	40-120	1213	3.24	25	
Naphthalene	1126	30	1333	0	84.5	40-105	924.3	19.7	25	
Pyrene	1262	30	1333	0	94.7	45-125	1201	4.95	25	
<i>Surr: 2,4,6-Tribromophenol</i>	1613	0	1667	0	96.8	34-140	1522	5.76	40	
<i>Surr: 2-Fluorobiphenyl</i>	1476	0	1667	0	88.5	12-100	1238	17.5	40	
<i>Surr: 2-Fluorophenol</i>	1384	0	1667	0	83.1	33-117	1112	21.8	40	
<i>Surr: 4-Terphenyl-d14</i>	1644	0	1667	0	98.6	25-137	1590	3.36	40	
<i>Surr: Nitrobenzene-d5</i>	1422	0	1667	0	85.3	37-107	1163	20	40	
<i>Surr: Phenol-d6</i>	1483	0	1667	0	89	40-106	1197	21.3	40	

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35213**      Instrument ID **SVMS4**      Method: **SW8270**

MS				Sample ID: <b>1108889-20B MS</b>			Units: <b>µg/Kg</b>		Analysis Date: <b>8/31/2011 11:30 AM</b>	
Client ID:				Run ID: <b>SVMS4_110831A</b>			SeqNo: <b>1722120</b>		Prep Date: <b>8/30/2011</b>	
							DF: <b>1</b>			
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Acenaphthene	1870	57	2546	0	73.5	45-110	0			
Anthracene	2013	57	2546	0	79.1	55-105	0			
Benzo(a)anthracene	1978	57	2546	0	77.7	50-110	0			
Benzo(a)pyrene	1713	57	2546	0	67.3	50-110	0			
Benzo(b)fluoranthene	1841	57	2546	0	72.3	45-115	0			
Benzo(g,h,i)perylene	1757	57	2546	0	69	40-125	0			
Benzo(k)fluoranthene	1815	57	2546	0	71.3	45-115	0			
Chrysene	2037	57	2546	0	80	55-110	0			
Dibenzo(a,h)anthracene	1818	57	2546	0	71.4	40-125	0			
Fluoranthene	1825	57	2546	0	71.7	55-115	0			
Fluorene	1981	57	2546	0	77.8	50-110	0			
Indeno(1,2,3-cd)pyrene	1776	57	2546	0	69.8	40-120	0			
Naphthalene	1759	57	2546	0	69.1	40-105	0			
Pyrene	1862	57	2546	0	73.2	45-125	0			
Surr: 2,4,6-Tribromophenol	2535	0	3182	0	79.7	34-140	0			
Surr: 2-Fluorobiphenyl	2092	0	3182	0	65.7	12-100	0			
Surr: 2-Fluorophenol	2086	0	3182	0	65.5	33-117	0			
Surr: 4-Terphenyl-d14	2140	0	3182	0	67.3	25-137	0			
Surr: Nitrobenzene-d5	2238	0	3182	0	70.3	37-107	0			
Surr: Phenol-d6	2202	0	3182	0	69.2	40-106	0			

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35213**      Instrument ID **SVMS4**      Method: **SW8270**

MSD				Sample ID: 1108889-20B MSD			Units: µg/Kg		Analysis Date: 8/31/2011 12:01 PM		
Client ID:		Run ID: SVMS4_110831A			SeqNo: 1722121		Prep Date: 8/30/2011		DF: 1		
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual	
Acenaphthene	1795	59	2643	0	67.9	45-110	1870	4.1	30		
Anthracene	1827	59	2643	0	69.1	55-105	2013	9.68	30		
Benzo(a)anthracene	1786	59	2643	0	67.6	50-110	1978	10.2	30		
Benzo(a)pyrene	1533	59	2643	0	58	50-110	1713	11.1	30		
Benzo(b)fluoranthene	1558	59	2643	0	59	45-115	1841	16.6	30		
Benzo(g,h,i)perylene	1546	59	2643	0	58.5	40-125	1757	12.7	30		
Benzo(k)fluoranthene	1821	59	2643	0	68.9	45-115	1815	0.336	30		
Chrysene	1799	59	2643	0	68.1	55-110	2037	12.4	30		
Dibenzo(a,h)anthracene	1560	59	2643	0	59	40-125	1818	15.3	30		
Fluoranthene	1657	59	2643	0	62.7	55-115	1825	9.61	30		
Fluorene	1877	59	2643	0	71	50-110	1981	5.42	30		
Indeno(1,2,3-cd)pyrene	1542	59	2643	0	58.4	40-120	1776	14.1	30		
Naphthalene	1737	59	2643	0	65.7	40-105	1759	1.25	30		
Pyrene	1657	59	2643	0	62.7	45-125	1862	11.7	30		
Surr: 2,4,6-Tribromophenol	2336	0	3304	0	70.7	34-140	2535	8.17	40		
Surr: 2-Fluorobiphenyl	2037	0	3304	0	61.6	12-100	2092	2.68	40		
Surr: 2-Fluorophenol	2061	0	3304	0	62.4	33-117	2086	1.21	40		
Surr: 4-Terphenyl-d14	1841	0	3304	0	55.7	25-137	2140	15	40		
Surr: Nitrobenzene-d5	2234	0	3304	0	67.6	37-107	2238	0.212	40		
Surr: Phenol-d6	2172	0	3304	0	65.7	40-106	2202	1.34	40		

The following samples were analyzed in this batch:      1108912-01B      1108912-02B      1108912-03B

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35235**      Instrument ID **SVMS6**      Method: **SW8270**

MBLK		Sample ID: <b>SBLKS1-35235-35235</b>				Units: <b>µg/Kg</b>		Analysis Date: <b>9/1/2011 11:58 AM</b>		
Client ID:		Run ID: <b>SVMS6_110901A</b>				SeqNo: <b>1724106</b>		Prep Date: <b>8/31/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Acenaphthene	ND	30								
Anthracene	ND	30								
Benzo(a)anthracene	ND	30								
Benzo(a)pyrene	ND	30								
Benzo(b)fluoranthene	ND	30								
Benzo(g,h,i)perylene	ND	30								
Benzo(k)fluoranthene	ND	30								
Chrysene	ND	30								
Dibenzo(a,h)anthracene	ND	30								
Fluoranthene	ND	30								
Fluorene	ND	30								
Indeno(1,2,3-cd)pyrene	ND	30								
Naphthalene	ND	30								
Pyrene	ND	30								
<hr/>										
<i>Surr: 2,4,6-Tribromophenol</i>	<i>1546</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>92.7</i>	<i>34-140</i>		<i>0</i>		
<i>Surr: 2-Fluorobiphenyl</i>	<i>1272</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>76.3</i>	<i>12-100</i>		<i>0</i>		
<hr/>										
<i>Surr: 2-Fluorophenol</i>	<i>1393</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>83.6</i>	<i>33-117</i>		<i>0</i>		
<i>Surr: 4-Terphenyl-d14</i>	<i>1410</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>84.6</i>	<i>25-137</i>		<i>0</i>		
<hr/>										
<i>Surr: Nitrobenzene-d5</i>	<i>1378</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>82.7</i>	<i>37-107</i>		<i>0</i>		
<i>Surr: Phenol-d6</i>	<i>1412</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>84.7</i>	<i>40-106</i>		<i>0</i>		

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35235**      Instrument ID **SVMS6**      Method: **SW8270**

LCS		Sample ID: <b>SLCSS1-35235-35235</b>				Units: <b>µg/Kg</b>		Analysis Date: <b>9/1/2011 11:04 AM</b>		
Client ID:		Run ID: <b>SVMS6_110901A</b>				SeqNo: <b>1724104</b>		Prep Date: <b>8/31/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Acenaphthene	1181	30	1333	0	88.6	45-110	0			
Anthracene	1242	30	1333	0	93.2	55-105	0			
Benzo(a)anthracene	1235	30	1333	0	92.6	50-110	0			
Benzo(a)pyrene	1314	30	1333	0	98.5	50-110	0			
Benzo(b)fluoranthene	1239	30	1333	0	92.9	45-115	0			
Benzo(g,h,i)perylene	1342	30	1333	0	101	40-125	0			
Benzo(k)fluoranthene	1309	30	1333	0	98.2	45-115	0			
Chrysene	1226	30	1333	0	92	55-110	0			
Dibenzo(a,h)anthracene	1367	30	1333	0	103	40-125	0			
Fluoranthene	1253	30	1333	0	94	55-115	0			
Fluorene	1191	30	1333	0	89.3	50-110	0			
Indeno(1,2,3-cd)pyrene	1361	30	1333	0	102	40-120	0			
Naphthalene	1164	30	1333	0	87.3	40-105	0			
Pyrene	1249	30	1333	0	93.7	45-125	0			
<i>Surr: 2,4,6-Tribromophenol</i>	<i>1689</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>101</i>	<i>34-140</i>	<i>0</i>			
<i>Surr: 2-Fluorobiphenyl</i>	<i>1305</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>78.3</i>	<i>12-100</i>	<i>0</i>			
<i>Surr: 2-Fluorophenol</i>	<i>1371</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>82.3</i>	<i>33-117</i>	<i>0</i>			
<i>Surr: 4-Terphenyl-d14</i>	<i>1483</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>89</i>	<i>25-137</i>	<i>0</i>			
<i>Surr: Nitrobenzene-d5</i>	<i>1395</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>83.7</i>	<i>37-107</i>	<i>0</i>			
<i>Surr: Phenol-d6</i>	<i>1345</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>80.7</i>	<i>40-106</i>	<i>0</i>			

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35235**      Instrument ID **SVMS6**      Method: **SW8270**

LCSD		Sample ID: <b>SLCSDS1-35235-35235</b>				Units: <b>µg/Kg</b>		Analysis Date: <b>9/1/2011 11:31 AM</b>		
Client ID:		Run ID: <b>SVMS6_110901A</b>				SeqNo: <b>1724105</b>		Prep Date: <b>8/31/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Acenaphthene	1092	30	1333	0	81.9	45-110	1181	7.8	25	
Anthracene	1188	30	1333	0	89.1	55-105	1242	4.44	25	
Benzo(a)anthracene	1204	30	1333	0	90.3	50-110	1235	2.49	25	
Benzo(a)pyrene	1270	30	1333	0	95.2	50-110	1314	3.41	25	
Benzo(b)fluoranthene	1348	30	1333	0	101	45-115	1239	8.43	25	
Benzo(g,h,i)perylene	1210	30	1333	0	90.7	40-125	1342	10.3	25	
Benzo(k)fluoranthene	1319	30	1333	0	98.9	45-115	1309	0.736	25	
Chrysene	1204	30	1333	0	90.3	55-110	1226	1.81	25	
Dibenzo(a,h)anthracene	1315	30	1333	0	98.6	40-125	1367	3.88	25	
Fluoranthene	1215	30	1333	0	91.2	55-115	1253	3.03	25	
Fluorene	1122	30	1333	0	84.1	50-110	1191	6	25	
Indeno(1,2,3-cd)pyrene	1300	30	1333	0	97.5	40-120	1361	4.53	25	
Naphthalene	1055	30	1333	0	79.1	40-105	1164	9.88	25	
Pyrene	1222	30	1333	0	91.7	45-125	1249	2.19	25	
<i>Surr: 2,4,6-Tribromophenol</i>	1595	0	1667	0	95.7	34-140	1689	5.72	40	
<i>Surr: 2-Fluorobiphenyl</i>	1199	0	1667	0	71.9	12-100	1305	8.47	40	
<i>Surr: 2-Fluorophenol</i>	1286	0	1667	0	77.2	33-117	1371	6.37	40	
<i>Surr: 4-Terphenyl-d14</i>	1471	0	1667	0	88.3	25-137	1483	0.812	40	
<i>Surr: Nitrobenzene-d5</i>	1287	0	1667	0	77.2	37-107	1395	8.06	40	
<i>Surr: Phenol-d6</i>	1263	0	1667	0	75.8	40-106	1345	6.31	40	

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.



**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35235**      Instrument ID **SVMS6**      Method: **SW8270**

MS				Sample ID: 1108926-27B MS			Units: µg/Kg		Analysis Date: 9/1/2011 02:59 PM		
Client ID:			Run ID: SVMS6_110901A			SeqNo: 1724444		Prep Date: 8/31/2011		DF: 1	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual	
Acenaphthene	2149	59	2604	0	82.5	45-110	0				
Anthracene	2305	59	2604	0	88.5	55-105	0				
Benzo(a)anthracene	2310	59	2604	0	88.7	50-110	0				
Benzo(a)pyrene	2418	59	2604	0	92.8	50-110	0				
Benzo(b)fluoranthene	2526	59	2604	0	97	45-115	0				
Benzo(g,h,i)perylene	2388	59	2604	0	91.7	40-125	0				
Benzo(k)fluoranthene	2621	59	2604	0	101	45-115	0				
Chrysene	2343	59	2604	0	90	55-110	0				
Dibenzo(a,h)anthracene	2505	59	2604	0	96.2	40-125	0				
Fluoranthene	2349	59	2604	0	90.2	55-115	0				
Fluorene	2201	59	2604	0	84.5	50-110	0				
Indeno(1,2,3-cd)pyrene	2452	59	2604	0	94.2	40-120	0				
Naphthalene	2088	59	2604	0	80.2	40-105	0				
Pyrene	2383	59	2604	0	91.5	45-125	0				
Surr: 2,4,6-Tribromophenol	3134	0	3256	0	96.3	34-140	0				
Surr: 2-Fluorobiphenyl	2329	0	3256	0	71.5	12-100	0				
Surr: 2-Fluorophenol	2506	0	3256	0	77	33-117	0				
Surr: 4-Terphenyl-d14	2667	0	3256	0	81.9	25-137	0				
Surr: Nitrobenzene-d5	2576	0	3256	0	79.1	37-107	0				
Surr: Phenol-d6	2487	0	3256	0	76.4	40-106	0				

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35235**      Instrument ID **SVMS6**      Method: **SW8270**

MSD				Sample ID: 1108926-27B MSD			Units: µg/Kg		Analysis Date: 9/1/2011 03:27 PM		
Client ID:			Run ID: SVMS6_110901A			SeqNo: 1724445		Prep Date: 8/31/2011		DF: 1	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual	
Acenaphthene	2190	60	2650	0	82.6	45-110	2149	1.9	30		
Anthracene	2305	60	2650	0	87	55-105	2305	0.00899	30		
Benzo(a)anthracene	2312	60	2650	0	87.3	50-110	2310	0.127	30		
Benzo(a)pyrene	2432	60	2650	0	91.8	50-110	2418	0.582	30		
Benzo(b)fluoranthene	2590	60	2650	0	97.7	45-115	2526	2.49	30		
Benzo(g,h,i)perylene	2439	60	2650	0	92	40-125	2388	2.1	30		
Benzo(k)fluoranthene	2420	60	2650	0	91.3	45-115	2621	7.97	30		
Chrysene	2320	60	2650	0	87.5	55-110	2343	0.986	30		
Dibenzo(a,h)anthracene	2516	60	2650	0	94.9	40-125	2505	0.439	30		
Fluoranthene	2340	60	2650	0	88.3	55-115	2349	0.382	30		
Fluorene	2230	60	2650	0	84.1	50-110	2201	1.3	30		
Indeno(1,2,3-cd)pyrene	2495	60	2650	0	94.1	40-120	2452	1.72	30		
Naphthalene	2120	60	2650	0	80	40-105	2088	1.53	30		
Pyrene	2352	60	2650	0	88.8	45-125	2383	1.3	30		
Surr: 2,4,6-Tribromophenol	3098	0	3313	0	93.5	34-140	3134	1.16	40		
Surr: 2-Fluorobiphenyl	2433	0	3313	0	73.4	12-100	2329	4.37	40		
Surr: 2-Fluorophenol	2566	0	3313	0	77.5	33-117	2506	2.37	40		
Surr: 4-Terphenyl-d14	2835	0	3313	0	85.6	25-137	2667	6.12	40		
Surr: Nitrobenzene-d5	2595	0	3313	0	78.3	37-107	2576	0.731	40		
Surr: Phenol-d6	2524	0	3313	0	76.2	40-106	2487	1.46	40		

The following samples were analyzed in this batch:    1108912-04B    1108912-05B    1108912-06B

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35319**      Instrument ID **SVMS4**      Method: **SW8270**

**MBLK**      Sample ID: **SBLKS1-35319-35319**      Units: **µg/Kg**      Analysis Date: **9/6/2011 06:49 PM**

Client ID:      Run ID: **SVMS4\_110906A**      SeqNo: **1726859**      Prep Date: **9/6/2011**      DF: **1**

Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Acenaphthene	ND	30								
Anthracene	ND	30								
Benzo(a)anthracene	ND	30								
Benzo(a)pyrene	ND	30								
Benzo(b)fluoranthene	ND	30								
Benzo(g,h,i)perylene	ND	30								
Benzo(k)fluoranthene	ND	30								
Chrysene	ND	30								
Dibenzo(a,h)anthracene	ND	30								
Fluoranthene	ND	30								
Fluorene	ND	30								
Indeno(1,2,3-cd)pyrene	ND	30								
Naphthalene	ND	30								
Pyrene	ND	30								
<i>Surr: 2,4,6-Tribromophenol</i>	<i>1431</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>85.8</i>	<i>34-140</i>	<i>0</i>			
<i>Surr: 2-Fluorobiphenyl</i>	<i>1104</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>66.2</i>	<i>12-100</i>	<i>0</i>			
<i>Surr: 2-Fluorophenol</i>	<i>1304</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>78.3</i>	<i>33-117</i>	<i>0</i>			
<i>Surr: 4-Terphenyl-d14</i>	<i>1694</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>102</i>	<i>25-137</i>	<i>0</i>			
<i>Surr: Nitrobenzene-d5</i>	<i>1128</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>67.7</i>	<i>37-107</i>	<i>0</i>			
<i>Surr: Phenol-d6</i>	<i>1226</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>73.6</i>	<i>40-106</i>	<i>0</i>			

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

# QC BATCH REPORT

Batch ID: **35319**      Instrument ID **SVMS4**      Method: **SW8270**

**LCS**      Sample ID: **SLCSS1-35319-35319**      Units: **µg/Kg**      Analysis Date: **9/6/2011 05:47 PM**

Client ID:      Run ID: **SVMS4\_110906A**      SeqNo: **1726857**      Prep Date: **9/6/2011**      DF: **1**

Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Acenaphthene	1139	30	1333	0	85.4	45-110	0			
Anthracene	1257	30	1333	0	94.3	55-105	0			
Benzo(a)anthracene	1363	30	1333	0	102	50-110	0			
Benzo(a)pyrene	1310	30	1333	0	98.3	50-110	0			
Benzo(b)fluoranthene	1232	30	1333	0	92.4	45-115	0			
Benzo(g,h,i)perylene	1274	30	1333	0	95.5	40-125	0			
Benzo(k)fluoranthene	1252	30	1333	0	93.9	45-115	0			
Chrysene	1277	30	1333	0	95.8	55-110	0			
Dibenzo(a,h)anthracene	1295	30	1333	0	97.2	40-125	0			
Fluoranthene	1261	30	1333	0	94.6	55-115	0			
Fluorene	1367	30	1333	0	103	50-110	0			
Indeno(1,2,3-cd)pyrene	1310	30	1333	0	98.3	40-120	0			
Naphthalene	1133	30	1333	0	85	40-105	0			
Pyrene	1306	30	1333	0	98	45-125	0			
<i>Surr: 2,4,6-Tribromophenol</i>	<i>1544</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>92.6</i>	<i>34-140</i>	<i>0</i>			
<i>Surr: 2-Fluorobiphenyl</i>	<i>1248</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>74.9</i>	<i>12-100</i>	<i>0</i>			
<i>Surr: 2-Fluorophenol</i>	<i>1197</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>71.8</i>	<i>33-117</i>	<i>0</i>			
<i>Surr: 4-Terphenyl-d14</i>	<i>1641</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>98.5</i>	<i>25-137</i>	<i>0</i>			
<i>Surr: Nitrobenzene-d5</i>	<i>1199</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>71.9</i>	<i>37-107</i>	<i>0</i>			
<i>Surr: Phenol-d6</i>	<i>1247</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>74.8</i>	<i>40-106</i>	<i>0</i>			

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

# QC BATCH REPORT

Batch ID: **35319**      Instrument ID **SVMS4**      Method: **SW8270**

LCSD				Sample ID: <b>SLCSDS1-35319-35319</b>			Units: <b>µg/Kg</b>		Analysis Date: <b>9/6/2011 06:18 PM</b>	
Client ID:				Run ID: <b>SVMS4_110906A</b>			SeqNo: <b>1726858</b>		Prep Date: <b>9/6/2011</b>	
							DF: <b>1</b>			
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Acenaphthene	1023	30	1333	0	76.7	45-110	1139	10.8	25	
Anthracene	1127	30	1333	0	84.5	55-105	1257	11	25	
Benzo(a)anthracene	1247	30	1333	0	93.5	50-110	1363	8.86	25	
Benzo(a)pyrene	1182	30	1333	0	88.7	50-110	1310	10.3	25	
Benzo(b)fluoranthene	1178	30	1333	0	88.4	45-115	1232	4.45	25	
Benzo(g,h,i)perylene	1158	30	1333	0	86.8	40-125	1274	9.54	25	
Benzo(k)fluoranthene	1239	30	1333	0	92.9	45-115	1252	1.02	25	
Chrysene	1148	30	1333	0	86.1	55-110	1277	10.6	25	
Dibenzo(a,h)anthracene	1179	30	1333	0	88.4	40-125	1295	9.43	25	
Fluoranthene	1120	30	1333	0	84	55-115	1261	11.9	25	
Fluorene	1241	30	1333	0	93.1	50-110	1367	9.67	25	
Indeno(1,2,3-cd)pyrene	1182	30	1333	0	88.6	40-120	1310	10.3	25	
Naphthalene	1070	30	1333	0	80.3	40-105	1133	5.69	25	
Pyrene	1181	30	1333	0	88.6	45-125	1306	10.1	25	
<i>Surr: 2,4,6-Tribromophenol</i>	<i>1425</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>85.5</i>	<i>34-140</i>	<i>1544</i>	<i>8.04</i>	<i>40</i>	
<i>Surr: 2-Fluorobiphenyl</i>	<i>1153</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>69.2</i>	<i>12-100</i>	<i>1248</i>	<i>7.97</i>	<i>40</i>	
<i>Surr: 2-Fluorophenol</i>	<i>1190</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>71.4</i>	<i>33-117</i>	<i>1197</i>	<i>0.531</i>	<i>40</i>	
<i>Surr: 4-Terphenyl-d14</i>	<i>1475</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>88.5</i>	<i>25-137</i>	<i>1641</i>	<i>10.7</i>	<i>40</i>	
<i>Surr: Nitrobenzene-d5</i>	<i>1140</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>68.4</i>	<i>37-107</i>	<i>1199</i>	<i>5.04</i>	<i>40</i>	
<i>Surr: Phenol-d6</i>	<i>1200</i>	<i>0</i>	<i>1667</i>	<i>0</i>	<i>72</i>	<i>40-106</i>	<i>1247</i>	<i>3.87</i>	<i>40</i>	

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35319**      Instrument ID **SVMS4**      Method: **SW8270**

MS				Sample ID: 1109075-05A MS			Units: µg/Kg		Analysis Date: 9/6/2011 07:20 PM		
Client ID:		Run ID: SVMS4_110906A			SeqNo: 1726860		Prep Date: 9/6/2011		DF: 1		
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual	
Acenaphthene	2186	60	2649	0	82.5	45-110	0				
Anthracene	2417	60	2649	0	91.2	55-105	0				
Benzo(a)anthracene	2616	60	2649	0	98.8	50-110	0				
Benzo(a)pyrene	2471	60	2649	0	93.3	50-110	0				
Benzo(b)fluoranthene	2380	60	2649	0	89.9	45-115	0				
Benzo(g,h,i)perylene	2593	60	2649	0	97.9	40-125	0				
Benzo(k)fluoranthene	2754	60	2649	0	104	45-115	0				
Chrysene	2462	60	2649	0	92.9	55-110	0				
Dibenzo(a,h)anthracene	2514	60	2649	0	94.9	40-125	0				
Fluoranthene	2419	60	2649	0	91.3	55-115	0				
Fluorene	2657	60	2649	0	100	50-110	0				
Indeno(1,2,3-cd)pyrene	2534	60	2649	0	95.7	40-120	0				
Naphthalene	2314	60	2649	0	87.4	40-105	0				
Pyrene	2500	60	2649	0	94.4	45-125	0				
Surr: 2,4,6-Tribromophenol	3166	0	3312	0	95.6	34-140	0				
Surr: 2-Fluorobiphenyl	2423	0	3312	0	73.2	12-100	0				
Surr: 2-Fluorophenol	2601	0	3312	0	78.5	33-117	0				
Surr: 4-Terphenyl-d14	3040	0	3312	0	91.8	25-137	0				
Surr: Nitrobenzene-d5	2491	0	3312	0	75.2	37-107	0				
Surr: Phenol-d6	2648	0	3312	0	80	40-106	0				

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35319**      Instrument ID **SVMS4**      Method: **SW8270**

MSD				Sample ID: <b>1109075-05A MSD</b>			Units: <b>µg/Kg</b>		Analysis Date: <b>9/6/2011 07:51 PM</b>	
Client ID:				Run ID: <b>SVMS4_110906A</b>			SeqNo: <b>1726861</b>		Prep Date: <b>9/6/2011</b>	
									DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Acenaphthene	2103	58	2574	0	81.7	45-110	2186	3.89	30	
Anthracene	2241	58	2574	0	87.1	55-105	2417	7.57	30	
Benzo(a)anthracene	2465	58	2574	0	95.8	50-110	2616	5.94	30	
Benzo(a)pyrene	2330	58	2574	0	90.5	50-110	2471	5.85	30	
Benzo(b)fluoranthene	2279	58	2574	0	88.6	45-115	2380	4.34	30	
Benzo(g,h,i)perylene	2391	58	2574	0	92.9	40-125	2593	8.1	30	
Benzo(k)fluoranthene	2499	58	2574	0	97.1	45-115	2754	9.69	30	
Chrysene	2288	58	2574	0	88.9	55-110	2462	7.34	30	
Dibenzo(a,h)anthracene	2326	58	2574	0	90.4	40-125	2514	7.77	30	
Fluoranthene	2219	58	2574	0	86.2	55-115	2419	8.6	30	
Fluorene	2516	58	2574	0	97.8	50-110	2657	5.43	30	
Indeno(1,2,3-cd)pyrene	2360	58	2574	0	91.7	40-120	2534	7.1	30	
Naphthalene	2283	58	2574	0	88.7	40-105	2314	1.35	30	
Pyrene	2349	58	2574	0	91.3	45-125	2500	6.22	30	
<i>Surr: 2,4,6-Tribromophenol</i>	2930	0	3218	0	91.1	34-140	3166	7.75	40	
<i>Surr: 2-Fluorobiphenyl</i>	2413	0	3218	0	75	12-100	2423	0.429	40	
<i>Surr: 2-Fluorophenol</i>	2640	0	3218	0	82.1	33-117	2601	1.5	40	
<i>Surr: 4-Terphenyl-d14</i>	2827	0	3218	0	87.9	25-137	3040	7.27	40	
<i>Surr: Nitrobenzene-d5</i>	2551	0	3218	0	79.3	37-107	2491	2.4	40	
<i>Surr: Phenol-d6</i>	2708	0	3218	0	84.2	40-106	2648	2.26	40	

The following samples were analyzed in this batch: | 1108912-06B |

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **R94139**      Instrument ID **VMS6**      Method: **SW8260**

<b>MBLK</b>	Sample ID: <b>VBLKW3-110831-R94139</b>				Units: <b>µg/L</b>		Analysis Date: <b>9/1/2011 05:32 AM</b>			
Client ID:	Run ID: <b>VMS6_110831B</b>				SeqNo: <b>1722685</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Benzene	ND	1.0								
Ethylbenzene	ND	1.0								
m,p-Xylene	ND	2.0								
o-Xylene	ND	1.0								
Toluene	ND	1.0								
Xylenes, Total	ND	2.0								
<i>Surr: 1,2-Dichloroethane-d4</i>	99.77	0	100	0	99.8	70-120	0			
<i>Surr: 4-Bromofluorobenzene</i>	98.59	0	100	0	98.6	75-120	0			
<i>Surr: Dibromofluoromethane</i>	98.56	0	100	0	98.6	85-115	0			
<i>Surr: Toluene-d8</i>	95.9	0	100	0	95.9	85-120	0			

<b>LCS</b>	Sample ID: <b>VLCSW2-110831-R94139</b>				Units: <b>µg/L</b>		Analysis Date: <b>9/1/2011 04:17 AM</b>			
Client ID:	Run ID: <b>VMS6_110831B</b>				SeqNo: <b>1722683</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Benzene	20.86	1.0	20	0	104	80-120	0			
Ethylbenzene	20.74	1.0	20	0	104	75-125	0			
m,p-Xylene	42.12	2.0	40	0	105	75-130	0			
o-Xylene	20.7	1.0	20	0	104	80-120	0			
Toluene	20.68	1.0	20	0	103	75-120	0			
Xylenes, Total	62.82	2.0	60	0	105	75-130	0			
<i>Surr: 1,2-Dichloroethane-d4</i>	99.47	0	100	0	99.5	70-120	0			
<i>Surr: 4-Bromofluorobenzene</i>	103.1	0	100	0	103	75-120	0			
<i>Surr: Dibromofluoromethane</i>	100.9	0	100	0	101	85-115	0			
<i>Surr: Toluene-d8</i>	99.09	0	100	0	99.1	85-120	0			

<b>LCSD</b>	Sample ID: <b>VLCSDW2-110831-R94139</b>				Units: <b>µg/L</b>		Analysis Date: <b>9/1/2011 04:42 AM</b>			
Client ID:	Run ID: <b>VMS6_110831B</b>				SeqNo: <b>1722684</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Benzene	19.4	1.0	20	0	97	80-120	20.86	7.25	30	
Ethylbenzene	19.48	1.0	20	0	97.4	75-125	20.74	6.27	30	
m,p-Xylene	38.81	2.0	40	0	97	75-130	42.12	8.18	30	
o-Xylene	19.44	1.0	20	0	97.2	80-120	20.7	6.28	30	
Toluene	19.37	1.0	20	0	96.8	75-120	20.68	6.54	30	
Xylenes, Total	58.25	2.0	60	0	97.1	75-130	62.82	7.55	30	
<i>Surr: 1,2-Dichloroethane-d4</i>	98.84	0	100	0	98.8	70-120	99.47	0.635	30	
<i>Surr: 4-Bromofluorobenzene</i>	103.4	0	100	0	103	75-120	103.1	0.339	30	
<i>Surr: Dibromofluoromethane</i>	101.3	0	100	0	101	85-115	100.9	0.396	30	
<i>Surr: Toluene-d8</i>	96.15	0	100	0	96.2	85-120	99.09	3.01	30	

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.



**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **R94139**      Instrument ID **VMS6**      Method: **SW8260**

MS				Units: <b>µg/Kg</b>			Analysis Date: <b>9/1/2011 01:57 PM</b>			
Sample ID: <b>1108912-01A MS</b>										
Client ID: <b>TR 22-20-597 N. Bottom</b>				Run ID: <b>VMS6_110831B</b>			SeqNo: <b>1723399</b>		Prep Date:	
									DF: <b>100</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Benzene	2028	100	2000	0	101	75-125	0			
Ethylbenzene	1894	200	2000	0	94.7	75-125	0			
m,p-Xylene	3911	200	4000	0	97.8	80-125	0			
o-Xylene	1975	100	2000	0	98.8	75-125	0			
Toluene	1865	150	2000	0	93.2	70-125	0			
Xylenes, Total	5886	300	6000	0	98.1	75-125	0			
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>10070</i>	<i>0</i>	<i>10000</i>	<i>0</i>	<i>101</i>	<i>70-120</i>	<i>0</i>			
<i>Surr: 4-Bromofluorobenzene</i>	<i>11130</i>	<i>0</i>	<i>10000</i>	<i>0</i>	<i>111</i>	<i>75-120</i>	<i>0</i>			
<i>Surr: Dibromofluoromethane</i>	<i>9675</i>	<i>0</i>	<i>10000</i>	<i>0</i>	<i>96.8</i>	<i>85-115</i>	<i>0</i>			
<i>Surr: Toluene-d8</i>	<i>9694</i>	<i>0</i>	<i>10000</i>	<i>0</i>	<i>96.9</i>	<i>85-115</i>	<i>0</i>			

MSD				Units: <b>µg/Kg</b>			Analysis Date: <b>9/1/2011 02:22 PM</b>			
Sample ID: <b>1108912-01A MSD</b>										
Client ID: <b>TR 22-20-597 N. Bottom</b>				Run ID: <b>VMS6_110831B</b>			SeqNo: <b>1723400</b>		Prep Date:	
									DF: <b>100</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Benzene	2005	100	2000	0	100	75-125	2028	1.14	30	
Ethylbenzene	1818	200	2000	0	90.9	75-125	1894	4.09	30	
m,p-Xylene	3747	200	4000	0	93.7	80-125	3911	4.28	30	
o-Xylene	1872	100	2000	0	93.6	75-125	1975	5.35	30	
Toluene	1858	150	2000	0	92.9	70-125	1865	0.376	30	
Xylenes, Total	5619	300	6000	0	93.6	75-125	5886	4.64	30	
<i>Surr: 1,2-Dichloroethane-d4</i>	<i>9675</i>	<i>0</i>	<i>10000</i>	<i>0</i>	<i>96.8</i>	<i>70-120</i>	<i>10070</i>	<i>4.04</i>	<i>30</i>	
<i>Surr: 4-Bromofluorobenzene</i>	<i>10210</i>	<i>0</i>	<i>10000</i>	<i>0</i>	<i>102</i>	<i>75-120</i>	<i>11130</i>	<i>8.6</i>	<i>30</i>	
<i>Surr: Dibromofluoromethane</i>	<i>9612</i>	<i>0</i>	<i>10000</i>	<i>0</i>	<i>96.1</i>	<i>85-115</i>	<i>9675</i>	<i>0.653</i>	<i>30</i>	
<i>Surr: Toluene-d8</i>	<i>9698</i>	<i>0</i>	<i>10000</i>	<i>0</i>	<i>97</i>	<i>85-115</i>	<i>9694</i>	<i>0.0413</i>	<i>30</i>	

The following samples were analyzed in this batch:

1108912-01A	1108912-02A	1108912-03A
1108912-04A	1108912-05A	1108912-06A

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35281**      Instrument ID **WETCHEM**      Method: **SW7196A**

<b>MBLK</b>	Sample ID: <b>MBLK-35281-35281</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/2/2011 01:00 PM</b>			
Client ID:	Run ID: <b>WETCHEM_110902C</b>				SeqNo: <b>1724443</b>		Prep Date: <b>9/1/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Chromium, Hexavalent	ND	0.49								

<b>LCS</b>	Sample ID: <b>LCS-35281-35281</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/2/2011 01:00 PM</b>			
Client ID:	Run ID: <b>WETCHEM_110902C</b>				SeqNo: <b>1724441</b>		Prep Date: <b>9/1/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Chromium, Hexavalent	2.054	0.48	1.931		0	106	75-110	0		

<b>LCSD</b>	Sample ID: <b>LCSD-35281-35281</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/2/2011 01:00 PM</b>			
Client ID:	Run ID: <b>WETCHEM_110902C</b>				SeqNo: <b>1724442</b>		Prep Date: <b>9/1/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Chromium, Hexavalent	2.078	0.49	1.953		0	106	75-110	2.054	1.17	20

<b>MS</b>	Sample ID: <b>1108868-01A MS</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/2/2011 01:00 PM</b>			
Client ID:	Run ID: <b>WETCHEM_110902C</b>				SeqNo: <b>1724424</b>		Prep Date: <b>9/1/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Chromium, Hexavalent	1.616	0.48	1.938	0.2326	71.4	60-130		0		

<b>MSD</b>	Sample ID: <b>1108868-01A MSD</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/2/2011 01:00 PM</b>			
Client ID:	Run ID: <b>WETCHEM_110902C</b>				SeqNo: <b>1724425</b>		Prep Date: <b>9/1/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Chromium, Hexavalent	1.806	0.50	1.984	0.2326	79.3	60-130	1.616	11.1	30	

The following samples were analyzed in this batch:

1108912-01B	1108912-02B	1108912-03B
1108912-04B	1108912-05B	1108912-06B

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **R94050**      Instrument ID **WETCHEM**      Method: **A4500-H B**

**DUP**      Sample ID: **1108914-01A DUP**      Units: **s.u.**      Analysis Date: **8/30/2011 10:15 AM**

Client ID:      Run ID: **WETCHEM\_110830F**      SeqNo: **1720591**      Prep Date:      DF: **1**

Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
pH	8.57	0	0	0	0	0-0	8.57	0	20	

**DUP**      Sample ID: **1108868-01A DUP**      Units: **s.u.**      Analysis Date: **8/30/2011 10:15 AM**

Client ID:      Run ID: **WETCHEM\_110830F**      SeqNo: **1720594**      Prep Date:      DF: **1**

Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
pH	8.26	0	0	0	0	0-0	8.26	0	20	

**DUP**      Sample ID: **1108912-05B DUP**      Units: **s.u.**      Analysis Date: **8/30/2011 10:15 AM**

Client ID: **TR 22-20-597 W. Wall**      Run ID: **WETCHEM\_110830F**      SeqNo: **1720606**      Prep Date:      DF: **1**

Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
pH	8.65	0	0	0	0	0-0	8.65	0	20	H

The following samples were analyzed in this batch:

1108912-01B	1108912-02B	1108912-03B
1108912-04B	1108912-05B	1108912-06B

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108912  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **R94091**      Instrument ID **MOIST**      Method: **A2540 G**

<b>MBLK</b>	Sample ID: <b>WBLKS1-R94091</b>				Units: % of sample			Analysis Date: <b>8/30/2011 02:49 PM</b>		
Client ID:	Run ID: <b>MOIST_110830C</b>				SeqNo: <b>1721824</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Moisture	ND	0.050								

<b>LCS</b>	Sample ID: <b>LCS-R94091</b>				Units: % of sample			Analysis Date: <b>8/30/2011 02:49 PM</b>		
Client ID:	Run ID: <b>MOIST_110830C</b>				SeqNo: <b>1721823</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Moisture	99.99	0.050	100	0	100	99.5-100.5	0			

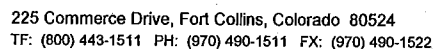
<b>DUP</b>	Sample ID: <b>1108898-01ADUP</b>				Units: % of sample			Analysis Date: <b>8/30/2011 02:49 PM</b>		
Client ID:	Run ID: <b>MOIST_110830C</b>				SeqNo: <b>1721813</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Moisture	0.75	0.050	0	0	0	0-0	0.76	1.32	20	

<b>DUP</b>	Sample ID: <b>1108912-01BDUP</b>				Units: % of sample			Analysis Date: <b>8/30/2011 02:49 PM</b>		
Client ID: <b>TR 22-20-597 N. Bottom</b>	Run ID: <b>MOIST_110830C</b>				SeqNo: <b>1721817</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Moisture	14.53	0.050	0	0	0	0-0	14.03	3.5	20	

The following samples were analyzed in this batch:

1108912-01B	1108912-02B	1108912-03B
1108912-04B	1108912-05B	1108912-06B

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.



1208912

Form 202r8

PAGE 1 of 1

DISPOSAL ☒ By Lab or ☐ Return to Client

\*Time Zone (Circle): EST CST MST PST Matrix: O = oil S = soil NS = non-soil solid W = water L = liquid E = extract F = filter

**For metals or anions, please detail analytes below.**

	SIGNATURE	PRINTED NAME	DATE	TIME
RELINQUISHED BY	<i>[Signature]</i>	<i>[Signature]</i>	8/29/11	6pm
RECEIVED BY	<i>[Signature]</i>	Diane F. Shaw	8/30/11	1030
RELINQUISHED BY				
RECEIVED BY				
RELINQUISHED BY				
RECEIVED BY				



Environmental

**Subcontractor:**A & L Great Lakes Agricultural La  
3505 Conestoga Dr

Ft. Wayne, IN 46808

TEL: (260) 483-4759

FAX: (260) 483-5274

Acct #: 91000

**CHAIN-OF-CUSTODY RECORD**

Page 1 of 1

Date: **30-Aug-11**COC ID: **3077**Due Date **07-Sep-11**

Customer Information		Project Information		Parameter/Method Request for Analysis										
Purchase Order		Project Name	1108912	A Subcontracted Analyses (SUBCONTRACT)										
Work Order		Project Number		B										
Company Name	ALS Group USA, Corp	Bill To Company	ALS Group USA, Corp	C										
Send Report To	Ann Preston	Inv Attn	Accounts Payable	D										
Address	3352 128th Avenue	Address	3352 128th Avenue	E										
				F										
City/State/Zip	Holland, Michigan 49424-9263	City/State/Zip	Holland, Michigan 49424-9263	G										
Phone	(616) 399-6070	Phone	(616) 399-6070	H										
Fax	(616) 399-6185	Fax	(616) 399-6185	I										
eMail Address	ann.preston@alsglobal.com	eMail CC		J										
Sample ID	Matrix	Collection Date 24hr	Bottle	A	B	C	D	E	F	G	H	I	J	
1108912-01C (TR 22-20-597 N. Bottom)	Soil	24/Aug/2011 14:20	(1) MISC	X										
1108912-02C (TR 22-20-597 N. Wall)	Soil	24/Aug/2011 13:45	(1) MISC	X										
1108912-03C (TR 22-20-597 S Bottom)	Soil	24/Aug/2011 14:15	(1) MISC	X										
1108912-04C (TR 22-20-597 S. Wall)	Soil	24/Aug/2011 14:00	(1) MISC	X										
1108912-05C (TR 22-20-597 W. Wall)	Soil	24/Aug/2011 13:50	(1) MISC	X										
1108912-06C (TR 22-20-597 E. Wall)	Soil	24/Aug/2011 14:10	(1) MISC	X										

**Comments:**Please analyze for SAR-EC. Email results to Ann Preston.

Relinquished by:	Date/Time	Received by:	Date/Time	Cooler IDs	Report/QC Level
					<b>Std</b>
Relinquished by:	Date/Time	Received by:	Date/Time		

Sample Receipt Checklist

Client Name: HRL

Date/Time Received: 30-Aug-11 10:30

Work Order: 1108912

Received by: DS

Checklist completed by Diane Shaw 30-Aug-11  
eSignature Date

Reviewed by: Ann Preston 30-Aug-11  
eSignature Date

Matrices: Soil

Carrier name: FedEx

Shipping container/cooler in good condition?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	Not Present <input type="checkbox"/>
Custody seals intact on shipping container/cooler?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	Not Present <input type="checkbox"/>
Custody seals intact on sample bottles?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	Not Present <input checked="" type="checkbox"/>
Chain of custody present?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Chain of custody signed when relinquished and received?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Chain of custody agrees with sample labels?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Samples in proper container/bottle?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Sample containers intact?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Sufficient sample volume for indicated test?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
All samples received within holding time?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Container/Temp Blank temperature in compliance?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Temperature(s)/Thermometer(s):	<u>4.4 c</u>		
Cooler(s)/Kit(s):			
Water - VOA vials have zero headspace?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	No VOA vials submitted <input checked="" type="checkbox"/>
Water - pH acceptable upon receipt?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	N/A <input checked="" type="checkbox"/>
pH adjusted?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	N/A <input checked="" type="checkbox"/>
pH adjusted by:			
Login Notes:			

-----

Client Contacted:

Date Contacted:

Person Contacted:

Contacted By:

Regarding:

Comments:

CorrectiveAction:

CUSTODY SEAL

DATE

SIGNATURE

QEC

Quality Environmental Containers  
800-255-3950 • 304-255-3900FedEx Express NEW Package  
US AirbillFedEx  
Tracking  
Number

8758 3471 3970

0200

Form  
ID No.

FedEx Retrieval Copy

## 1 From

Date

8/29/11

Sender's FedEx  
Account NumberSender's  
Name

Reed, David

Phone

710 433-3271

Company

HCS

Address

744 N. 2nd St. 140

City

Grand Junction

State

CO

ZIP

81506

## 2 Your Internal Billing Reference

## 3 To

Recipient's  
Name

Samuel Believing

Phone

616 379-6070

Company

ALS Group

Address

2352 12th Ave

We cannot deliver to P.O. boxes or P.O. ZIP codes.

Address

Use this line for the HOLD location address or for continuation of your shipping address.

City

Halters

State

MT

ZIP

49424

01

HOLD Weekday

FedEx location address

REQUIRED. NOT available for

FedEx First Overnight.

31

HOLD Saturday

FedEx location address

REQUIRED. Available ONLY for

FedEx Priority Overnight and

FedEx 2Day to select locations.

## 4 Express Package Service

NOTE: Service order has changed. Please select carefully.

## Next Business Day

06

FedEx First Overnight

Earliest next business morning delivery to select

locations. Friday shipments will be delivered on

Monday unless SATURDAY Delivery is selected.

01

FedEx Priority Overnight

Next business morning. \* Friday shipments will be

delivered on Monday unless SATURDAY Delivery

is selected.

05

FedEx Standard Overnight

Next business afternoon. \*

Saturday Delivery NOT available.

## 5 Packaging

\* Declared value limit \$500.

06

FedEx Envelope\*

02

FedEx Pak\*

03

FedEx Box\*

04

FedEx Tube

01

Other

## 6 Special Handling and Delivery Signature Options

## 03 SATURDAY DELIVERY

No Signature Required

Package may be left without

obtaining a signature for delivery.

10

Direct Signature

Someone at recipient's address

may sign for delivery. Fee applies.

34

Indirect Signature

If no one is available at recipient's

address, someone at a neighboring

address may sign for delivery. For

residential deliveries only. Fee applies.

Does this shipment contain dangerous goods?

One box must be checked.

No 04

Yes

As per attached

Shipper's Declaration.

Yes

Shipper's Declaration

not required.

06

Dry Ice

Dry Ice, 3, UN 1845

x

kg

Cargo Aircraft Only

Dangerous goods (including dry ice) cannot be shipped in FedEx packaging

or placed in a FedEx Express Drop Box.

7 Payment Bill to:

1

Sender

Acct. No. in Section

1 will be billed.

Enter FedEx Acct. No. or Credit Card No. below.

2

Recipient

3

Third Party

4

Credit Card

Obtain recip.

Acct. No.

5

Cash/Check

Total Packages

1

Total Weight

72

lbs.

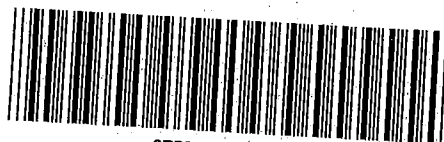
Credit Card Auth.

Your liability is limited to \$100 unless you declare a higher value. See the current FedEx Service Guide for details.

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8758 3471 3970

612



**Appendix 2: North, South Pit Bottom and Southern Pit Wall Additional Excavation Raw  
Analytical Data**



28-Sep-2011

Kris Rowe  
HRL Compliance Solutions  
744 Horizon Ct. Suite 140  
Grand Junction, CO 81506

Re: **Williams TR 22-20-597 Pad LOE 9/19/11**

Work Order: **1109654**

Dear Kris,

ALS Environmental received 3 samples on 21-Sep-2011 10:30 AM for the analyses presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested.

QC sample results for this data met laboratory specifications. Any exceptions are noted in the Case Narrative, or noted with qualifiers in the report or QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained from ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

The total number of pages in this report is 11.

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

Sincerely,

A handwritten signature in cursive script that reads "Ann Preston".

Electronically approved by: Ann Preston

Ann Preston  
Project Manager



Certificate No: IL100452

ADDRESS 3352 128th Avenue Holland, Michigan 49424-9263 | PHONE (616) 399-6070 | FAX (616) 399-6185

ALS GROUP USA, CORP Part of the ALS Laboratory Group A Campbell Brothers Limited Company

Environmental The ALS logo, a stylized blue triangle with a yellow flame inside.

[www.alsglobal.com](http://www.alsglobal.com)

RIGHT SOLUTIONS RIGHT PARTNER

---

**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Pad LOE 9/19/11  
**Work Order:** 1109654

---

**Work Order Sample Summary**

---

<u>Lab Samp ID</u>	<u>Client Sample ID</u>	<u>Matrix</u>	<u>Tag Number</u>	<u>Collection Date</u>	<u>Date Received</u>	<u>Hold</u>
1109654-01	TR 22-20-597 N. Bottom	Soil		9/19/2011 15:15	9/21/2011 10:30	<input type="checkbox"/>
1109654-02	TR 22-20-597 S. Bottom	Soil		9/19/2011 15:20	9/21/2011 10:30	<input type="checkbox"/>
1109654-03	TR 22-20-597 S. Wall	Soil		9/19/2011 15:25	9/21/2011 10:30	<input type="checkbox"/>

---

**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Pad LOE 9/19/11  
**WorkOrder:** 1109654

**QUALIFIERS,  
ACRONYMS, UNITS**

<b><u>Qualifier</u></b>	<b><u>Description</u></b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL

<b><u>Acronym</u></b>	<b><u>Description</u></b>
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
TDL	Target Detection Limit

<b><u>Units Reported</u></b>	<b><u>Description</u></b>
% of sample	Percent of Sample
mg/Kg-dry	Milligrams per Kilogram Dry Weight

# ALS Group USA, Corp

Date: 28-Sep-11

**Client:** HRL Compliance Solutions

**Project:** Williams TR 22-20-597 Pad LOE 9/19/11

**Work Order:** 1109654

**Sample ID:** TR 22-20-597 N. Bottom

**Lab ID:** 1109654-01

**Collection Date:** 9/19/2011 03:15 PM

**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<hr/>						
<b>DIESEL RANGE ORGANICS BY GC-FID</b>			<b>SW8015M</b>		Prep Date: <b>9/21/2011</b>	Analyst: <b>RM</b>
<b>DRO (C10-C28)</b>	<b>31</b>		<b>5.4</b>	<b>mg/Kg-dry</b>	<b>1</b>	9/22/2011 05:43 PM
<i>Surr: 4-Terphenyl-d14</i>	<i>63.1</i>		<i>39-115</i>	<i>%REC</i>	<i>1</i>	9/22/2011 05:43 PM
<b>MOISTURE</b>			<b>A2540 G</b>			Analyst: <b>CG</b>
<b>Moisture</b>	<b>24</b>		<b>0.050</b>	<b>% of sample</b>	<b>1</b>	9/21/2011 04:33 PM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

# ALS Group USA, Corp

Date: 28-Sep-11

**Client:** HRL Compliance Solutions

**Project:** Williams TR 22-20-597 Pad LOE 9/19/11

**Work Order:** 1109654

**Sample ID:** TR 22-20-597 S. Bottom

**Lab ID:** 1109654-02

**Collection Date:** 9/19/2011 03:20 PM

**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<hr/>						
<b>DIESEL RANGE ORGANICS BY GC-FID</b>			<b>SW8015M</b>		Prep Date: <b>9/21/2011</b>	Analyst: <b>RM</b>
<b>DRO (C10-C28)</b>	<b>71</b>		<b>4.7</b>	<b>mg/Kg-dry</b>	<b>1</b>	9/22/2011 05:43 PM
<i>Surr: 4-Terphenyl-d14</i>	39.3		39-115	%REC	1	9/22/2011 05:43 PM
<b>MOISTURE</b>			<b>A2540 G</b>			Analyst: <b>CG</b>
<b>Moisture</b>	<b>13</b>		<b>0.050</b>	<b>% of sample</b>	<b>1</b>	9/21/2011 04:33 PM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

# ALS Group USA, Corp

Date: 28-Sep-11

**Client:** HRL Compliance Solutions

**Project:** Williams TR 22-20-597 Pad LOE 9/19/11

**Work Order:** 1109654

**Sample ID:** TR 22-20-597 S. Wall

**Lab ID:** 1109654-03

**Collection Date:** 9/19/2011 03:25 PM

**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<hr/>						
<b>DIESEL RANGE ORGANICS BY GC-FID</b>			<b>SW8015M</b>		Prep Date: <b>9/21/2011</b>	Analyst: <b>RM</b>
<b>DRO (C10-C28)</b>	<b>110</b>		<b>5.2</b>	<b>mg/Kg-dry</b>	<b>1</b>	9/22/2011 06:05 PM
<i>Surr: 4-Terphenyl-d14</i>	<i>70.4</i>		<i>39-115</i>	<i>%REC</i>	<i>1</i>	9/22/2011 06:05 PM
<b>MOISTURE</b>			<b>A2540 G</b>			Analyst: <b>CG</b>
<b>Moisture</b>	<b>22</b>		<b>0.050</b>	<b>% of sample</b>	<b>1</b>	9/21/2011 04:33 PM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

Client: HRL Compliance Solutions

# QC BATCH REPORT

Work Order: 1109654

Project: Williams TR 22-20-597 Pad LOE 9/19/11

Batch ID: 35736

Instrument ID GC8

Method: SW8015M

<b>MBLK</b>		Sample ID: <b>DBLKS1-35736-35736</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/22/2011 02:25 PM</b>		
Client ID:		Run ID: <b>GC8_110922A</b>				SeqNo: <b>1744437</b>		Prep Date: <b>9/21/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	ND	4.2								
Surr: 4-Terphenyl-d14	1.473	0	1.667	0	88.4	39-115	0			

<b>LCS</b>		Sample ID: <b>DLCSS1-35736-35736</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/22/2011 01:19 PM</b>		
Client ID:		Run ID: <b>GC8_110922A</b>				SeqNo: <b>1744435</b>		Prep Date: <b>9/21/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	147.1	4.2	166.7	0	88.3	60-130	0			
Surr: 4-Terphenyl-d14	1.243	0	1.667	0	74.6	39-115	0			

<b>LCSD</b>		Sample ID: <b>DLCSDS1-35736-35736</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/22/2011 01:19 PM</b>		
Client ID:		Run ID: <b>GC8_110922A</b>				SeqNo: <b>1744441</b>		Prep Date: <b>9/21/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	150.8	4.2	166.7	0	90.5	60-130	147.1	2.45	30	
Surr: 4-Terphenyl-d14	1.043	0	1.667	0	62.6	39-115	1.243	17.6	30	

<b>MS</b>		Sample ID: <b>1109625-21B MS</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/22/2011 01:41 PM</b>		
Client ID:		Run ID: <b>GC8_110922A</b>				SeqNo: <b>1744436</b>		Prep Date: <b>9/21/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	285.5	8.0	321.1	6.641	86.8	60-130	0			
Surr: 4-Terphenyl-d14	2.285	0	3.211	0	71.1	39-115	0			

<b>MSD</b>		Sample ID: <b>1109625-21B MSD</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/22/2011 01:41 PM</b>		
Client ID:		Run ID: <b>GC8_110922A</b>				SeqNo: <b>1744442</b>		Prep Date: <b>9/21/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	267.2	7.9	316.7	6.641	82.3	60-130	285.5	6.61	30	
Surr: 4-Terphenyl-d14	1.911	0	3.167	0	60.4	39-115	2.285	17.8	30	

The following samples were analyzed in this batch:

1109654-01A 1109654-02A 1109654-03A



**Client:** HRL Compliance Solutions  
**Work Order:** 1109654  
**Project:** Williams TR 22-20-597 Pad LOE 9/19/11

## QC BATCH REPORT

Batch ID: **R94973** Instrument ID **MOIST** Method: **A2540 G**

<b>MBLK</b>		Sample ID: <b>WBLKS1-R94973</b>				Units: % of sample			Analysis Date: <b>9/21/2011 04:33 PM</b>		
Client ID:		Run ID: <b>MOIST_110921E</b>				SeqNo: <b>1743850</b>			Prep Date:		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual	

Moisture ND 0.050

<b>LCS</b>		Sample ID: <b>LCS-R94973</b>				Units: % of sample			Analysis Date: <b>9/21/2011 04:33 PM</b>		
Client ID:		Run ID: <b>MOIST_110921E</b>				SeqNo: <b>1743849</b>			Prep Date:		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual	

Moisture 100 0.050 100 0 100 99.5-100.5 0

<b>DUP</b>		Sample ID: <b>1109618-03BDUP</b>				Units: % of sample			Analysis Date: <b>9/21/2011 04:33 PM</b>		
Client ID:		Run ID: <b>MOIST_110921E</b>				SeqNo: <b>1743824</b>			Prep Date:		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual	

Moisture 77.53 0.050 0 0 0 0-0 78.11 0.745 20

<b>DUP</b>		Sample ID: <b>1109623-03BDUP</b>				Units: % of sample			Analysis Date: <b>9/21/2011 04:33 PM</b>		
Client ID:		Run ID: <b>MOIST_110921E</b>				SeqNo: <b>1743829</b>			Prep Date:		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual	

Moisture 9.59 0.050 0 0 0 0-0 8.98 6.57 20

The following samples were analyzed in this batch:

1109654-01A	1109654-02A	1109654-03A
-------------	-------------	-------------

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.



# ALS Laboratory Group

225 Commerce Drive, Fort Collins, Colorado 80524  
TF: (800) 443-1511 PH: (970) 490-1511 FX: (970) 490-1522

## Chain-of-Custody

Form 202r8

WORKORDER  
#

1109654

PAGE

1 of 1

DISPOSAL

By Lab or Return to Client

SAMPLER

Reed Wold

DATE

9/20/2011

PROJECT NAME

Williams TR 22-20-597 Pad LOE

SITE ID

TR 22-20-597

TURNAROUND

Standard

PROJECT No.

EDD FORMAT

PURCHASE ORDER

COMPANY NAME

HRL Compliance

BILL TO COMPANY

Williams

SEND REPORT TO

Kris Rowe

INVOICE ATTN TO

Karolina Blaney

ADDRESS

744 Horizon Ct Ste. 140

ADDRESS

1058 Co. Rd. 215

CITY / STATE / ZIP

Grand Junction, CO 81506

CITY / STATE / ZIP

Parachute, CO 81635

PHONE

970-243-3271

PHONE

970-683-2295

FAX

970-243-3280

FAX

E-MAIL

Krowe@hrlcomp.com

E-MAIL

Karolina.blaney@williams.com

Lab ID

Field ID

Matrix

Sample  
Date

Sample  
Time

#  
Bottles

Pres.

QC

01

TR 22-20-597 N. Bottom

SO

9/19/2011

3:15

1

8

X

02

TR 22-20-597 S. Bottom

SO

9/19/2011

3:20

1

8

X

03

TR 22-20-597 S.Wall

SO

9/19/2011

3:25

1

8

X

\*Time Zone (Circle): EST CST MST PST Matrix: O = oil S = soil NS = non-soil solid W = water L = liquid E = extract F = filler

For metals or anions, please detail analytes below.

Comments:

JRP  
5.2°C

QC PACKAGE (check below)

X

LEVEL II (Standard QC)

LEVEL III (Std QC + forms)

LEVEL IV (Std QC + forms  
+ raw data)

Preservative Key: 1-HCl 2-HNO3 3-H2SO4 4-NaOH 5-NaHSO4 7-Other 8-4 degrees C 9-5035

SIGNATURE

PRINTED NAME

DATE

TIME

RELINQUISHED BY

Reed Wold

Reed Wold

9/20/11

4:30pm

RECEIVED BY

Alex J. (Sister)

Alex J. (Sister)

9/21/11

10:30

RELINQUISHED BY

RECEIVED BY

RELINQUISHED BY

RECEIVED BY

Sample Receipt Checklist

Client Name: **HRL**

Date/Time Received: **21-Sep-11 10:30**

Work Order: **1109654**

Received by: **AC**

Checklist completed by Alex Coaszar  
eSignature

21-Sep-11  
Date

Reviewed by: Ann Preston  
eSignature

22-Sep-11  
Date

Matrices: **Soil**

Carrier name: **FedEx**

Shipping container/cooler in good condition?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	Not Present <input type="checkbox"/>
Custody seals intact on shipping container/cooler?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	Not Present <input type="checkbox"/>
Custody seals intact on sample bottles?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	Not Present <input checked="" type="checkbox"/>
Chain of custody present?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Chain of custody signed when relinquished and received?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Chain of custody agrees with sample labels?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Samples in proper container/bottle?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Sample containers intact?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Sufficient sample volume for indicated test?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
All samples received within holding time?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Container/Temp Blank temperature in compliance?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Temperature(s)/Thermometer(s):	<u>5.2 degrees C</u>		
Cooler(s)/Kit(s):			
Water - VOA vials have zero headspace?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	No VOA vials submitted <input checked="" type="checkbox"/>
Water - pH acceptable upon receipt?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	N/A <input checked="" type="checkbox"/>
pH adjusted?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	N/A <input checked="" type="checkbox"/>
pH adjusted by:			
Login Notes:			

-----

Client Contacted:

Date Contacted:

Person Contacted:

Contacted By:

Regarding:

Comments:

CorrectiveAction:

**FedEx** Express **NEW Package**  
**US Airbill**

FedEx  
Tracking  
Number

8758 3471 3683

0200

Form  
10.10c

FedEx Retrieval Copy

**1 From**

Date 7/20/11 Sender's FedEx  
Account Number 1

Sender's  
Name Deed Delle

Phone 714 200 3071

Company MTI Compliance

Address 744 Harbor Ct Ste 114

City Long Beach State CA ZIP 90801

Dept./Floor/State/Room

**2 Your Internal Billing Reference**

**3 To**

Recipient's  
Name Sample Recipient

Phone 614 599 6070

Company MTI Group

Address 3150 12th Ave

We cannot deliver to P.O. boxes or P.O. ZIP codes

Dept./Floor/State/Room

Address

Use this line for the HOLD location address or for continuation of your shipping address.

City Hanford State MT ZIP 49424

**01** **HOLD Weekday**  
FedEx location address  
REQUIRED NOT available for  
FedEx First Overnight

**31** **HOLD Saturday**  
FedEx location address  
REQUIRED Available ONLY for  
FedEx Priority Overnight and  
FedEx 2Day to select locations

**4 Express Package Service**

\*To most locations

NOTE: Service order has changed. Please select carefully.

Packages up to 150 lbs.  
For packages over 150 lbs., use a different  
FedEx service. Foreign US Airmail

**06** **FedEx First Overnight**  
Earliest next business morning delivery to select  
locations. Friday shipments will be delivered on  
Monday unless SATURDAY Delivery is selected

**01** **FedEx Priority Overnight**  
Next business morning delivery to select  
locations. Friday shipments will be  
delivered on Monday unless SATURDAY Delivery is  
selected

**05** **FedEx Standard Overnight**  
Next business afternoon  
Saturday Delivery, FedEx ground

**49** **NEW FedEx 2Day A.M.**  
Second business morning  
Saturday - Third day, 7:00 A.M. delivery

**03** **FedEx 2Day**  
Second business morning delivery to select  
locations. Friday shipments will be  
delivered on Monday unless SATURDAY  
Delivery is selected

**20** **FedEx Express Saver**  
Third business morning  
Saturday - Third day, 12:00 P.M. delivery

**5 Packaging**

\*Declared value limit \$500

**06** **FedEx Envelope** **02** **FedEx Pak** **03** **FedEx Box** **04** **FedEx Tube** **11** **Other**

**6 Special Handling and Delivery Signature Options**

**03 SATURDAY DELIVERY**

**No Signature Required**  
Package may be left without  
signature for delivery

**Direct Signature**  
Signature of shipper, addressee,  
or other designated person  
required for delivery. Fee applies

**Indirect Signature**  
If not this is a delivery to recipient  
address, addressee's signature  
not required. Addressee may sign for delivery. If  
no delivery address, recipient's signature required

Does this shipment contain dangerous goods?

One box must be checked

**No** **04** **Yes**

As per attached  
Shipping Declaration

**Yes**

Shipper's declaration  
not required

**06** **Dry Ice**

Dry Ice, 9, 10, 105

**Cargo Aircraft Only**

**7 Payment Bill to:**

Enter FedEx Acct. No. or Credit Card No. below

Payment type:  
Acct. No.

**1** **Sender** **2** **Recipient** **3** **Third Party** **4** **Credit Card** **5** **Cash/Check**

Total Packages

Total Weight

Check Care Arch

If an invoice is issued to you in 300 orders you receive a higher value. See the current FedEx Service Guide for details

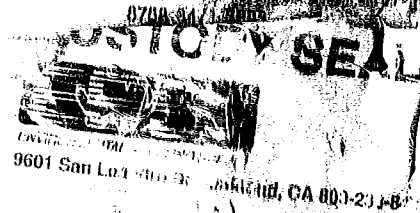
Dom. Rate 11/10 • Post #103120 • ©1994-2010 FedEx • PRINTED IN U.S.A. SBY

Date: 7/20/11

Signature: Reed Delle

fedex.com 1.800.GoFedEx 1.800.463.3339

fedex.com 1.800.GoFedEx 1.800.463.3339



5.2°C

### **Appendix 3: Background Raw Analytical Data**



08-Sep-2011

Kris Rowe  
HRL Compliance Solutions  
744 Horizon Ct. Suite 140  
Grand Junction, CO 81506

Re: **Williams TR 22-20-597 Pad LOE 8/24/11**

Work Order: **1108948**

Dear Kris,

ALS Environmental received 3 samples on 31-Aug-2011 09:00 AM for the analyses presented in the following report.

The analytical data provided relates directly to the samples received by ALS Environmental and for only the analyses requested.

QC sample results for this data met laboratory specifications. Any exceptions are noted in the Case Narrative, or noted with qualifiers in the report or QC batch information. Should this laboratory report need to be reproduced, it should be reproduced in full unless written approval has been obtained from ALS Environmental. Samples will be disposed in 30 days unless storage arrangements are made.

The total number of pages in this report is 17.

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

Sincerely,

A handwritten signature in cursive script that reads "Ann Preston".

Electronically approved by: Ann Preston

Ann Preston  
Project Manager



Certificate No: IL100452

ADDRESS 3352 128th Avenue Holland, Michigan 49424-9263 | PHONE (616) 399-6070 | FAX (616) 399-6185

ALS GROUP USA, CORP Part of the ALS Laboratory Group A Campbell Brothers Limited Company

Environmental

[www.alsglobal.com](http://www.alsglobal.com)

RIGHT SOLUTIONS RIGHT PARTNER

---

**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11  
**Work Order:** 1108948

---

**Work Order Sample Summary**

---

<u>Lab Samp ID</u>	<u>Client Sample ID</u>	<u>Matrix</u>	<u>Tag Number</u>	<u>Collection Date</u>	<u>Date Received</u>	<u>Hold</u>
1108948-01	TR 22-20-597 BG 1	Soil		8/24/2011 13:15	8/31/2011 09:00	<input type="checkbox"/>
1108948-02	TR 22-20-597 BG 2	Soil		8/24/2011 13:20	8/31/2011 09:00	<input type="checkbox"/>
1108948-03	TR 22-20-597 BG 3	Soil		8/24/2011 13:25	8/31/2011 09:00	<input type="checkbox"/>

---

## ALS Group USA, Corp

*Date: 08-Sep-11*

---

**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11  
**Work Order:** 1108948

---

### Case Narrative

Batch 35255 MS/MSD data for Metals is not related to this project's samples.



**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11  
**WorkOrder:** 1108948

**QUALIFIERS,  
ACRONYMS, UNITS**

<b><u>Qualifier</u></b>	<b><u>Description</u></b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL

<b><u>Acronym</u></b>	<b><u>Description</u></b>
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
MQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
SD	Serial Dilution
TDL	Target Detection Limit

<b><u>Units Reported</u></b>	<b><u>Description</u></b>
% of sample as noted	Percent of Sample
mg/Kg-dry s.u.	Milligrams per Kilogram Dry Weight Standard Units

**ALS Group USA, Corp****Date:** 08-Sep-11**Client:** HRL Compliance Solutions**Project:** Williams TR 22-20-597 Pad LOE 8/24/11**Work Order:** 1108948**Sample ID:** TR 22-20-597 BG 1**Lab ID:** 1108948-01**Collection Date:** 8/24/2011 01:15 PM**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<b>METALS BY ICP-MS</b>			<b>SW6020A</b>		Prep Date: <b>8/31/2011</b>	Analyst: <b>CES</b>
Arsenic	5.4		0.95	mg/Kg-dry	2	9/1/2011 11:04 PM
<b>SUBCONTRACTED ANALYSES</b>			<b>SUBCONTRACT</b>			Analyst: <b>A&amp;LGL</b>
Subcontracted Analyses	Rcvd 9/6/11		as noted		1	9/6/2011
<b>MOISTURE</b>			<b>A2540 G</b>			Analyst: <b>CG</b>
Moisture	21		0.050	% of sample	1	8/31/2011 01:03 PM
<b>PH</b>			<b>SW9045D</b>			Analyst: <b>JJG</b>
pH	7.02	H		s.u.	1	8/31/2011 08:30 AM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

**ALS Group USA, Corp****Date:** 08-Sep-11**Client:** HRL Compliance Solutions**Project:** Williams TR 22-20-597 Pad LOE 8/24/11**Work Order:** 1108948**Sample ID:** TR 22-20-597 BG 2**Lab ID:** 1108948-02**Collection Date:** 8/24/2011 01:20 PM**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<b>METALS BY ICP-MS</b>			<b>SW6020A</b>		Prep Date: <b>8/31/2011</b>	Analyst: <b>CES</b>
Arsenic	5.2		0.84	mg/Kg-dry	2	9/1/2011 11:10 PM
<b>MOISTURE</b>			<b>A2540 G</b>			Analyst: <b>CG</b>
Moisture	6.2		0.050	% of sample	1	8/31/2011 01:03 PM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

**ALS Group USA, Corp****Date:** 08-Sep-11**Client:** HRL Compliance Solutions**Project:** Williams TR 22-20-597 Pad LOE 8/24/11**Work Order:** 1108948**Sample ID:** TR 22-20-597 BG 3**Lab ID:** 1108948-03**Collection Date:** 8/24/2011 01:25 PM**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<b>METALS BY ICP-MS</b>			<b>SW6020A</b>		Prep Date: <b>9/1/2011</b>	Analyst: <b>CES</b>
Arsenic	5.7		0.73	mg/Kg-dry	2	9/6/2011 05:44 PM
<b>MOISTURE</b>			<b>A2540 G</b>			Analyst: <b>CG</b>
Moisture	7.0		0.050	% of sample	1	8/31/2011 01:03 PM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

Report Number: F11244-0175

Account Number: 91000

# A & L GREAT LAKES LABORATORIES, INC.

3505 Conestoga Drive • Fort Wayne, Indiana 46808-4413 • Phone 260-483-4759 • Fax 260-483-5274

www.algreatlakes.com • lab@algreatlakes.com



**QUALITY ANALYSES FOR INFORMED DECISIONS**

TO: ALS LABORATORY GROUP  
3352 128TH AVE  
HOLLAND, MI 49424-9263

RE: 1108948

DATE RECEIVED: 09/01/2011

DATE REPORTED: 09/06/2011

PAGE: 1

P.O. NUMBER: 20-122010675

ATTN: ANN PRESTON

## REPORT OF ANALYSIS

LAB NO.	SAMPLE ID	ANALYSIS	RESULT	UNIT	METHOD
82828	01B	Sat'd Paste Extraction with DIW	1		USDA Handbook 60
		Conductivity (ECe)	3.01	mmho/cm	USDA Handbook 60
		Calcium (Sat'd Paste)	275	ppm	USDA Handbook 60
		Magnesium (Sat'd Paste)	39	ppm	USDA Handbook 60
		Sodium (Sat'd Paste)	3137	ppm	USDA Handbook 60
		Sodium Adsorption Ratio	46.8	-	USDA Handbook 60

Client: HRL Compliance Solutions

Work Order: 1108948

Project: Williams TR 22-20-597 Pad LOE 8/24/11

# QC BATCH REPORT

Batch ID: **35255** Instrument ID **ICPMS1** Method: **SW6020A**

<b>MBLK</b>	Sample ID: <b>MBLK-35255-35255</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/1/2011 07:48 PM</b>			
Client ID:	Run ID: <b>ICPMS1_110901A</b>				SeqNo: <b>1723857</b>		Prep Date: <b>8/31/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	ND	0.25								

<b>LCS</b>	Sample ID: <b>LCS-35255-35255</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/1/2011 07:53 PM</b>			
Client ID:	Run ID: <b>ICPMS1_110901A</b>				SeqNo: <b>1723858</b>		Prep Date: <b>8/31/2011</b>		DF: <b>2</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	4.757	0.50	5	0	95.1	80-120	0			

<b>LCSD</b>	Sample ID: <b>LCSD-35255-35255</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/1/2011 08:20 PM</b>			
Client ID:	Run ID: <b>ICPMS1_110901A</b>				SeqNo: <b>1723861</b>		Prep Date: <b>8/31/2011</b>		DF: <b>2</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	4.786	0.50	5	0	95.7	80-120	4.757	0.608	20	

<b>MS</b>	Sample ID: <b>1108926-21BMS</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/1/2011 08:41 PM</b>			
Client ID:	Run ID: <b>ICPMS1_110901A</b>				SeqNo: <b>1723865</b>		Prep Date: <b>8/31/2011</b>		DF: <b>4</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	12.04	1.4	7.123	7.383	65.4	80-120	0			S

<b>MS</b>	Sample ID: <b>1108926-27BMS</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/1/2011 09:55 PM</b>			
Client ID:	Run ID: <b>ICPMS1_110901A</b>				SeqNo: <b>1723877</b>		Prep Date: <b>8/31/2011</b>		DF: <b>4</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	14.1	1.6	7.924	7.02	89.3	80-120	0			

<b>MSD</b>	Sample ID: <b>1108926-21BMSD</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/1/2011 08:46 PM</b>			
Client ID:	Run ID: <b>ICPMS1_110901A</b>				SeqNo: <b>1723866</b>		Prep Date: <b>8/31/2011</b>		DF: <b>4</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	12.02	1.4	6.993	7.383	66.3	80-120	12.04	0.169	25	S

<b>MSD</b>	Sample ID: <b>1108926-27BMSD</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/1/2011 10:01 PM</b>			
Client ID:	Run ID: <b>ICPMS1_110901A</b>				SeqNo: <b>1723878</b>		Prep Date: <b>8/31/2011</b>		DF: <b>4</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	13.98	1.6	7.752	7.02	89.8	80-120	14.1	0.855	25	

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108948  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

---

Batch ID: **35255** Instrument ID **ICPMS1** Method: **SW6020A**

---

**The following samples were analyzed in this batch:**

1108948-01A	1108948-02A
-------------	-------------

---

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108948  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **35276**      Instrument ID **ICPMS1**      Method: **SW6020A**

<b>MBLK</b>	Sample ID: <b>MBLK-35276-35276</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/6/2011 02:22 PM</b>			
Client ID:	Run ID: <b>ICPMS1_110906A</b>				SeqNo: <b>1725875</b>		Prep Date: <b>9/1/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	ND	0.25								

<b>LCS</b>	Sample ID: <b>LCS-35276-35276</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/6/2011 02:28 PM</b>			
Client ID:	Run ID: <b>ICPMS1_110906A</b>				SeqNo: <b>1725876</b>		Prep Date: <b>9/1/2011</b>		DF: <b>2</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	4.617	0.50	5	0	92.3	80-120	0			

<b>LCSD</b>	Sample ID: <b>LCSD-35276-35276</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/6/2011 02:54 PM</b>			
Client ID:	Run ID: <b>ICPMS1_110906A</b>				SeqNo: <b>1726029</b>		Prep Date: <b>9/1/2011</b>		DF: <b>2</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	4.606	0.50	5	0	92.1	80-120	4.617	0.239	20	

<b>MS</b>	Sample ID: <b>1108982-28BMS</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/6/2011 03:36 PM</b>			
Client ID:	Run ID: <b>ICPMS1_110906A</b>				SeqNo: <b>1726037</b>		Prep Date: <b>9/1/2011</b>		DF: <b>4</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	14.63	1.4	6.897	8.826	84.2	80-120	0			

<b>MSD</b>	Sample ID: <b>1108982-28BMSD</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>9/6/2011 03:41 PM</b>			
Client ID:	Run ID: <b>ICPMS1_110906A</b>				SeqNo: <b>1726038</b>		Prep Date: <b>9/1/2011</b>		DF: <b>4</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	15.41	1.3	6.64	8.826	99.1	80-120	14.63	5.15	25	

The following samples were analyzed in this batch:

1108948-03A

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.



**Client:** HRL Compliance Solutions  
**Work Order:** 1108948  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **R94122** Instrument ID **WETCHEM** Method: **SW9045D**

<b>DUP</b>	Sample ID: <b>1108840-01A DUP</b>				Units: <b>s.u.</b>			Analysis Date: <b>8/31/2011 08:30 AM</b>		
Client ID:	Run ID: <b>WETCHEM_110831E</b>				SeqNo: <b>1722332</b>			Prep Date:		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
pH	7.51	0	0	0	0	0-0	7.51	0	20	

<b>DUP</b>	Sample ID: <b>1108951-01D DUP</b>				Units: <b>s.u.</b>			Analysis Date: <b>8/31/2011 08:30 AM</b>		
Client ID:	Run ID: <b>WETCHEM_110831E</b>				SeqNo: <b>1722340</b>			Prep Date:		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
pH	8.32	0	0	0	0	0-0	8.32	0	20	

The following samples were analyzed in this batch:

1108948-01A

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Client:** HRL Compliance Solutions  
**Work Order:** 1108948  
**Project:** Williams TR 22-20-597 Pad LOE 8/24/11

## QC BATCH REPORT

Batch ID: **R94148**      Instrument ID **MOIST**      Method: **A2540 G**

<b>MBLK</b>	Sample ID: <b>WBLKS1-R94148</b>				Units: % of sample			Analysis Date: <b>8/31/2011 01:03 PM</b>		
Client ID:	Run ID: <b>MOIST_110831C</b>				SeqNo: <b>1723039</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Moisture	ND	0.050								

<b>LCS</b>	Sample ID: <b>LCS-R94148</b>				Units: % of sample			Analysis Date: <b>8/31/2011 01:03 PM</b>		
Client ID:	Run ID: <b>MOIST_110831C</b>				SeqNo: <b>1723038</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Moisture	100	0.050	100	0	100	99.5-100.5	0			

<b>DUP</b>	Sample ID: <b>1108926-07BDUP</b>				Units: % of sample			Analysis Date: <b>8/31/2011 01:03 PM</b>		
Client ID:	Run ID: <b>MOIST_110831C</b>				SeqNo: <b>1723023</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Moisture	10.85	0.050	0	0	0	0-0	9.48	13.5	20	

<b>DUP</b>	Sample ID: <b>1108926-16BDUP</b>				Units: % of sample			Analysis Date: <b>8/31/2011 01:03 PM</b>		
Client ID:	Run ID: <b>MOIST_110831C</b>				SeqNo: <b>1723033</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Moisture	9.2	0.050	0	0	0	0-0	7.85	15.8	20	

The following samples were analyzed in this batch:

1108948-01A	1108948-02A	1108948-03A
-------------	-------------	-------------

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.



# ALS Laboratory Group

225 Commerce Drive, Fort Collins, Colorado 80524  
TF: (800) 443-1511 PH: (970) 490-1511 FX: (970) 490-1522

## Chain-of-Custody

Form 202r8

WORKORDER  
#

1108948

PROJECT NAME		TR 22-20-597 Pad LOE		SAMPLER		Reed Wold		DATE		8/30/2011		PAGE		1 of 1	
PROJECT No.				SITE ID		TR 22-20-597		TURNAROUND		Standard		DISPOSAL		By Lab or Return to Client	
COMPANY NAME		HRL COMPLIANCE SOLUTIONS Inc.		BILL TO COMPANY		Williams		Arsenic SAR/EC/PH							
SEND REPORT TO		KRIS ROWE		INVOICE ATTN TO		Karolia Blaney									
ADDRESS		744 HORIZON CT SUITE 140		ADDRESS		1058 co rd 215									
CITY / STATE / ZIP		GRAND JUNCTION CO 81506		CITY / STATE / ZIP		Parachute CO 81635									
PHONE		970-243-3271		PHONE		970-683-2295									
FAX		970-243-3280		FAX		970-285-9573									
E-MAIL		Krowe@hrlcomp.com		E-MAIL		Karolia.blaney@williams.com									
Lab ID	Field ID	Matrix	Sample Date	Sample Time	# Bottles	Pres.	QC								
01	TR 22-20-597 BG 1	SO	8/24/2011	1:15	2	8		X	X						
02	TR 22-20-597 BG 2	SO	8/24/2011	1:20	1	8		X							
03	TR 22-20-597 BG 3	SO	8/24/2011	1:25	1	8		X							

\*Time Zone (Circle): EST CST MST PST Matrix: O = oil S = soil NS = non-soil solid W = water L = liquid E = extract F = filler

For metals or anions, please detail analytes below.

Comments:	QC PACKAGE (check below)	
	X	LEVEL II (Standard QC)
		LEVEL III (Std QC + forms)
		LEVEL IV (Std QC + forms + raw data)
Preservative Key: 1-HCl 2-HNO3 3-H2SO4 4-NaOH 5-NaHSO4 7-Other 8-4 degrees C 9-5035		

	SIGNATURE	PRINTED NAME	DATE	TIME
RELINQUISHED BY	<i>Reed Wold</i>	Reed Wold	8/31/11	5:30
RECEIVED BY	<i>Diane F. Shaw</i>	Diane F. Shaw	8/31/11	0900
RELINQUISHED BY				
RECEIVED BY				
RELINQUISHED BY				
RECEIVED BY				

**Subcontractor:**

A &amp; L Great Lakes Agricultural La

3505 Conestoga Dr

TEL: (260) 483-4759

FAX: (260) 483-5274

Acct #: 91000

Ft. Wayne, IN 46808

**CHAIN-OF-CUSTODY RECORD**

Page 1 of 1

Date: **31-Aug-11**COC ID: **3081**Due Date **07-Sep-11**

Environmental

Customer Information		Project Information		Parameter/Method Request for Analysis										
Purchase Order		Project Name	1108948	A Subcontracted Analyses (SUBCONTRACT)										
Work Order		Project Number		B										
Company Name	ALS Group USA, Corp	Bill To Company	ALS Group USA, Corp	C										
Send Report To	Ann Preston	Inv Attn	Accounts Payable	D										
Address	3352 128th Avenue	Address	3352 128th Avenue	E										
				F										
City/State/Zip	Holland, Michigan 49424-9263	City/State/Zip	Holland, Michigan 49424-9263	G										
Phone	(616) 399-6070	Phone	(616) 399-6070	H										
Fax	(616) 399-6185	Fax	(616) 399-6185	I										
eMail Address	ann.preston@alsglobal.com	eMail CC		J										
<b>Sample ID</b>	<b>Matrix</b>	<b>Collection Date 24hr</b>	<b>Bottle</b>	<b>A</b>	<b>B</b>	<b>C</b>	<b>D</b>	<b>E</b>	<b>F</b>	<b>G</b>	<b>H</b>	<b>I</b>	<b>J</b>	
1108948-01B (TR 22-20-597 BG 1)	Soil	24/Aug/2011 13:15	(1) MISC	X										

**Comments:**Please analyze for SAR-EC. Email results to Ann Preston.

Relinquished by:	Date/Time	Received by:	Date/Time	Cooler IDs	Report/QC Level
					<b>Std</b>
Relinquished by:	Date/Time	Received by:	Date/Time		

Sample Receipt Checklist

Client Name: HRL

Date/Time Received: 31-Aug-11 09:00

Work Order: 1108948

Received by: DS

Checklist completed by Diane Shaw 31-Aug-11  
eSignature Date

Reviewed by: Ann Preston 02-Sep-11  
eSignature Date

Matrices: Soil

Carrier name: FedEx

Shipping container/cooler in good condition?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	Not Present <input type="checkbox"/>
Custody seals intact on shipping container/cooler?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	Not Present <input type="checkbox"/>
Custody seals intact on sample bottles?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	Not Present <input checked="" type="checkbox"/>
Chain of custody present?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Chain of custody signed when relinquished and received?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Chain of custody agrees with sample labels?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Samples in proper container/bottle?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Sample containers intact?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Sufficient sample volume for indicated test?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
All samples received within holding time?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Container/Temp Blank temperature in compliance?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Temperature(s)/Thermometer(s):	<u>3.2 c</u>		
Cooler(s)/Kit(s):			
Water - VOA vials have zero headspace?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	No VOA vials submitted <input checked="" type="checkbox"/>
Water - pH acceptable upon receipt?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	N/A <input checked="" type="checkbox"/>
pH adjusted?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	N/A <input checked="" type="checkbox"/>
pH adjusted by:			
Login Notes:			

-----

Client Contacted:

Date Contacted:

Person Contacted:

Contacted By:

Regarding:

Comments:

CorrectiveAction:

MAINTENANCE

DATE

SIGNATURE

QEC

Quality Environmental Containers  
800-266-3880 • 804-266-3880

**FedEx** **NEW YORK**  
**Express** **US Airbill**

Number

8758 3471 3889

0200

FedEx Retrieval Copy

1 From

Date

2/30/11

Sender's FedEx  
Account Number

Sender's  
Name

NEW YORK

Phone

212 466 1001

Company

1051

Address

1000 Avenue of the Americas

Dept./Floor/Suite/Room

City

New York

State

NY

ZIP

10001

2 Your Internal Billing Reference

3 To

Recipient's  
Name

People Resources

Phone

212 299 1000

Company

1051

Address

2352 1st Ave

Dept./Floor/Suite/Room

Address

Use this line for the HOLD location address or for sending it to your shipping address

City

New York

State

NY

ZIP

10001

HOLD Weekday  
FedEx location address  
REQUIRED. NOT available for  
FedEx First Overnight.

HOLD Saturday  
FedEx location address  
REQUIRED. Available ONLY for  
FedEx Priority Overnight and  
FedEx 2Day to select locations.

4 Express Package Service

For most locations,  
NEXT business day. Please select carefully.

001 FedEx First Overnight

Next business day morning delivery to select  
locations. Delivery times vary by location and  
package weight. SATURDAY delivery not available.

011 FedEx Priority Overnight

Next business day morning delivery to select  
locations. Delivery times vary by location and  
package weight. SATURDAY delivery not available.

051 FedEx Standard Overnight

Next business day morning delivery to select  
locations. SATURDAY delivery not available.

5 Packaging

\* Declared value limit \$500

061 FedEx Envelope

021 FedEx Pak

03 FedEx Box

04 FedEx Tube

011 FedEx Mailer

6 Special Handling and Delivery Signature Options

031 SATURDAY DELIVERY

No Signature Required  
Package may be left without  
obtaining a signature for delivery.

101 Direct Signature  
Signature at time of delivery.  
No return address required.

341 Indirect Signature  
Signature at time of delivery.  
Signature required at time of delivery.  
Signature required at time of delivery.

Does this shipment contain dangerous goods?

One box must be checked

No 041 Yes  
As per attached  
Shipper's Declaration

Yes  
Shipper's Declaration  
is required

Dangerous goods (including dry ice) cannot be shipped via FedEx Home Delivery.

Pay to  
Order of

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7 Payment Bill to:

Enter FedEx Acct. No. or Credit Card No. below

1 Sender  
Acct. No. in Section  
1 must be filled

2 Recipient

3 Third Party

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5 Cash/Check  
Acct. No.

Total Packages

101

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#### **Appendix 4: Treatment Cell Confirmation Raw Analytical Data**



15-Dec-2011

Kris Rowe  
HRL Compliance Solutions  
744 Horizon Ct. Suite 140  
Grand Junction, CO 81506

Re: **Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/11**

Work Order: **1111567**  
Revision: **1**

Dear Kris,

ALS Environmental received 2 samples on 16-Nov-2011 10:00 AM for the analyses presented in the following report.

This is a REVISED REPORT. The Case Narrative provides information discussing the reason for issuing a revised report. The total number of pages in this revision is 28.

If you have any questions regarding these test results, please feel free to contact me.

Sincerely,

A handwritten signature in cursive script that reads "Ann Preston".

Electronically approved by: Alex Cszaszar

Ann Preston  
Project Manager



Certificate No: IL100452

ADDRESS 3352 128th Avenue Holland, Michigan 49424-9263 | PHONE (616) 399-6070 | FAX (616) 399-6185

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Environmental 

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RIGHT SOLUTIONS RIGHT PARTNER



---

**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/11  
**Work Order:** 1111567

---

**Work Order Sample Summary**

---

<u>Lab Samp ID</u>	<u>Client Sample ID</u>	<u>Matrix</u>	<u>Tag Number</u>	<u>Collection Date</u>	<u>Date Received</u>	<u>Hold</u>
1111567-01	Treatment Cell North Side	Soil		11/8/2011 12:30	11/16/2011 10:00	<input type="checkbox"/>
1111567-02	Treatment Cell South Side	Soil		11/8/2011 12:35	11/16/2011 10:00	<input type="checkbox"/>

---

**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/11  
**Work Order:** 1111567

---

**Case Narrative**

Batch 37526 MS/MSD data for DRO is not related to this project's samples.

Batch 37537 MS/MSD data for Metals is not related to this project's samples.

Batch R97849 samples 1111567-01 and 1111567-02 samples for pH were received after the hold time had expired.

Batch R97862 samples 1111567-01 and 1111567-02 samples for % Moisture were received after the hold time had expired.

Revised report issued per client request. Project name and sample IDs were changed.

**Client:** HRL Compliance Solutions  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/11  
**WorkOrder:** 1111567

## **QUALIFIERS, ACRONYMS, UNITS**

<b><u>Qualifier</u></b>	<b><u>Description</u></b>
*	Value exceeds Regulatory Limit
a	Not accredited
B	Analyte detected in the associated Method Blank above the Reporting Limit
E	Value above quantitation range
H	Analyzed outside of Holding Time
J	Analyte detected below quantitation limit
n	Not offered for accreditation
ND	Not Detected at the Reporting Limit
O	Sample amount is > 4 times amount spiked
P	Dual Column results percent difference > 40%
R	RPD above laboratory control limit
S	Spike Recovery outside laboratory control limits
U	Analyzed but not detected above the MDL

<b><u>Acronym</u></b>	<b><u>Description</u></b>
DUP	Method Duplicate
LCS	Laboratory Control Sample
LCSD	Laboratory Control Sample Duplicate
MBLK	Method Blank
MDL	Method Detection Limit
SQL	Method Quantitation Limit
MS	Matrix Spike
MSD	Matrix Spike Duplicate
PDS	Post Digestion Spike
PQL	Practical Quantitation Limit
RPD	Relative Percent Difference
SD	Serial Dilution
TDL	Target Detection Limit

<b><u>Units Reported</u></b>	<b><u>Description</u></b>
% of sample	Percent of Sample
µg/Kg-dry as noted	Micrograms per Kilogram Dry Weight
mg/Kg-dry	Milligrams per Kilogram Dry Weight
s.u.	Standard Units

# ALS Group USA, Corp

Date: 15-Dec-11

**Client:** HRL Compliance Solutions

**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/11

**Work Order:** 1111567

**Sample ID:** Treatment Cell North Side

**Lab ID:** 1111567-01

**Collection Date:** 11/8/2011 12:30 PM

**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<b>DIESEL RANGE ORGANICS BY GC-FID</b>						
			<b>SW8015M</b>		Prep Date: <b>11/17/2011</b>	Analyst: <b>RM</b>
<b>DRO (C10-C28)</b>	<b>96</b>		<b>5.3</b>	<b>mg/Kg-dry</b>	1	11/18/2011 08:11 PM
Surr: 4-Terphenyl-d14	60.8		39-115	%REC	1	11/18/2011 08:11 PM
<b>GASOLINE RANGE ORGANICS BY GC-FID</b>						
			<b>SW8015</b>			Analyst: <b>RM</b>
GRO (C6-C10)	ND		6.5	mg/Kg-dry	100	11/21/2011 08:46 AM
Surr: Toluene-d8	112		50-150	%REC	100	11/21/2011 08:46 AM
<b>MERCURY BY CVAA</b>						
			<b>SW7471</b>		Prep Date: <b>11/18/2011</b>	Analyst: <b>LR</b>
Mercury	ND		0.026	mg/Kg-dry	1	11/21/2011 01:35 PM
<b>METALS BY ICP-MS</b>						
			<b>SW6020A</b>		Prep Date: <b>11/17/2011</b>	Analyst: <b>CES</b>
Arsenic	2.7		0.83	mg/Kg-dry	2	11/18/2011 08:49 PM
Barium	500		8.3	mg/Kg-dry	20	11/22/2011 02:54 AM
Cadmium	ND		0.33	mg/Kg-dry	2	11/18/2011 08:49 PM
Chromium	17		0.83	mg/Kg-dry	2	11/18/2011 08:49 PM
Copper	10		0.83	mg/Kg-dry	2	11/18/2011 08:49 PM
Lead	22		0.83	mg/Kg-dry	2	11/18/2011 08:49 PM
Nickel	12		0.83	mg/Kg-dry	2	11/18/2011 08:49 PM
Selenium	ND		0.83	mg/Kg-dry	2	11/18/2011 08:49 PM
Silver	ND		0.83	mg/Kg-dry	2	11/18/2011 08:49 PM
Zinc	48		1.7	mg/Kg-dry	2	11/18/2011 08:49 PM
<b>SUBCONTRACTED ANALYSES</b>						
Subcontracted Analyses	Rcvd 11/21/11		<b>SUBCONTRACT</b>			Analyst: <b>A&amp;LGL</b>
			as noted		1	11/21/2011
<b>SEMI-VOLATILE ORGANIC COMPOUNDS</b>						
			<b>SW8270</b>		Prep Date: <b>11/17/2011</b>	Analyst: <b>CW</b>
Acenaphthene	ND		38	µg/Kg-dry	1	11/20/2011 10:21 AM
Anthracene	ND		38	µg/Kg-dry	1	11/20/2011 10:21 AM
Benzo(a)anthracene	ND		38	µg/Kg-dry	1	11/20/2011 10:21 AM
Benzo(a)pyrene	ND		38	µg/Kg-dry	1	11/20/2011 10:21 AM
Benzo(b)fluoranthene	ND		38	µg/Kg-dry	1	11/20/2011 10:21 AM
Benzo(g,h,i)perylene	ND		38	µg/Kg-dry	1	11/20/2011 10:21 AM
Benzo(k)fluoranthene	ND		38	µg/Kg-dry	1	11/20/2011 10:21 AM
Chrysene	ND		38	µg/Kg-dry	1	11/20/2011 10:21 AM
Dibenzo(a,h)anthracene	ND		38	µg/Kg-dry	1	11/20/2011 10:21 AM
Fluoranthene	ND		38	µg/Kg-dry	1	11/20/2011 10:21 AM
Fluorene	ND		38	µg/Kg-dry	1	11/20/2011 10:21 AM
Indeno(1,2,3-cd)pyrene	ND		38	µg/Kg-dry	1	11/20/2011 10:21 AM
Naphthalene	ND		38	µg/Kg-dry	1	11/20/2011 10:21 AM
Pyrene	ND		38	µg/Kg-dry	1	11/20/2011 10:21 AM
Surr: 2,4,6-Tribromophenol	69.8		34-140	%REC	1	11/20/2011 10:21 AM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

Revision: 1

# ALS Group USA, Corp

Date: 15-Dec-11

**Client:** HRL Compliance Solutions

**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/11

**Work Order:** 1111567

**Sample ID:** Treatment Cell North Side

**Lab ID:** 1111567-01

**Collection Date:** 11/8/2011 12:30 PM

**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<i>Surr: 2-Fluorobiphenyl</i>	64.1		12-100	%REC	1	11/20/2011 10:21 AM
<i>Surr: 2-Fluorophenol</i>	72.9		33-117	%REC	1	11/20/2011 10:21 AM
<i>Surr: 4-Terphenyl-d14</i>	85.9		25-137	%REC	1	11/20/2011 10:21 AM
<i>Surr: Nitrobenzene-d5</i>	69.6		37-107	%REC	1	11/20/2011 10:21 AM
<i>Surr: Phenol-d6</i>	74.0		40-106	%REC	1	11/20/2011 10:21 AM
<b>VOLATILE ORGANIC COMPOUNDS</b>			<b>SW8260</b>			Analyst: <b>CW</b>
Benzene	ND		130	µg/Kg-dry	100	11/20/2011 05:48 AM
Ethylbenzene	ND		260	µg/Kg-dry	100	11/20/2011 05:48 AM
m,p-Xylene	ND		260	µg/Kg-dry	100	11/20/2011 05:48 AM
o-Xylene	ND		130	µg/Kg-dry	100	11/20/2011 05:48 AM
Toluene	ND		190	µg/Kg-dry	100	11/20/2011 05:48 AM
Xylenes, Total	ND		390	µg/Kg-dry	100	11/20/2011 05:48 AM
<i>Surr: 1,2-Dichloroethane-d4</i>	94.7		70-120	%REC	100	11/20/2011 05:48 AM
<i>Surr: 4-Bromofluorobenzene</i>	92.2		75-120	%REC	100	11/20/2011 05:48 AM
<i>Surr: Dibromofluoromethane</i>	93.5		85-115	%REC	100	11/20/2011 05:48 AM
<i>Surr: Toluene-d8</i>	106		85-115	%REC	100	11/20/2011 05:48 AM
<b>CHROMIUM, TRIVALENT</b>			<b>CALCULATION</b>			Analyst: <b>JJG</b>
Chromium, Trivalent	17		0.65	mg/Kg-dry	1	11/22/2011 03:20 PM
<b>CHROMIUM, HEXAVALENT</b>			<b>SW7196A</b>		Prep Date: 11/20/2011	Analyst: <b>MB</b>
Chromium, Hexavalent	ND		0.65	mg/Kg-dry	1	11/22/2011 01:00 PM
<b>MOISTURE</b>			<b>A2540 G</b>			Analyst: <b>CG</b>
Moisture	23	H	0.050	% of sample	1	11/16/2011 02:34 PM
<b>PH</b>			<b>SW9045D</b>			Analyst: <b>JJG</b>
pH	8.95	H		s.u.	1	11/17/2011 08:40 AM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

**Revision: 1**

# ALS Group USA, Corp

Date: 15-Dec-11

**Client:** HRL Compliance Solutions

**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/11

**Work Order:** 1111567

**Sample ID:** Treatment Cell South Side

**Lab ID:** 1111567-02

**Collection Date:** 11/8/2011 12:35 PM

**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<b>DIESEL RANGE ORGANICS BY GC-FID</b>						
			<b>SW8015M</b>		Prep Date: <b>11/17/2011</b>	Analyst: <b>RM</b>
<b>DRO (C10-C28)</b>	<b>47</b>		<b>5.0</b>	<b>mg/Kg-dry</b>	1	11/18/2011 03:07 PM
Surr: 4-Terphenyl-d14	67.7		39-115	%REC	1	11/18/2011 03:07 PM
<b>GASOLINE RANGE ORGANICS BY GC-FID</b>						
			<b>SW8015</b>			Analyst: <b>RM</b>
GRO (C6-C10)	ND		6.1	mg/Kg-dry	100	11/21/2011 09:11 AM
Surr: Toluene-d8	111		50-150	%REC	100	11/21/2011 09:11 AM
<b>MERCURY BY CVAA</b>						
			<b>SW7471</b>		Prep Date: <b>11/18/2011</b>	Analyst: <b>LR</b>
Mercury	ND		0.023	mg/Kg-dry	1	11/21/2011 01:37 PM
<b>METALS BY ICP-MS</b>						
			<b>SW6020A</b>		Prep Date: <b>11/17/2011</b>	Analyst: <b>CES</b>
Arsenic	3.9		0.99	mg/Kg-dry	2	11/18/2011 08:54 PM
Barium	540		9.9	mg/Kg-dry	20	11/22/2011 02:59 AM
Cadmium	ND		0.40	mg/Kg-dry	2	11/18/2011 08:54 PM
Chromium	16		0.99	mg/Kg-dry	2	11/18/2011 08:54 PM
Copper	10		0.99	mg/Kg-dry	2	11/18/2011 08:54 PM
Lead	20		0.99	mg/Kg-dry	2	11/18/2011 08:54 PM
Nickel	11		0.99	mg/Kg-dry	2	11/18/2011 08:54 PM
Selenium	ND		0.99	mg/Kg-dry	2	11/18/2011 08:54 PM
Silver	ND		0.99	mg/Kg-dry	2	11/18/2011 08:54 PM
Zinc	44		2.0	mg/Kg-dry	2	11/18/2011 08:54 PM
<b>SUBCONTRACTED ANALYSES</b>						
Subcontracted Analyses	Rcvd 11/21/11		<b>SUBCONTRACT</b>			Analyst: <b>A&amp;LGL</b>
			as noted		1	11/21/2011
<b>SEMI-VOLATILE ORGANIC COMPOUNDS</b>						
			<b>SW8270</b>		Prep Date: <b>11/17/2011</b>	Analyst: <b>CW</b>
Acenaphthene	ND		36	µg/Kg-dry	1	11/20/2011 10:56 AM
Anthracene	ND		36	µg/Kg-dry	1	11/20/2011 10:56 AM
Benzo(a)anthracene	ND		36	µg/Kg-dry	1	11/20/2011 10:56 AM
Benzo(a)pyrene	ND		36	µg/Kg-dry	1	11/20/2011 10:56 AM
Benzo(b)fluoranthene	ND		36	µg/Kg-dry	1	11/20/2011 10:56 AM
Benzo(g,h,i)perylene	ND		36	µg/Kg-dry	1	11/20/2011 10:56 AM
Benzo(k)fluoranthene	ND		36	µg/Kg-dry	1	11/20/2011 10:56 AM
Chrysene	ND		36	µg/Kg-dry	1	11/20/2011 10:56 AM
Dibenzo(a,h)anthracene	ND		36	µg/Kg-dry	1	11/20/2011 10:56 AM
Fluoranthene	ND		36	µg/Kg-dry	1	11/20/2011 10:56 AM
Fluorene	ND		36	µg/Kg-dry	1	11/20/2011 10:56 AM
Indeno(1,2,3-cd)pyrene	ND		36	µg/Kg-dry	1	11/20/2011 10:56 AM
Naphthalene	ND		36	µg/Kg-dry	1	11/20/2011 10:56 AM
Pyrene	ND		36	µg/Kg-dry	1	11/20/2011 10:56 AM
Surr: 2,4,6-Tribromophenol	74.1		34-140	%REC	1	11/20/2011 10:56 AM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

Revision: 1

# ALS Group USA, Corp

Date: 15-Dec-11

**Client:** HRL Compliance Solutions

**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/11

**Work Order:** 1111567

**Sample ID:** Treatment Cell South Side

**Lab ID:** 1111567-02

**Collection Date:** 11/8/2011 12:35 PM

**Matrix:** SOIL

Analyses	Result	Qual	Report Limit	Units	Dilution Factor	Date Analyzed
<i>Surr: 2-Fluorobiphenyl</i>	64.3		12-100	%REC	1	11/20/2011 10:56 AM
<i>Surr: 2-Fluorophenol</i>	69.9		33-117	%REC	1	11/20/2011 10:56 AM
<i>Surr: 4-Terphenyl-d14</i>	89.7		25-137	%REC	1	11/20/2011 10:56 AM
<i>Surr: Nitrobenzene-d5</i>	65.5		37-107	%REC	1	11/20/2011 10:56 AM
<i>Surr: Phenol-d6</i>	70.1		40-106	%REC	1	11/20/2011 10:56 AM
<b>VOLATILE ORGANIC COMPOUNDS</b>			<b>SW8260</b>			Analyst: <b>CW</b>
Benzene	ND		120	µg/Kg-dry	100	11/20/2011 06:12 AM
Ethylbenzene	ND		240	µg/Kg-dry	100	11/20/2011 06:12 AM
m,p-Xylene	ND		240	µg/Kg-dry	100	11/20/2011 06:12 AM
o-Xylene	ND		120	µg/Kg-dry	100	11/20/2011 06:12 AM
Toluene	ND		180	µg/Kg-dry	100	11/20/2011 06:12 AM
Xylenes, Total	ND		370	µg/Kg-dry	100	11/20/2011 06:12 AM
<i>Surr: 1,2-Dichloroethane-d4</i>	95.8		70-120	%REC	100	11/20/2011 06:12 AM
<i>Surr: 4-Bromofluorobenzene</i>	94.6		75-120	%REC	100	11/20/2011 06:12 AM
<i>Surr: Dibromofluoromethane</i>	96.4		85-115	%REC	100	11/20/2011 06:12 AM
<i>Surr: Toluene-d8</i>	100		85-115	%REC	100	11/20/2011 06:12 AM
<b>CHROMIUM, TRIVALENT</b>			<b>CALCULATION</b>			Analyst: <b>JJG</b>
Chromium, Trivalent	16		0.61	mg/Kg-dry	1	11/22/2011 03:20 PM
<b>CHROMIUM, HEXAVALENT</b>			<b>SW7196A</b>		Prep Date: 11/20/2011	Analyst: <b>MB</b>
Chromium, Hexavalent	ND		0.60	mg/Kg-dry	1	11/22/2011 01:00 PM
<b>MOISTURE</b>			<b>A2540 G</b>			Analyst: <b>CG</b>
Moisture	18	H	0.050	% of sample	1	11/16/2011 02:34 PM
<b>PH</b>			<b>SW9045D</b>			Analyst: <b>JJG</b>
pH	8.93	H		s.u.	1	11/17/2011 08:40 AM

**Note:** See Qualifiers page for a list of qualifiers and their definitions.

**Revision: 1**

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Report Number: F11321-0533

Account Number: 91000

# A & L GREAT LAKES LABORATORIES, INC.

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**QUALITY ANALYSES FOR INFORMED DECISIONS**

TO: ALS LABORATORY GROUP  
3352 128TH AVE  
HOLLAND, MI 49424-9263

RE: 1111567

DATE RECEIVED: 11/17/2011

DATE REPORTED: 11/21/2011

PAGE: 1

P.O. NUMBER: 20-122011243

ATTN: ANN PRESTON

## REPORT OF ANALYSIS

LAB NO.	SAMPLE ID	ANALYSIS	RESULT	UNIT	METHOD
67963	01C	Sat'd Paste Extraction with DIW	1		USDA Handbook 60
		Conductivity (ECe)	3.03	mmho/cm	USDA Handbook 60
		Calcium (Sat'd Paste)	103	ppm	USDA Handbook 60
		Magnesium (Sat'd Paste)	13	ppm	USDA Handbook 60
		Sodium (Sat'd Paste)	1229	ppm	USDA Handbook 60
		Sodium Adsorption Ratio (SAR)	30.3	-	USDA Handbook 60
67964	02C	Sat'd Paste Extraction with DIW	1		USDA Handbook 60
		Conductivity (ECe)	2.22	mmho/cm	USDA Handbook 60
		Calcium (Sat'd Paste)	59	ppm	USDA Handbook 60
		Magnesium (Sat'd Paste)	9	ppm	USDA Handbook 60
		Sodium (Sat'd Paste)	834	ppm	USDA Handbook 60
		Sodium Adsorption Ratio (SAR)	26.7	-	USDA Handbook 60



# ALS Group USA, Corp

Date: 15-Dec-11

**Client:** HRL Compliance Solutions

## QC BATCH REPORT

**Work Order:** 1111567

**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/1

Batch ID: **37526**

Instrument ID **GC8**

Method: **SW8015M**

<b>MBLK</b>		Sample ID: <b>DBLKS1-37526-37526</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>11/18/2011 05:54 PM</b>		
Client ID:		Run ID: <b>GC8_111118A</b>				SeqNo: <b>1830261</b>		Prep Date: <b>11/17/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	ND	4.2								
<i>Surr: 4-Terphenyl-d14</i>	1.086	0	1.667	0	65.1	39-115	0			

<b>LCS</b>		Sample ID: <b>DLCSS1-37526-37526</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>11/18/2011 04:46 PM</b>		
Client ID:		Run ID: <b>GC8_111118A</b>				SeqNo: <b>1830258</b>		Prep Date: <b>11/17/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	156	4.2	166.7	0	93.6	60-130	0			
<i>Surr: 4-Terphenyl-d14</i>	0.902	0	1.667	0	54.1	39-115	0			

<b>LCSD</b>		Sample ID: <b>DLCSDS1-37526-37526</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>11/18/2011 05:09 PM</b>		
Client ID:		Run ID: <b>GC8_111118A</b>				SeqNo: <b>1830237</b>		Prep Date: <b>11/17/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	137	4.2	166.7	0	82.2	60-130	156	12.9	30	
<i>Surr: 4-Terphenyl-d14</i>	0.9867	0	1.667	0	59.2	39-115	0.902	8.97	30	

<b>MS</b>		Sample ID: <b>1111465-01A MS</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>11/18/2011 05:09 PM</b>		
Client ID:		Run ID: <b>GC8_111118A</b>				SeqNo: <b>1830259</b>		Prep Date: <b>11/17/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	446.1	4.1	164.1	417	17.7	60-130	0			S
<i>Surr: 4-Terphenyl-d14</i>	1.239	0	1.641	0	75.5	39-115	0			

<b>MSD</b>		Sample ID: <b>1111465-01A MSD</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>11/18/2011 05:32 PM</b>		
Client ID:		Run ID: <b>GC8_111118A</b>				SeqNo: <b>1830238</b>		Prep Date: <b>11/17/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
DRO (C10-C28)	414.7	4.1	165.2	417	-1.42	60-130	446.1	7.3	30	S
<i>Surr: 4-Terphenyl-d14</i>	1.139	0	1.652	0	69	39-115	1.239	8.43	30	

The following samples were analyzed in this batch:

1111567-01B 1111567-02B

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Revision: 1**

QC Page: 1 of 15

**Client:** HRL Compliance Solutions  
**Work Order:** 1111567  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/1

## QC BATCH REPORT

Batch ID: **R98155** Instrument ID **GC9** Method: **SW8015**

<b>MBLK</b>	Sample ID: <b>MBLK-R98155-R98155</b>					Units: <b>µg/L</b>		Analysis Date: <b>11/21/2011 01:12 AM</b>		
Client ID:	Run ID: <b>GC9_111120A</b>				SeqNo: <b>1831029</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
GRO (C6-C10)	ND	200								
<i>Surr: Toluene-d8</i>	111.9	0	100	0	112	70-130	0			

<b>LCS</b>	Sample ID: <b>LCS-R98155-R98155</b>					Units: <b>µg/L</b>		Analysis Date: <b>11/20/2011 11:56 PM</b>		
Client ID:	Run ID: <b>GC9_111120A</b>				SeqNo: <b>1831028</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
GRO (C6-C10)	22740	200	25000	0	91	70-130	0			
<i>Surr: Toluene-d8</i>	92.23	0	100	0	92.2	70-130	0			

<b>LCSD</b>	Sample ID: <b>LCSD-R98155-R98155</b>					Units: <b>µg/L</b>		Analysis Date: <b>11/21/2011 12:21 PM</b>		
Client ID:	Run ID: <b>GC9_111120A</b>				SeqNo: <b>1831030</b>		Prep Date:		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
GRO (C6-C10)	22390	200	25000	0	89.6	70-130	22740	1.56	30	
<i>Surr: Toluene-d8</i>	96.89	0	100	0	96.9	70-130	92.23	4.93	30	

<b>MS</b>	Sample ID: <b>1111590-13A MS</b>					Units: <b>µg/Kg</b>		Analysis Date: <b>11/21/2011 10:02 AM</b>		
Client ID:	Run ID: <b>GC9_111120A</b>				SeqNo: <b>1831051</b>		Prep Date:		DF: <b>50</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
GRO (C6-C10)	1220000	2,500	1250000	0	97.6	70-130	0			
<i>Surr: Toluene-d8</i>	4619	0	5000	0	92.4	50-150	0			

<b>MSD</b>	Sample ID: <b>1111590-13A MSD</b>					Units: <b>µg/Kg</b>		Analysis Date: <b>11/21/2011 10:27 AM</b>		
Client ID:	Run ID: <b>GC9_111120A</b>				SeqNo: <b>1831052</b>		Prep Date:		DF: <b>50</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
GRO (C6-C10)	1201000	2,500	1250000	0	96.1	70-130	1220000	1.57	30	
<i>Surr: Toluene-d8</i>	4582	0	5000	0	91.6	50-150	4619	0.804	30	

The following samples were analyzed in this batch: | 1111567-01A | 1111567-02A |

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Revision: 1**

**Client:** HRL Compliance Solutions  
**Work Order:** 1111567  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/1

## QC BATCH REPORT

Batch ID: **37601**      Instrument ID **HG1**      Method: **SW7471**

<b>MBLK</b>	Sample ID: <b>MBLK-37601-37601</b>					Units: <b>mg/Kg</b>		Analysis Date: <b>11/21/2011 01:05 PM</b>		
Client ID:	Run ID: <b>HG1_111121A</b>				SeqNo: <b>1828453</b>		Prep Date: <b>11/18/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual

Mercury      ND      0.020

<b>LCS</b>	Sample ID: <b>LCS-37601-37601</b>					Units: <b>mg/Kg</b>		Analysis Date: <b>11/21/2011 01:13 PM</b>		
Client ID:	Run ID: <b>HG1_111121A</b>				SeqNo: <b>1828454</b>		Prep Date: <b>11/18/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual

Mercury      0.1586      0.020      0.1665      0      95.2      80-120      0

<b>LCSD</b>	Sample ID: <b>LCSD-37601-37601</b>					Units: <b>mg/Kg</b>		Analysis Date: <b>11/21/2011 01:15 PM</b>		
Client ID:	Run ID: <b>HG1_111121A</b>				SeqNo: <b>1828455</b>		Prep Date: <b>11/18/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual

Mercury      0.1491      0.020      0.1665      0      89.5      80-120      0.1586      6.18      20

<b>MS</b>	Sample ID: <b>1111558-02BMS</b>					Units: <b>mg/Kg</b>		Analysis Date: <b>11/21/2011 01:26 PM</b>		
Client ID:	Run ID: <b>HG1_111121A</b>				SeqNo: <b>1828460</b>		Prep Date: <b>11/18/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual

Mercury      0.1237      0.017      0.1382      0.01107      81.5      75-125      0

<b>MSD</b>	Sample ID: <b>1111558-02BMSD</b>					Units: <b>mg/Kg</b>		Analysis Date: <b>11/21/2011 01:28 PM</b>		
Client ID:	Run ID: <b>HG1_111121A</b>				SeqNo: <b>1828461</b>		Prep Date: <b>11/18/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual

Mercury      0.1219      0.016      0.1337      0.01107      82.9      75-125      0.1237      1.49      35

The following samples were analyzed in this batch:

1111567-01B      1111567-02B

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Revision: 1**

QC Page: 3 of 15

**Client:** HRL Compliance Solutions  
**Work Order:** 1111567  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/1

## QC BATCH REPORT

Batch ID: **37537** Instrument ID **ICPMS1** Method: **SW6020A**

<b>MBLK</b>		Sample ID: <b>MBLK-37537-37537</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>11/18/2011 07:50 PM</b>		
Client ID:		Run ID: <b>ICPMS1_111117A</b>				SeqNo: <b>1828236</b>		Prep Date: <b>11/17/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	ND	0.25								
Barium	ND	0.25								
Cadmium	ND	0.10								
Chromium	ND	0.25								
Copper	0.02006	0.25								J
Lead	ND	0.25								
Nickel	ND	0.25								
Selenium	ND	0.25								
Silver	ND	0.25								
Zinc	ND	0.50								

<b>LCS</b>		Sample ID: <b>LCS-37537-37537</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>11/18/2011 07:56 PM</b>		
Client ID:		Run ID: <b>ICPMS1_111117A</b>				SeqNo: <b>1828237</b>		Prep Date: <b>11/17/2011</b>		DF: <b>2</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	4.145	0.50	5	0	82.9	80-120	0			
Barium	4.558	0.50	5	0	91.2	80-120	0			
Cadmium	4.736	0.20	5	0	94.7	80-120	0			
Chromium	4.384	0.50	5	0	87.7	80-120	0			
Copper	4.382	0.50	5	0	87.6	80-120	0			
Lead	4.536	0.50	5	0	90.7	80-120	0			
Nickel	4.343	0.50	5	0	86.9	80-120	0			
Selenium	4.291	0.50	5	0	85.8	80-120	0			
Silver	4.21	0.50	5	0	84.2	80-120	0			
Zinc	4.678	1.0	5	0	93.6	80-120	0			

<b>LCSD</b>		Sample ID: <b>LCSD-37537-37537</b>				Units: <b>mg/Kg</b>		Analysis Date: <b>11/18/2011 08:01 PM</b>		
Client ID:		Run ID: <b>ICPMS1_111117A</b>				SeqNo: <b>1828238</b>		Prep Date: <b>11/17/2011</b>		DF: <b>2</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Arsenic	4.306	0.50	5	0	86.1	80-120	4.145	3.81	20	
Barium	4.527	0.50	5	0	90.5	80-120	4.558	0.682	20	
Cadmium	4.796	0.20	5	0	95.9	80-120	4.736	1.26	20	
Chromium	4.381	0.50	5	0	87.6	80-120	4.384	0.0685	20	
Copper	4.422	0.50	5	0	88.4	80-120	4.382	0.909	20	
Lead	4.602	0.50	5	0	92	80-120	4.536	1.44	20	
Nickel	4.362	0.50	5	0	87.2	80-120	4.343	0.437	20	
Selenium	4.457	0.50	5	0	89.1	80-120	4.291	3.8	20	
Silver	4.24	0.50	5	0	84.8	80-120	4.21	0.71	20	
Zinc	4.452	1.0	5	0	89	80-120	4.678	4.95	20	

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Revision: 1**

**Client:** HRL Compliance Solutions  
**Work Order:** 1111567  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/1

## QC BATCH REPORT

Batch ID: **37537**      Instrument ID **ICPMS1**      Method: **SW6020A**

MS					Sample ID: 1111569-04AMS		Units: mg/Kg		Analysis Date: 11/18/2011 09:25 PM		
Client ID:			Run ID: ICPMS1_111117A			SeqNo: 1828252		Prep Date: 11/17/2011		DF: 2	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual	
Arsenic	8.15	0.80	8	1.708	80.5	80-120	0				
Barium	673	0.80	8	507.7	2070	80-120	0			SEO	
Cadmium	7.17	0.32	8	0.1583	87.6	80-120	0				
Chromium	27.06	0.80	8	22.79	53.3	80-120	0			S	
Copper	12.67	0.80	8	7.653	62.7	80-120	0			S	
Lead	14.62	0.80	8	7.673	86.8	80-120	0				
Nickel	14.52	0.80	8	10.07	55.7	80-120	0			S	
Selenium	6.395	0.80	8	0.3981	75	80-120	0			S	
Silver	5.403	0.80	8	0.02799	67.2	80-120	0			S	
Zinc	35.2	1.6	8	31.22	49.8	80-120	0			S	

MSD					Sample ID: 1111569-04AMSD		Units:mg/Kg		Analysis Date: 11/18/2011 11:17 PM		
Client ID:			Run ID: ICPMS1_111117A			SeqNo:1828266		Prep Date: 11/17/2011		DF: 2	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual	
Arsenic	9.214	0.76	7.587	1.708	98.9	80-120	8.15	12.2	25		
Barium	488.6	0.76	7.587	507.7	-251	80-120	673	31.7	25	SREO	
Cadmium	6.954	0.30	7.587	0.1583	89.6	80-120	7.17	3.05	25		
Chromium	30.74	0.76	7.587	22.79	105	80-120	27.06	12.8	25		
Copper	14.44	0.76	7.587	7.653	89.4	80-120	12.67	13	25		
Lead	15.36	0.76	7.587	7.673	101	80-120	14.62	4.94	25		
Nickel	18.12	0.76	7.587	10.07	106	80-120	14.52	22	25		
Selenium	6.53	0.76	7.587	0.3981	80.8	80-120	6.395	2.08	25		
Silver	5.31	0.76	7.587	0.02799	69.6	80-120	5.403	1.75	25	S	
Zinc	39.7	1.5	7.587	31.22	112	80-120	35.2	12	25	O	

The following samples were analyzed in this batch:

1111567-01B      1111567-02B

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Revision: 1**

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**Client:** HRL Compliance Solutions  
**Work Order:** 1111567  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/1

## QC BATCH REPORT

Batch ID: **37525**      Instrument ID **SVMS7**      Method: **SW8270**

<b>MBLK</b>		Sample ID: <b>SBLKS1-37525-37525</b>				Units: <b>µg/Kg</b>		Analysis Date: <b>11/18/2011 10:08 AM</b>		
Client ID:		Run ID: <b>SVMS7_111118A</b>				SeqNo: <b>1826696</b>		Prep Date: <b>11/17/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Acenaphthene	ND	30								
Anthracene	ND	30								
Benzo(a)anthracene	ND	30								
Benzo(a)pyrene	ND	30								
Benzo(b)fluoranthene	ND	30								
Benzo(g,h,i)perylene	ND	30								
Benzo(k)fluoranthene	ND	30								
Chrysene	ND	30								
Dibenzo(a,h)anthracene	ND	30								
Fluoranthene	ND	30								
Fluorene	ND	30								
Indeno(1,2,3-cd)pyrene	ND	30								
Naphthalene	ND	30								
Pyrene	ND	30								
<hr/>										
Surr: 2,4,6-Tribromophenol	1338	0	1667	0	80.3	34-140		0		
Surr: 2-Fluorobiphenyl	1062	0	1667	0	63.7	12-100		0		
Surr: 2-Fluorophenol	1255	0	1667	0	75.3	33-117		0		
Surr: 4-Terphenyl-d14	1209	0	1667	0	72.5	25-137		0		
Surr: Nitrobenzene-d5	1171	0	1667	0	70.2	37-107		0		
Surr: Phenol-d6	1278	0	1667	0	76.7	40-106		0		

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Revision: 1**

**Client:** HRL Compliance Solutions  
**Work Order:** 1111567  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/1

## QC BATCH REPORT

Batch ID: **37525**      Instrument ID **SVMS7**      Method: **SW8270**

<b>LCS</b>		Sample ID: <b>SLCSS1-37525-37525</b>				Units: <b>µg/Kg</b>		Analysis Date: <b>11/18/2011 10:37 AM</b>		
Client ID:		Run ID: <b>SVMS7_111118A</b>				SeqNo: <b>1826697</b>		Prep Date: <b>11/17/2011</b>		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Acenaphthene	1026	30	1333	0	76.9	45-110	0			
Anthracene	1070	30	1333	0	80.3	55-105	0			
Benzo(a)anthracene	998	30	1333	0	74.9	50-110	0			
Benzo(a)pyrene	1100	30	1333	0	82.5	50-110	0			
Benzo(b)fluoranthene	968.3	30	1333	0	72.6	45-115	0			
Benzo(g,h,i)perylene	1128	30	1333	0	84.6	40-125	0			
Benzo(k)fluoranthene	1073	30	1333	0	80.5	45-115	0			
Chrysene	1031	30	1333	0	77.3	55-110	0			
Dibenzo(a,h)anthracene	1153	30	1333	0	86.5	40-125	0			
Fluoranthene	1040	30	1333	0	78	55-115	0			
Fluorene	1106	30	1333	0	82.9	50-110	0			
Indeno(1,2,3-cd)pyrene	1159	30	1333	0	86.9	40-120	0			
Naphthalene	1006	30	1333	0	75.5	40-105	0			
Pyrene	1128	30	1333	0	84.6	45-125	0			
Surr: 2,4,6-Tribromophenol	1375	0	1667	0	82.5	34-140	0			
Surr: 2-Fluorobiphenyl	1116	0	1667	0	66.9	12-100	0			
Surr: 2-Fluorophenol	1255	0	1667	0	75.3	33-117	0			
Surr: 4-Terphenyl-d14	1295	0	1667	0	77.7	25-137	0			
Surr: Nitrobenzene-d5	1271	0	1667	0	76.3	37-107	0			
Surr: Phenol-d6	1230	0	1667	0	73.8	40-106	0			

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Revision: 1**

**Client:** HRL Compliance Solutions  
**Work Order:** 1111567  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/1

## QC BATCH REPORT

Batch ID: **37525**      Instrument ID **SVMS7**      Method: **SW8270**

LCSD				Sample ID: SLCSDS1-37525-37525			Units: µg/Kg		Analysis Date: 11/18/2011 11:06 AM		
Client ID:			Run ID: SVMS7_111118A			SeqNo: 1826698		Prep Date: 11/17/2011		DF: 1	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual	
Acenaphthene	1032	30	1333	0	77.4	45-110	1026	0.583	25		
Anthracene	1076	30	1333	0	80.7	55-105	1070	0.59	25		
Benzo(a)anthracene	1007	30	1333	0	75.5	50-110	998	0.865	25		
Benzo(a)pyrene	1107	30	1333	0	83	50-110	1100	0.574	25		
Benzo(b)fluoranthene	1073	30	1333	0	80.5	45-115	968.3	10.3	25		
Benzo(g,h,i)perylene	1128	30	1333	0	84.6	40-125	1128	0.0295	25		
Benzo(k)fluoranthene	932	30	1333	0	69.9	45-115	1073	14.1	25		
Chrysene	1043	30	1333	0	78.2	55-110	1031	1.16	25		
Dibenzo(a,h)anthracene	1169	30	1333	0	87.7	40-125	1153	1.32	25		
Fluoranthene	1053	30	1333	0	79	55-115	1040	1.27	25		
Fluorene	1110	30	1333	0	83.3	50-110	1106	0.391	25		
Indeno(1,2,3-cd)pyrene	1167	30	1333	0	87.5	40-120	1159	0.659	25		
Naphthalene	1001	30	1333	0	75.1	40-105	1006	0.465	25		
Pyrene	1130	30	1333	0	84.7	45-125	1128	0.177	25		
Surr: 2,4,6-Tribromophenol	1419	0	1667	0	85.1	34-140	1375	3.15	40		
Surr: 2-Fluorobiphenyl	1130	0	1667	0	67.8	12-100	1116	1.28	40		
Surr: 2-Fluorophenol	1263	0	1667	0	75.8	33-117	1255	0.662	40		
Surr: 4-Terphenyl-d14	1312	0	1667	0	78.7	25-137	1295	1.25	40		
Surr: Nitrobenzene-d5	1265	0	1667	0	75.9	37-107	1271	0.447	40		
Surr: Phenol-d6	1258	0	1667	0	75.5	40-106	1230	2.28	40		

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Revision: 1**



**Client:** HRL Compliance Solutions  
**Work Order:** 1111567  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/1

## QC BATCH REPORT

Batch ID: **37525**      Instrument ID **SVMS7**      Method: **SW8270**

MS				Sample ID: <b>1111465-01A MS</b>			Units: <b>µg/Kg</b>		Analysis Date: <b>11/19/2011 10:46 PM</b>	
Client ID:				Run ID: <b>SVMS7_111119A</b>			SeqNo: <b>1827080</b>		Prep Date: <b>11/17/2011</b>	
							DF: <b>1</b>			
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Acenaphthene	965.6	30	1317	0	73.3	45-110	0			
Anthracene	916.2	30	1317	0	69.6	55-105	0			
Benzo(a)anthracene	998.2	30	1317	0	75.8	50-110	0			
Benzo(a)pyrene	991.6	30	1317	0	75.3	50-110	0			
Benzo(b)fluoranthene	1101	30	1317	0	83.6	45-115	0			
Benzo(g,h,i)perylene	1019	30	1317	0	77.4	40-125	0			
Benzo(k)fluoranthene	878.7	30	1317	0	66.7	45-115	0			
Chrysene	979.1	30	1317	0	74.4	55-110	0			
Dibenzo(a,h)anthracene	887.9	30	1317	0	67.4	40-125	0			
Fluoranthene	1009	30	1317	0	76.6	55-115	0			
Fluorene	999.9	30	1317	0	75.9	50-110	0			
Indeno(1,2,3-cd)pyrene	911.3	30	1317	0	69.2	40-120	0			
Naphthalene	1305	30	1317	137.1	88.7	40-105	0			
Pyrene	1057	30	1317	0	80.3	45-125	0			
Surr: 2,4,6-Tribromophenol	1191	0	1646	0	72.4	34-140	0			
Surr: 2-Fluorobiphenyl	997.2	0	1646	0	60.6	12-100	0			
Surr: 2-Fluorophenol	1026	0	1646	0	62.3	33-117	0			
Surr: 4-Terphenyl-d14	1252	0	1646	0	76.1	25-137	0			
Surr: Nitrobenzene-d5	1058	0	1646	0	64.3	37-107	0			
Surr: Phenol-d6	1071	0	1646	0	65.1	40-106	0			

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Revision: 1**

**Client:** HRL Compliance Solutions  
**Work Order:** 1111567  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/1

## QC BATCH REPORT

Batch ID: **37525**      Instrument ID **SVMS7**      Method: **SW8270**

MSD				Sample ID: 1111465-01A MSD			Units: µg/Kg		Analysis Date: 11/19/2011 11:16 PM		
Client ID:		Run ID: SVMS7_111119A			SeqNo: 1827081		Prep Date: 11/17/2011		DF: 1		
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual	
Acenaphthene	1043	30	1327	0	78.6	45-110	965.6	7.68	30		
Anthracene	941.5	30	1327	0	70.9	55-105	916.2	2.72	30		
Benzo(a)anthracene	1029	30	1327	0	77.6	50-110	998.2	3.08	30		
Benzo(a)pyrene	1031	30	1327	0	77.7	50-110	991.6	3.94	30		
Benzo(b)fluoranthene	1136	30	1327	0	85.6	45-115	1101	3.19	30		
Benzo(g,h,i)perylene	957.8	30	1327	0	72.2	40-125	1019	6.22	30		
Benzo(k)fluoranthene	942.5	30	1327	0	71	45-115	878.7	7.01	30		
Chrysene	1021	30	1327	0	76.9	55-110	979.1	4.14	30		
Dibenzo(a,h)anthracene	869.2	30	1327	0	65.5	40-125	887.9	2.13	30		
Fluoranthene	1060	30	1327	0	79.9	55-115	1009	4.99	30		
Fluorene	1064	30	1327	0	80.2	50-110	999.9	6.21	30		
Indeno(1,2,3-cd)pyrene	884.4	30	1327	0	66.6	40-120	911.3	2.99	30		
Naphthalene	1332	30	1327	137.1	90	40-105	1305	2.08	30		
Pyrene	1113	30	1327	0	83.9	45-125	1057	5.22	30		
Surr: 2,4,6-Tribromophenol	1226	0	1659	0	73.9	34-140	1191	2.85	40		
Surr: 2-Fluorobiphenyl	1097	0	1659	0	66.1	12-100	997.2	9.54	40		
Surr: 2-Fluorophenol	1141	0	1659	0	68.8	33-117	1026	10.7	40		
Surr: 4-Terphenyl-d14	1293	0	1659	0	77.9	25-137	1252	3.22	40		
Surr: Nitrobenzene-d5	1156	0	1659	0	69.7	37-107	1058	8.83	40		
Surr: Phenol-d6	1153	0	1659	0	69.5	40-106	1071	7.31	40		

The following samples were analyzed in this batch:      1111567-01B      1111567-02B

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Revision: 1**

**Client:** HRL Compliance Solutions  
**Work Order:** 1111567  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/1

## QC BATCH REPORT

Batch ID: **R97963**      Instrument ID **VMS9**      Method: **SW8260**

<b>MBLK</b>		Sample ID: <b>VBLKW2-111119-R97963</b>				Units: <b>µg/L</b>		Analysis Date: <b>11/20/2011 03:46 AM</b>		
Client ID:		Run ID: <b>VMS9_111119B</b>				SeqNo: <b>1827243</b>		Prep Date:		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Benzene	ND	1.0								
Ethylbenzene	ND	1.0								
m,p-Xylene	ND	2.0								
o-Xylene	ND	1.0								
Toluene	ND	1.0								
Xylenes, Total	ND	3.0								
Surr: 1,2-Dichloroethane-d4	95.32	0	100	0	95.3	70-120	0			
Surr: 4-Bromofluorobenzene	97.33	0	100	0	97.3	75-120	0			
Surr: Dibromofluoromethane	97.63	0	100	0	97.6	85-115	0			
Surr: Toluene-d8	95.35	0	100	0	95.4	85-120	0			

<b>LCS</b>		Sample ID: <b>VLCSW2-111119-R97963</b>				Units: <b>µg/L</b>		Analysis Date: <b>11/20/2011 02:32 AM</b>		
Client ID:		Run ID: <b>VMS9_111119B</b>				SeqNo: <b>1827241</b>		Prep Date:		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Benzene	22.51	1.0	20	0	113	80-120	0			
Ethylbenzene	22.53	1.0	20	0	113	75-125	0			
m,p-Xylene	45.18	2.0	40	0	113	75-130	0			
o-Xylene	22.33	1.0	20	0	112	80-120	0			
Toluene	22.43	1.0	20	0	112	75-120	0			
Xylenes, Total	67.51	3.0	60	0	113	75-130	0			
Surr: 1,2-Dichloroethane-d4	94.82	0	100	0	94.8	70-120	0			
Surr: 4-Bromofluorobenzene	95.12	0	100	0	95.1	75-120	0			
Surr: Dibromofluoromethane	95.49	0	100	0	95.5	85-115	0			
Surr: Toluene-d8	99.48	0	100	0	99.5	85-120	0			

<b>LCSD</b>		Sample ID: <b>VLCSDW2-111119-R97963</b>				Units: <b>µg/L</b>		Analysis Date: <b>11/20/2011 02:57 AM</b>		
Client ID:		Run ID: <b>VMS9_111119B</b>				SeqNo: <b>1827242</b>		Prep Date:		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual
Benzene	21.95	1.0	20	0	110	80-120	22.51	2.52	30	
Ethylbenzene	22.13	1.0	20	0	111	75-125	22.53	1.79	30	
m,p-Xylene	44.38	2.0	40	0	111	75-130	45.18	1.79	30	
o-Xylene	22.06	1.0	20	0	110	80-120	22.33	1.22	30	
Toluene	22.1	1.0	20	0	110	75-120	22.43	1.48	30	
Xylenes, Total	66.44	3.0	60	0	111	75-130	67.51	1.6	30	
Surr: 1,2-Dichloroethane-d4	94.69	0	100	0	94.7	70-120	94.82	0.137	30	
Surr: 4-Bromofluorobenzene	96.16	0	100	0	96.2	75-120	95.12	1.09	30	
Surr: Dibromofluoromethane	96.34	0	100	0	96.3	85-115	95.49	0.886	30	
Surr: Toluene-d8	99.55	0	100	0	99.6	85-120	99.48	0.0703	30	

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Revision: 1**

**Client:** HRL Compliance Solutions  
**Work Order:** 1111567  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/1

## QC BATCH REPORT

Batch ID: **R97963**      Instrument ID **VMS9**      Method: **SW8260**

MS				Sample ID: 1111525-02A MS				Units: µg/L			Analysis Date: 11/20/2011 12:19 PM			
Client ID:				Run ID: VMS9_111119B				SeqNo: 1827256			Prep Date:		DF: 1	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual				
Benzene	22.62	1.0	20	0	113	80-120	0							
Ethylbenzene	22.99	1.0	20	0	115	75-125	0							
m,p-Xylene	45.64	2.0	40	0	114	75-130	0							
o-Xylene	22.35	1.0	20	0	112	80-120	0							
Toluene	23.07	1.0	20	0	115	75-120	0							
Xylenes, Total	67.99	3.0	60	0	113	75-130	0							
Surr: 1,2-Dichloroethane-d4	95.65	0	100	0	95.6	70-120	0							
Surr: 4-Bromofluorobenzene	95.15	0	100	0	95.2	75-120	0							
Surr: Dibromofluoromethane	96.24	0	100	0	96.2	85-115	0							
Surr: Toluene-d8	100.6	0	100	0	101	85-120	0							

MSD				Sample ID: 1111525-02A MSD				Units: µg/L		Analysis Date: 11/20/2011 12:44 PM	
Client ID:			Run ID: VMS9_111119B			SeqNo: 1827257		Prep Date:		DF: 1	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual	
Benzene	21.69	1.0	20	0	108	80-120	22.62	4.2	30		
Ethylbenzene	21.8	1.0	20	0	109	75-125	22.99	5.31	30		
m,p-Xylene	43.3	2.0	40	0	108	75-130	45.64	5.26	30		
o-Xylene	21.44	1.0	20	0	107	80-120	22.35	4.16	30		
Toluene	21.74	1.0	20	0	109	75-120	23.07	5.94	30		
Xylenes, Total	64.74	3.0	60	0	108	75-130	67.99	4.9	30		
Surr: 1,2-Dichloroethane-d4	95.19	0	100	0	95.2	70-120	95.65	0.482	30		
Surr: 4-Bromofluorobenzene	96.45	0	100	0	96.4	75-120	95.15	1.36	30		
Surr: Dibromofluoromethane	96.29	0	100	0	96.3	85-115	96.24	0.0519	30		
Surr: Toluene-d8	99.67	0	100	0	99.7	85-120	100.6	0.949	30		

The following samples were analyzed in this batch:

1111567-01A      1111567-02A

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Revision: 1**

**Client:** HRL Compliance Solutions  
**Work Order:** 1111567  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/1

## QC BATCH REPORT

Batch ID: **37655**      Instrument ID **WETCHEM**      Method: **SW7196A**

<b>MBLK</b>	Sample ID: <b>MBLK-37655-37655</b>					Units: <b>mg/Kg</b>		Analysis Date: <b>11/22/2011 01:00 PM</b>		
Client ID:	Run ID: <b>WETCHEM_111122L</b>				SeqNo: <b>1830803</b>		Prep Date: <b>11/20/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual

Chromium, Hexavalent      ND      0.50

<b>LCS</b>	Sample ID: <b>LCS-37655-37655</b>					Units: <b>mg/Kg</b>		Analysis Date: <b>11/22/2011 01:00 PM</b>		
Client ID:	Run ID: <b>WETCHEM_111122L</b>				SeqNo: <b>1830804</b>		Prep Date: <b>11/20/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual

Chromium, Hexavalent      1.928      0.50      2      0      96.4      75-110      0

<b>LCSD</b>	Sample ID: <b>LCSD-37655-37655</b>					Units: <b>mg/Kg</b>		Analysis Date: <b>11/22/2011 01:00 PM</b>		
Client ID:	Run ID: <b>WETCHEM_111122L</b>				SeqNo: <b>1830817</b>		Prep Date: <b>11/20/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual

Chromium, Hexavalent      1.98      0.50      2      0      99      75-110      1.928      2.66      20

<b>MS</b>	Sample ID: <b>1111615-01A MS</b>					Units: <b>mg/Kg</b>		Analysis Date: <b>11/22/2011 01:00 PM</b>		
Client ID:	Run ID: <b>WETCHEM_111122L</b>				SeqNo: <b>1830814</b>		Prep Date: <b>11/20/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual

Chromium, Hexavalent      1.243      0.50      1.992      0      62.4      60-130      0

<b>MSD</b>	Sample ID: <b>1111615-01A MSD</b>					Units: <b>mg/Kg</b>		Analysis Date: <b>11/22/2011 01:00 PM</b>		
Client ID:	Run ID: <b>WETCHEM_111122L</b>				SeqNo: <b>1830815</b>		Prep Date: <b>11/20/2011</b>		DF: <b>1</b>	
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual

Chromium, Hexavalent      1.243      0.50      1.992      0      62.4      60-130      1.243      0      30

The following samples were analyzed in this batch:

1111567-01B      1111567-02B

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Revision: 1**

QC Page: 13 of 15

**Client:** HRL Compliance Solutions  
**Work Order:** 1111567  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/1

## QC BATCH REPORT

Batch ID: **R97849** Instrument ID **WETCHEM** Method: **SW9045D**

LCS				Sample ID: LCS-R97849-R97849				Units: s.u.			Analysis Date: 11/17/2011 08:40 AM			
Client ID:				Run ID: WETCHEM_111117A				SeqNo: 1823641			Prep Date:		DF: 1	
Analyte		Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual			
pH		4.39	0	4.4	0	99.8	90-110	0						

DUP		Sample ID: 1111557-01B DUP				Units: s.u.		Analysis Date: 11/17/2011 08:40 AM			
Client ID:		Run ID: WETCHEM_111117A			SeqNo: 1823643		Prep Date:		DF: 1		
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual	
pH	8.75	0	0	0	0	0-0	8.75	0	20	H	

The following samples were analyzed in this batch:

1111567-01B 1111567-02B

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Revision: 1**

QC Page: 14 of 15

**Client:** HRL Compliance Solutions  
**Work Order:** 1111567  
**Project:** Williams TR 22-20-597 Treat. Cell Pad LOE 11/8/1

## QC BATCH REPORT

Batch ID: **R97862**      Instrument ID **MOIST**      Method: **A2540 G**

<b>MBLK</b>	Sample ID: <b>WBLKS1-R97862</b>				Units: % of sample			Analysis Date: <b>11/16/2011 02:34 PM</b>		
Client ID:	Run ID: <b>MOIST_111116B</b>				SeqNo: <b>1823914</b>			Prep Date:		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual

Moisture      ND      0.050

<b>LCS</b>	Sample ID: <b>LCS-R97862</b>				Units: % of sample			Analysis Date: <b>11/16/2011 02:34 PM</b>		
Client ID:	Run ID: <b>MOIST_111116B</b>				SeqNo: <b>1823913</b>			Prep Date:		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual

Moisture      100      0.050      100      0      100      99.5-100.5      0

<b>DUP</b>	Sample ID: <b>1111558-02BDUP</b>				Units: % of sample			Analysis Date: <b>11/16/2011 02:34 PM</b>		
Client ID:	Run ID: <b>MOIST_111116B</b>				SeqNo: <b>1823897</b>			Prep Date:		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual

Moisture      14.84      0.050      0      0      0      0-0      14.8      0.27      20

<b>DUP</b>	Sample ID: <b>1111566-01ADUP</b>				Units: % of sample			Analysis Date: <b>11/16/2011 02:34 PM</b>		
Client ID:	Run ID: <b>MOIST_111116B</b>				SeqNo: <b>1823904</b>			Prep Date:		DF: <b>1</b>
Analyte	Result	PQL	SPK Val	SPK Ref Value	%REC	Control Limit	RPD Ref Value	%RPD	RPD Limit	Qual

Moisture      16.06      0.050      0      0      0      0-0      16.97      5.51      20      H

The following samples were analyzed in this batch:

1111567-01B      1111567-02B

**Note:** See Qualifiers Page for a list of Qualifiers and their explanation.

**Revision: 1**

QC Page: 15 of 15

**Subcontractor:**A & L Great Lakes Agricultural La  
3505 Conestoga Dr

TEL: (260) 483-4759

FAX: (260) 483-5274

Ft. Wayne, IN 46808

Acct #: 91000

**CHAIN-OF-CUSTODY RECORD**Date: **16-Nov-11**COC ID: **3309**Due Da **22-Nov-11**

Page 1 of 1

**Environmental**

Salesperson

Debbie Fazio

Customer Information		Project Information		Parameter/Method Request for Analysis												
Purchase Order	20-122011243	Project Name	1111567	A	Subcontracted Analyses (SUBCONTRACT) SAR-EC											
Work Order		Project Number		B												
Company Name	ALS Group USA, Corp	Bill To Company	ALS Group USA, Corp	C												
Send Report To	Ann Preston	Inv Attn	Accounts Payable	D												
Address	3352 128th Avenue	Address	3352 128th Avenue	E												
				F												
City/State/Zip	Holland, Michigan 49424-9263	City/State/Zip	Holland, Michigan 49424-9263	G												
Phone	(616) 399-6070	Phone	(616) 399-6070	H												
Fax	(616) 399-6185	Fax	(616) 399-6185	I												
eMail Address	ann.preston@alsglobal.com	eMail CC		J												
<b>Sample ID</b>	<b>Matrix</b>	<b>Collection Date 24hr</b>	<b>Bottle</b>	<b>A</b>	<b>B</b>	<b>C</b>	<b>D</b>	<b>E</b>	<b>F</b>	<b>G</b>	<b>H</b>	<b>I</b>	<b>J</b>			
1111567-01C	Soil	8/Nov/2011 12:30	(1) MISC	X												
1111567-02C	Soil	8/Nov/2011 12:35	(1) MISC	X												

**Comments:**Please analyze for SAR-EC. Email results to Ann Preston.

Relinquished by:

Date/Time

Received by:

Date/Time

Cooler IDs

Report/QC Level

Std

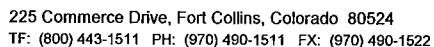
Relinquished by:

Date/Time

Received by:

Date/Time







1111567

1 of 1

By Lab or Return to Client

\*Time Zone (Circle): EST CST MST PST Matrix: O = oil S = soil NS = non-soil solid W = water L = liquid E = extract F = filler

**For metals or anions, please detail analytes below.**

	SIGNATURE	PRINTED NAME	DATE	TIME
RELINQUISHED BY		Paul W. H.	11/14/11	5:00
RECEIVED BY		Diane F. Shaw	11/16/11	1000
RELINQUISHED BY				
RECEIVED BY				
RELINQUISHED BY				
RECEIVED BY				

Sample Receipt Checklist

Client Name: HRL

Date/Time Received: 16-Nov-11 10:00

Work Order: 1111567

Received by: DS

Checklist completed by Diane Shaw 16-Nov-11  
eSignature Date

Reviewed by: Alex Csaszar 16-Nov-11  
eSignature Date

Matrices: Soil

Carrier name: FedEx

Shipping container/cooler in good condition?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	Not Present <input type="checkbox"/>
Custody seals intact on shipping container/cooler?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	Not Present <input type="checkbox"/>
Custody seals intact on sample bottles?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	Not Present <input checked="" type="checkbox"/>
Chain of custody present?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Chain of custody signed when relinquished and received?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Chain of custody agrees with sample labels?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Samples in proper container/bottle?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Sample containers intact?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Sufficient sample volume for indicated test?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
All samples received within holding time?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Container/Temp Blank temperature in compliance?	Yes <input checked="" type="checkbox"/>	No <input type="checkbox"/>	
Temperature(s)/Thermometer(s):	<u>5.8 c</u>		
Cooler(s)/Kit(s):			
Water - VOA vials have zero headspace?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	No VOA vials submitted <input checked="" type="checkbox"/>
Water - pH acceptable upon receipt?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	N/A <input checked="" type="checkbox"/>
pH adjusted?	Yes <input type="checkbox"/>	No <input type="checkbox"/>	N/A <input checked="" type="checkbox"/>
pH adjusted by:			

Login Notes:

-----

Client Contacted:

Date Contacted:

Person Contacted:

Contacted By:

Regarding:

Comments:

CorrectiveAction:

Revision: 1

CUSTODY SEAL

DATE

SIGNATURE

Quality Environmental Containers  
800-255-3950 • 304-255-3900

Express US AIRBILL

Tracking Number

8769 1479 5625

0200

FedEx Retrieval Copy

## 1 From

Date

11/14/11

Sender's FedEx  
Account NumberSender's  
Name

Bridgeway

Phone

72 242-3771

Company

HLS

Address

744 Harrison St. 14

City

Westerville

State

OH

ZIP 43086

## 2 Your Internal Billing Reference

## 3 To

Recipient's  
Name

SMPH Group

Phone

614 299-6072

Company

HLS Group

Address

3352 Locust Ave

Address

Use this line for the HOLD location address or for contribution of your shipping address.

City

Hillsdale

State

MI

ZIP

48124

HOLD Weekday  
FedEx location address  
REQUIRED. NOT available for  
FedEx First OvernightHOLD Saturday  
FedEx location address  
REQUIRED. Available ONLY for  
FedEx Priority Overnight and  
FedEx 2Day to select markets

## 4 Express Package Service

To select location.

NOTE: Service order has changed. Please select carefully.

Packages up to 150 lbs.  
For packages over 150 lbs., use the new  
FedEx Express Freight US Airbill.

06

FedEx First Overnight  
Earliest next business morning delivery to select  
locations. Friday shipments will be delivered on  
Monday unless SATURDAY Delivery is selected.

01

FedEx Priority Overnight  
Next business morning. Friday shipments will be  
delivered on Monday unless SATURDAY Delivery  
is selected.

05

FedEx Standard Overnight  
Next business day. Friday shipments will be  
delivered on Monday unless SATURDAY Delivery  
is selected.

## 5 Packaging

Additional value limit \$50.

05

FedEx envelope

06

FedEx Pak

03

FedEx  
Box

04

FedEx  
Tube

01

Other

## 6 Special Handling and Delivery Signature Options

## 03 SATURDAY DELIVERY

No Signature Required

Direct Signature

Indirect Signature

Does this shipment contain dangerous goods?

No (1)

Yes

Yes

Yes

Yes

Yes

Yes

Yes

Yes

Yes

Yes

Yes

Yes

Yes

Yes

## 7 Payment Bill to:

Sender

Account

Third Party

Credit Card

Cash/Check

Total Packages

Total Weight

Event Code

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8769 1479 5625

612

**Appendix 5: Sundry Notice Form 4 for Background Arsenic Consideration**

State of Colorado  
Oil and Gas Conservation Commission

1120 Lincoln Street, Suite 801, Denver, Colorado 80203 Phone: (303)894-2100 Fax: (303)894-2109



SUNDRY NOTICE

Submit original plus one copy. This form is to be used for general, technical and environmental sundry information. For proposed or completed operations, describe in full on Technical Information Page (Page 2 of this form.) Identify well or other facility by API Number or by OGCC Facility ID. Operator shall send an informational copy of all sundry notices for wells located in High Density Areas to the Local Government Designee (Rule 603b.)

1. OGCC Operator Number: 96850	4. Contact Name Karolina Blaney	Complete the Attachment Checklist  OP OGCC
2. Name of Operator: Williams Production RMT Company	Phone: 970-683-2295	
3. Address: 1058 County Road 215 City: Parachute State: CO Zip: 81635	Fax: 970-285-9573	
5. API Number 05- N/A	OGCC Facility ID Number 284697	Survey Plat
6. Well/Facility Name: Chevron TR 21-20-597	7. Well/Facility Number 21-20-597	Directional Survey
8. Location (Qtr/Sec, Twp, Rng, Meridian): SENW, Sec 20, TSS, R97W, 6th PM		Surface Eqmpt Diagram
9. County: Garfield	10. Field Name: Trail Ridge	Technical Info Page <input checked="" type="checkbox"/>
11. Federal, Indian or State Lease Number:		Other

General Notice

<input type="checkbox"/> CHANGE OF LOCATION: Attach New Survey Plat (a change of surface qtr/qtr is substantive and requires a new permit)	
Change of Surface Footage from Exterior Section Lines:	<input type="checkbox"/> FNLFSL <input type="checkbox"/> FELFWL
Change of Surface Footage to Exterior Section Lines:	<input type="checkbox"/>
Change of Bottomhole Footage from Exterior Section Lines:	<input type="checkbox"/>
Change of Bottomhole Footage to Exterior Section Lines:	<input type="checkbox"/> attach directional survey
Bottomhole location Qtr/Sec, Twp, Rng, Mer	
Latitude	Distance to nearest property line
Longitude	Distance to nearest bldg, public rd, utility or RR
Ground Elevation	Distance to nearest lease line
	Is location in a High Density Area (rule 603b)? Yes/No
	Distance to nearest well same formation
	Surface owner consultation date:
GPS DATA:	
Date of Measurement	PDOP Reading
	Instrument Operator's Name
<input type="checkbox"/> CHANGE SPACING UNIT	<input type="checkbox"/> Remove from surface bond
Formation	Formation Code
Spacing order number	Unit Acreage
Unit configuration	Signed surface use agreement attached
<input type="checkbox"/> CHANGE OF OPERATOR (prior to drilling):	<input type="checkbox"/> CHANGE WELL NAME
Effective Date:	From:
Plugging Bond: <input type="checkbox"/> Blanket <input type="checkbox"/> Individual	To:
	Effective Date:
<input type="checkbox"/> ABANDONED LOCATION:	<input type="checkbox"/> NOTICE OF CONTINUED SHUT IN STATUS
Was location ever built? <input type="checkbox"/> Yes <input type="checkbox"/> No	Date well shut in or temporarily abandoned:
Is site ready for inspection? <input type="checkbox"/> Yes <input type="checkbox"/> No	Has Production Equipment been removed from site? <input type="checkbox"/> Yes <input type="checkbox"/> No
Date Ready for Inspection:	MIT required if shut in longer than two years. Date of last MIT
<input type="checkbox"/> SPUD DATE:	<input type="checkbox"/> REQUEST FOR CONFIDENTIAL STATUS (6 mos from date casing set)
<input type="checkbox"/> SUBSEQUENT REPORT OF STAGE, SQUEEZE OR REMEDIAL CEMENT WORK	*submit cbl and cement job summaries
Method used	Cementing tool setting/perf depth
Cement volume	Cement top
Cement bottom	Date
<input type="checkbox"/> RECLAMATION: Attach technical page describing final reclamation procedures per Rule 1004.	
Final reclamation will commence on approximately	<input type="checkbox"/> Final reclamation is completed and site is ready for inspection.

Technical Engineering/Environmental Notice

<input type="checkbox"/> Notice of Intent	<input type="checkbox"/> Report of Work Done
Approximate Start Date:	Date Work Completed:
Details of work must be described in full on Technical Information Page (Page 2 must be submitted.)	
<input type="checkbox"/> Intent to Recomplete (submit form 2)	<input type="checkbox"/> Request to Vent or Flare
<input type="checkbox"/> Change Drilling Plans	<input type="checkbox"/> Repair Well
<input type="checkbox"/> Gross Interval Changed?	<input type="checkbox"/> Rule 502 variance requested
<input type="checkbox"/> Casing/Cementing Program Change	<input checked="" type="checkbox"/> Other: Background
<input type="checkbox"/> E&P Waste Disposal	<input type="checkbox"/> Beneficial Reuse of E&P Waste
<input type="checkbox"/> Status Update/Change of Remediation Plans	for Spills and Releases

I hereby certify that the statements made in this form are, to the best of my knowledge, true, correct and complete.

Signed: Karolina Blaney Date: 12/15/2011 Email: Karolina.Blaney@williams.com  
Print Name: Karolina Blaney Title: Environmental Specialist

COGCC Approved: *Chris Canfield* Title: FOR Date: 01/26/2012  
CONDITIONS OF APPROVAL, IF ANY: Chris Canfield  
EPS NW Region

REN # 5949

TECHNICAL INFORMATION PAGE



FOR OGCC USE ONLY

1. OGCC Operator Number: 96850 API Number: N/A
2. Name of Operator: Williams Production RMT OGCC Facility ID # 284697
3. Well/Facility Name: Chevron TR 21-20-597 Well/Facility Number: 21-20-597
4. Location (QtrQtr, Sec, Twp, Rng, Meridian): SENW, Sec 20, T5S, R97W, 6th PM

This form is to be completed whenever a Sundry Notice is submitted requiring detailed report of work to be performed or completed. This form shall be transmitted within 30 days of work completed as a "subsequent" report and must accompany Form 4, page 1.

5. **DESCRIBE PROPOSED OR COMPLETED OPERATIONS**

This COGCC Form 4 is being submitted as a request to consider the background concentration levels for arsenic at the Chevron TR 21-20-597 well pad relative to production pit closure at the subject facility in accordance with footnote 1 to the COGCC Table 9101-1.

The request is based on the analytical results below (see attached analytical)

Eight (6) grab samples were collected from locations within the pit footprint at depths of approximately 18' to 18.6' below pad grade to ascertain the arsenic concentrations of the facility.

North Pit Bottom - 2 mg/kg  
South Pit Bottom - 2.2 mg/kg  
East Wall - 2.8 mg/kg  
South Wall - East Half - 2.0 mg/kg  
West Wall - 4.1 mg/kg  
North Wall - 1.7 mg/kg

Average Concentration: 2.46 mg/kg

Three (3) grab samples were collected from nearby non-impacted, native soil from surface to 6" below to establish the background arsenic concentrations.

BKGD 1 - 5.4 mg/kg  
BKGD 2 - 5.2 mg/kg  
BKGD 3 - 5.7 mg/kg

Average Concentration: 5.43 mg/kg

Williams is requesting this approval in order to proceed with closure and reclamation of the production pit on the Chevron TR 21-20-597 well pad.