

DATA VALIDATION REPORT - Level III Review

SDG No.: 218081812+13+14 **Analysis:** Explosives and Metals
Laboratory: GCAL **Project:** Williston LTA
Reviewer: Naoum Tavantzis **Date:** October 24th, 2018

This report presents the findings of a review of the referenced data. The report consists of this summary, a listing of the samples included in the review, copies of data reports with data qualifying flags applied, data review worksheets, supporting documentation, and an explanation of the data qualifying flags employed. The review performed is based on the specifics of the analytical method referenced and provisions of the approved project-specific QAPP; and, qualified according to the USEPA CLP National Functional Guidelines for Organic and Inorganic (January 2017) Superfund Data Review. Modifications reflect the level of review requested, the specifications of the project-specific QAPP, and the specifics of the analytical methods employed.

Major

Anomalies: During the explosives analysis, the following laboratory control spike pairs (LCS/LCSD) displayed percent recoveries outside the laboratory quality control (QC) limits:

| Preparation Batch | QC Limits (%) | Analyte | LCS Recovery (%) | LCSD Recovery (%) |
|-------------------|---------------|----------------------------|------------------|-------------------|
| 642698 | 80-116 | 1,3,5-Trinitrobenzene | 47 | 41 |
| | 73-119 | 1,3-Dinitrobenzene | 50 | 43 |
| | 71-120 | 2,4,6-Trinitrobenzene | 42 | 39 |
| | 75-121 | 2,4-Dinitrotoluene | 26 | 32 |
| | 79-117 | 2,6-Dinitrotoluene | 105 | 56 |
| | 71-123 | 2-Amino-4,6-dinitrotoluene | 37 | 34 |
| | 84-120 | 2-Nitrotoluene | 44 | 35 |
| | 86-118 | 3,5-Dinitroaniline | 39 | 34 |
| | 67-129 | 3-Nitrotoluene | 42 | 35 |
| | 64-127 | 4-Amino-2,6-dinitrotoluene | 38 | 34 |
| | 71-124 | 4-Nitrotoluene | 47 | 60 |
| | 74-124 | HMX | 39 | 34 |
| | 80-128 | Nitrobenzene | 39 | 34 |
| | 73-124 | Nitroglycerin | 37 | 32 |
| | 72-128 | PETN | 31 | 30 |
| | 67-129 | RDX | 42 | 42 |
| 68-135 | Tetryl | 0 | 0 | |
| 643342 | 80-116 | 1,3,5-Trinitrobenzene | 121 | 119 |
| | 73-119 | 1,3-Dinitrobenzene | 126 | 111 |
| | 67-129 | 3-Nitrotoluene | 113 | 145 |
| | 71-124 | 4-Nitrotoluene | 108 | 354 |
| | 72-128 | PETN | 116 | 166 |
| | 68-135 | Tetryl | 0 | 0 |

The field sample results associated with the 0% recovery for tetryl were non-detect and were qualified R,I. The field sample results associated with the remaining percent recoveries less than the lower QC limits were non-detect and were qualified UJ,I. The field sample results associated with the positive biases were non-detect; no data qualifying action was taken. In addition, the LCS/LCSD displayed several relative percent difference (RPD) anomalies greater than the laboratory QC limit of 20%. The associated field sample results were non-detect; no data qualifying action was required.

Minor Anomalies:

During the explosives analysis, the LCSD prepared in QC batch 642698 displayed a percent recovery less than the lower QC limit of 50% for surrogate 1,2-dinitrobenzene at 45%. All field samples displayed surrogate percent recoveries within control limits; no data qualifying action is taken based on QC sample surrogate percent recovery anomalies. The matrix spike (MS) performed on field sample WIL02DA02A, prepared in QC batch 643342, displayed a percent recovery less than the lower QC limit of 74% for HMX at 61%. The associated parent sample result was non-detect and was qualified UJ,m. In addition, the matrix spike pairs (MS/MSD) performed on field sample WIL02DA02A in QC batches 642698 and 643342 displayed several RPDs greater than the QC laboratory QC limit of 20%. The associated parent sample results were non-detect; no data qualifying action was required. The field sample results associated with the percent recoveries less than the lower QC limits were re-extracted after the technical holding time of 14 days had expired. The associated field sample results were non-detect and were qualified UJ,h, unless previously qualified due to a LCS or MS percent recovery anomaly.

During the metals analysis, the method blank prepared in QC batch 642222 displayed a detection greater than the limit of detection for zinc at 518 µg/Kg. The associated field sample results were positive and were greater than five times the concentration found in the blank; no data qualifying action was required. The equipment blank WIL03IS00 displayed concentrations greater than the DL for antimony at 0.74 µg/L and copper at 0.26 µg/L. The associated field sample results were positive and were greater than five times the concentration found in the blank; no data qualifying action was required. The following MS/MSD displayed percent recoveries outside the quality QC limits:

| Parent Sample | Preparation Batch | Analyte | QC Limits (%) | MS Recovery (%) | MSD Recovery (%) |
|---------------|-------------------|----------|---------------|-----------------|------------------|
| WIL02DA02A | 642222 | Antimony | 72-124 | 7 | 5 |
| | | | | 11 | 12 |
| WIL01IS02 | 642442 | Copper | 84-119 | 132 | 115 |
| | 642531 | Antimony | 72-124 | 7 | 8 |

The post-digestion spikes performed on these parent samples displayed percent recoveries within laboratory QC limits. The field sample results associated with the positive bias were positive and were qualified J+,m. The field sample results associated with the negative biases were non-detect and were qualified UJ,m. The serial dilution performed on field sample WIL02DA02A in analytical sequence 642309 displayed a percent difference greater than the QC limit of 10% for zinc at 10.9%. The associated field sample results were positive and were qualified J,s.

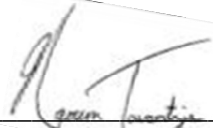
Correctable Anomalies:

None.

Comments:

On the basis of this evaluation, the laboratory appears to have followed the specified method, with the exception of anomalies discussed previously. If a given fraction was not discussed, all quality control criteria reviewed were within acceptable limits. Except for those data flagged "R", all data are usable, as qualified, for their intended purpose based on the data reviewed.

Signed:


 Naoum Tavantzis

Williston LTA

Job: 60520956

Laboratory: GCAL
SDG#: 217102040

| Sample ID | Client ID | Sample Type | Sample Date | Matrix | 8330B | Metals |
|-------------|------------|------------------|-------------|------------------|-------|--------|
| 21808181206 | WIL02DA02A | Field Sample | 8/16/2018 | Discrete Soil | x | x |
| 21808181213 | | | | | | |
| 21808181209 | WIL02DA01A | Field Sample | 8/16/2018 | Discrete Soil | x | x |
| 21808181216 | | | | | | |
| 21808181210 | WIL02DA01B | Field Duplicate | 8/16/2018 | Discrete Soil | x | x |
| 21808181217 | | | | | | |
| 21808181301 | WIL03IS01 | Field Sample | 8/14/2018 | Incremental Soil | | x |
| 21808181302 | WIL01IS02 | Field Duplicate | 8/14/2018 | Incremental Soil | | x |
| 21808181303 | WIL03IS00 | Equipment Blank | 8/16/2018 | Incremental Soil | | x |
| 21808181304 | WIL04IS03 | Field Triplicate | 8/15/2018 | Incremental Soil | | x |
| 21808181305 | WIL04IS01 | Field Sample | 8/15/2018 | Incremental Soil | | x |
| 21808181306 | WIL04IS02 | Field Duplicate | 8/15/2018 | Incremental Soil | | x |
| 21808181307 | WIL03IS03 | Field Triplicate | 8/14/2018 | Incremental Soil | | x |
| 21808181308 | WIL03IS02 | Field Duplicate | 8/14/2018 | Incremental Soil | | x |
| 21808181401 | WIL02IS01 | Field Sample | 8/15/2018 | Incremental Soil | x | x |
| 21808181404 | WIL02IS02 | Field Duplicate | 8/15/2018 | Incremental Soil | x | x |
| 21808181405 | WIL01IS03 | Field Triplicate | 8/14/2018 | Incremental Soil | | x |
| 21808181406 | WIL01IS01 | Field Sample | 8/14/2018 | Incremental Soil | | x |
| 21808181407 | WIL02IS03 | Field Triplicate | 8/15/2018 | Incremental Soil | x | x |

Williston Triplicate Results

| Client Sample ID: | | WIL01IS01 | | WIL01IS02 | | WIL01IS03 | | | | | | | | | |
|-------------------|-------|-----------|-----------|-----------|------|-------------------|-----------------|---------|-------|----------------------|--------------|---------------|-------|------|--|
| Date | | 8/14/18 | | 8/14/18 | | 8/14/18 | | | | | | | | | |
| | Units | RL | 5x LOQ | Sample | Conc | Duplicate Conc | Triplicate Conc | Average | % RSD | Average Deviation | 2x LOQ | Pass/ Fail | | | |
| Metals | | | | | | | | | | | | | | | |
| Antimony | mg/Kg | 0.855 | 4.28 | 0.422 | U | 0.427 | U | 0.383 | U | 0.411 | 4.82% | 0.0185 | 1.71 | Pass | |
| Copper | mg/Kg | 0.427 | 2.14 | 23.8 | | 21.0 | | 24.3 | | 23.0 | 6.30% | 1.36 | 0.854 | Pass | |
| Lead | mg/Kg | 0.427 | 2.14 | 46.5 | | 63.5 | | 69.1 | | 59.7 | 16.1% | 8.80 | 0.854 | Pass | |
| Zinc | mg/Kg | 8.55 | 42.8 | 67.3 | | 61.2 | | 64.5 | | 64.3 | 3.88% | 2.09 | 17.1 | Pass | |

Control limit [sample Average]>5xLOQ use 30%
 [sample Average]<5xLOQ use Average Deviation <2xRL

Williston Triplicate Results

| Client Sample ID: | | WIL01IS01 | | WIL01IS02 | | WIL01IS03 | | | | | | | | |
|----------------------|-------|--|-----------|-----------|------|-------------------|-----------------|----------------|-------|----------------------|--------------|----------------|-------|------|
| Date | | 8/14/18 | | 8/14/18 | | 8/14/18 | | | | | | | | |
| | Units | RL | 5x LOQ | Sample | Conc | Duplicate Conc | Triplicate Conc | Average | % RSD | Average Deviation | 2x LOQ | Pass/ Fail | | |
| Explosives | | | | | | | | All non-detect | | | | | | |
| Metals | | | | | | | | | | | | | | |
| Antimony | mg/Kg | 0.937 | 4.69 | 0.469 | U | 0.453 | U | 0.468 | U | 0.463 | 1.58% | 0.00689 | 1.874 | Pass |
| Copper | mg/Kg | 0.469 | 2.35 | 38.4 | | 33.9 | | 35.6 | | 36.0 | 5.16% | 1.62 | 0.938 | Pass |
| Lead | mg/Kg | 0.469 | 2.35 | 15.9 | | 15.1 | | 15.7 | | 15.6 | 2.18% | 0.311 | 0.938 | Pass |
| Zinc | mg/Kg | 9.37 | 46.9 | 88.5 | | 77.4 | | 81.4 | | 82.4 | 5.57% | 4.04 | 18.74 | Pass |
| Control limit | | [sample Average]>5xLOQ use 30% | | | | | | | | | | | | |
| | | [sample Average]<5xLOQ use Average Deviation <2xRL | | | | | | | | | | | | |

Williston Triplicate Results

| Client Sample ID: | | WIL03IS01 | | WIL03IS02 | | WIL03IS03 | | | | | | | | |
|-------------------|-------|-----------|-----------|----------------|---|-------------------|---|--------------------|---------|-------|----------------------|---------------|---------------|------|
| Date | | 8/14/18 | | 8/14/18 | | 8/14/18 | | | | | | | | |
| | Units | RL | 5x LOQ | Sample Conc | | Duplicate Conc | | Triplicate Conc | Average | % RSD | Average Deviation | 2x LOQ | Pass/ Fail | |
| Metals | | | | | | | | | | | | | | |
| Antimony | mg/Kg | 1.13 | 5.65 | 0.417 | U | 0.423 | U | 0.564 | U | 0.468 | 14.5% | 0.0640 | 2.26 | Pass |
| Copper | mg/Kg | 0.564 | 2.82 | 27.3 | | 23.1 | | 35.7 | | 28.7 | 18.3% | 4.67 | 1.13 | Pass |
| Lead | mg/Kg | 0.564 | 2.82 | 14.7 | | 14.3 | | 22.7 | | 17.2 | 22.5% | 3.64 | 1.13 | Pass |
| Zinc | mg/Kg | 11.3 | 56.5 | 72.9 | | 66.9 | | 81.7 | | 73.8 | 8.23% | 5.24 | 22.6 | Pass |

Control limit [sample Average]>5xLOQ use 30%
 [sample Average]<5xLOQ use Average Deviation <2xRL

Williston Triplicate Results

| Client Sample ID: | | WIL04IS01 | | WIL04IS02 | | WIL04IS03 | | | | | | | | |
|-------------------|-------|-----------|-----------|-----------|------|-------------------|-----------------|---------|-------|----------------------|--------------|----------------|-------|------|
| Date | | 8/15/18 | | 8/15/18 | | 8/15/18 | | | | | | | | |
| | Units | RL | 5x LOQ | Sample | Conc | Duplicate Conc | Triplicate Conc | Average | % RSD | Average Deviation | 2x LOQ | Pass/ Fail | | |
| Metals | | | | | | | | | | | | | | |
| Antimony | mg/Kg | 0.773 | 3.87 | 0.379 | U | 0.387 | U | 0.386 | U | 0.384 | 0.927% | 0.00333 | 1.546 | Pass |
| Copper | mg/Kg | 0.387 | 1.94 | 14.9 | | 14.4 | | 15.4 | | 14.9 | 2.74% | 0.333 | 0.774 | Pass |
| Lead | mg/Kg | 0.387 | 1.94 | 7.11 | | 6.66 | | 7.26 | | 7.01 | 3.64% | 0.233 | 0.774 | Pass |
| Zinc | mg/Kg | 7.73 | 38.7 | 56.4 | | 51.1 | | 56.5 | | 54.7 | 4.61% | 2.38 | 15.46 | Pass |

Control limit [sample Average]>5xLOQ use 30%
 [sample Average]<5xLOQ use Average Deviation <2xRL

**Williston
218081812
Duplicate**

Client Sample ID: WI02SA02A WI02SA02B
Date Sampled: 8/16/18 8/16/18

| | Units | LOQ | 5x LOQ | Sample Conc | | Duplicate Conc | | %RPD | Delta | 2x LOQ | Pass/ Fail |
|-------------------|----------------|-------|-----------|----------------|---|-------------------|---|-------------|--------------|-----------|---------------|
| Explosives | All Non-detect | | | | | | | | | | |
| Metals | | | | | | | | | | | |
| Antimony | mg/Kg | 1.15 | 5.8 | 0.574 | U | 0.535 | U | 7.0% | 0.039 | 2.3 | Pass |
| Copper | mg/Kg | 0.574 | 2.9 | 35.4 | | 38.4 | | 8.1% | 3.0 | 1.1 | Pass |
| Lead | mg/Kg | 0.574 | 2.9 | 18.9 | | 17.3 | | 8.8% | 1.6 | 1.1 | Pass |
| Zinc | mg/Kg | 11.5 | 58 | 101 | | 98.50 | | 2.5% | 2.5 | 23 | Pass |

Control limit [sample]>5xLOQ use 30%
[sample]<5xLOQ use Delta<2xLOQ

1D
ORGANICS ANALYSIS DATA SHEET

| | | | |
|------------------|------------------------------------|--------------------|-----------------------------------|
| Report No: | <u>218081812</u> | Client Sample ID: | <u>WIL02DA02A</u> |
| Collect Date: | <u>08/16/18</u> Time: <u>1020</u> | GCAL Sample ID: | <u>21808181206</u> |
| Matrix: | <u>Solid</u> % Moisture: <u>NA</u> | Instrument ID: | <u>HPLC3</u> |
| Sample Amt: | <u>10.3</u> g | Lab File ID: | <u>2180912VA64</u> |
| Injection Vol.: | <u>1.0</u> (µL) | GC Column: | <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: | <u>40000</u> (µL) | Dilution Factor: | <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: | <u>08/25/18</u> | Analysis Date: | <u>09/13/18</u> Time: <u>0540</u> |
| Prep Batch: | <u>642698</u> | Analytical Batch: | <u>643776</u> |
| Prep Method: | <u>8330B</u> | Analytical Method: | <u>EPA 8330B</u> |

CONCENTRATION UNITS: ug/kg

| CAS | ANALYTE | RESULT | Q | DL | LOD | LOQ |
|------------|-------------------------------------|--------|-----|------|------|-----|
| 99-35-4 | 1,3,5-Trinitrobenzene <i>U.S. 1</i> | 97.1 | UQ | 40.8 | 97.1 | 194 |
| 99-65-0 | 1,3-Dinitrobenzene | 97.1 | UQ | 74.8 | 97.1 | 194 |
| 118-96-7 | 2,4,6-Trinitrotoluene | 97.1 | UQ | 49.5 | 97.1 | 194 |
| 121-14-2 | 2,4-Dinitrotoluene | 97.1 | UQ | 96.1 | 97.1 | 194 |
| 606-20-2 | 2,6-Dinitrotoluene | 97.1 | UQ | 59.2 | 97.1 | 194 |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | 97.1 | UQ | 95.1 | 97.1 | 194 |
| 88-72-2 | 2-Nitrotoluene | 97.1 | UQJ | 62.1 | 97.1 | 194 |
| 618-87-1 | 3,5-Dinitroaniline | 97.1 | UQ | 80.6 | 97.1 | 194 |
| 99-08-1 | 3-Nitrotoluene | 146 | UQ | 121 | 146 | 194 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | 97.1 | UQ | 74.8 | 97.1 | 194 |
| 99-99-0 | 4-Nitrotoluene | 97.1 | UQJ | 74.8 | 97.1 | 194 |
| 2691-41-0 | HMX | 97.1 | UQ | 25.2 | 97.1 | 194 |
| 98-95-3 | Nitrobenzene | 97.1 | UQ | 35.0 | 97.1 | 194 |
| 55-63-0 | Nitroglycerin | 97.1 | UQ | 71.8 | 97.1 | 194 |
| 78-11-5 | Pentaerythritol Tetranitrate | 146 | UQ | 118 | 146 | 194 |
| 121-82-4 | RDX | 97.1 | UQ | 17.5 | 97.1 | 194 |
| 479-45-8 | Tetryl <i>B.I</i> | 97.1 | UQJ | 39.8 | 97.1 | 194 |

FORM I ORG-1

1D
ORGANICS ANALYSIS DATA SHEET

| | | | |
|------------------|------------------------------------|--------------------|-----------------------------------|
| Report No: | <u>218081812</u> | Client Sample ID: | <u>WIL02DA02A (RE)</u> |
| Collect Date: | <u>08/16/18</u> Time: <u>1020</u> | GCAL Sample ID: | <u>21808181213</u> |
| Matrix: | <u>Solid</u> % Moisture: <u>NA</u> | Instrument ID: | <u>HPLC3</u> |
| Sample Amt: | <u>10.1</u> g | Lab File ID: | <u>2180912A72</u> |
| Injection Vol.: | <u>1.0</u> (µL) | GC Column: | <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: | <u>40000</u> (µL) | Dilution Factor: | <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: | <u>09/05/18</u> | Analysis Date: | <u>09/13/18</u> Time: <u>0818</u> |
| Prep Batch: | <u>643342</u> | Analytical Batch: | <u>643776</u> |
| Prep Method: | <u>8330B</u> | Analytical Method: | <u>EPA 8330B</u> |

CONCENTRATION UNITS: ug/kg

| CAS | ANALYTE | RESULT | Q | DL | LOD | LOQ |
|------------|------------------------------|--------|----|------|------|-----|
| 99-35-4 | 1,3,5-Trinitrobenzene | 99.0 | UQ | 41.6 | 99.0 | 198 |
| 99-65-0 | 1,3-Dinitrobenzene | 99.0 | U | 76.2 | 99.0 | 198 |
| 118-96-7 | 2,4,6-Trinitrotoluene | 99.0 | U | 50.5 | 99.0 | 198 |
| 121-14-2 | 2,4-Dinitrotoluene | 99.0 | U | 98.0 | 99.0 | 198 |
| 606-20-2 | 2,6-Dinitrotoluene | 99.0 | UJ | 60.4 | 99.0 | 198 |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | 99.0 | U | 97.0 | 99.0 | 198 |
| 88-72-2 | 2-Nitrotoluene | 99.0 | U | 63.4 | 99.0 | 198 |
| 618-87-1 | 3,5-Dinitroaniline | 99.0 | U | 82.2 | 99.0 | 198 |
| 99-08-1 | 3-Nitrotoluene | 149 | UQ | 124 | 149 | 198 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | 99.0 | U | 76.2 | 99.0 | 198 |
| 99-99-0 | 4-Nitrotoluene | 99.0 | UQ | 76.2 | 99.0 | 198 |
| 2691-41-0 | HMX | 99.0 | UJ | 25.7 | 99.0 | 198 |
| 98-95-3 | Nitrobenzene | 99.0 | U | 35.6 | 99.0 | 198 |
| 55-83-0 | Nitroglycerin | 99.0 | U | 73.3 | 99.0 | 198 |
| 78-11-5 | Pentaerythritol Tetranitrate | 149 | UQ | 121 | 149 | 198 |
| 121-82-4 | RDX | 99.0 | U | 17.8 | 99.0 | 198 |
| 479-45-8 | Tetryl | 99.0 | UQ | 40.6 | 99.0 | 198 |

*NAT
10/24/18
Use initial
results*

1D
ORGANICS ANALYSIS DATA SHEET

| | | | |
|------------------|------------------------------------|--------------------|-----------------------------------|
| Report No: | <u>218081812</u> | Client Sample ID: | <u>WIL02DA01A</u> |
| Collect Date: | <u>08/16/18</u> Time: <u>0930</u> | GCAL Sample ID: | <u>21808181209</u> |
| Matrix: | <u>Solid</u> % Moisture: <u>NA</u> | Instrument ID: | <u>HPLC3</u> |
| Sample Amt: | <u>10.5</u> g | Lab File ID: | <u>2180912VA67</u> |
| Injection Vol.: | <u>1.0</u> (µL) | GC Column: | <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: | <u>40000</u> (µL) | Dilution Factor: | <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: | <u>08/25/18</u> | Analysis Date: | <u>09/13/18</u> Time: <u>0639</u> |
| Prep Batch: | <u>642698</u> | Analytical Batch: | <u>643776</u> |
| Prep Method: | <u>8330B</u> | Analytical Method: | <u>EPA 8330B</u> |

CONCENTRATION UNITS: ug/kg

| CAS | ANALYTE | RESULT | Q | DL | LOD | LOQ |
|------------|-------------------------------------|--------|----|------|------|-----|
| 99-35-4 | 1,3,5-Trinitrobenzene <i>U.S. 1</i> | 95.2 | UQ | 40.0 | 95.2 | 190 |
| 99-65-0 | 1,3-Dinitrobenzene | 95.2 | UQ | 73.3 | 95.2 | 190 |
| 118-96-7 | 2,4,6-Trinitrotoluene | 95.2 | UQ | 48.6 | 95.2 | 190 |
| 121-14-2 | 2,4-Dinitrotoluene | 95.2 | UQ | 94.3 | 95.2 | 190 |
| 606-20-2 | 2,6-Dinitrotoluene | 95.2 | UQ | 58.1 | 95.2 | 190 |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | 95.2 | UQ | 93.3 | 95.2 | 190 |
| 88-72-2 | 2-Nitrotoluene | 95.2 | UQ | 61.0 | 95.2 | 190 |
| 618-87-1 | 3,5-Dinitroaniline | 95.2 | UQ | 79.0 | 95.2 | 190 |
| 99-08-1 | 3-Nitrotoluene | 143 | UQ | 119 | 143 | 190 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | 95.2 | UQ | 73.3 | 95.2 | 190 |
| 99-99-0 | 4-Nitrotoluene | 95.2 | UQ | 73.3 | 95.2 | 190 |
| 2691-41-0 | HMX | 95.2 | UQ | 24.8 | 95.2 | 190 |
| 98-95-3 | Nitrobenzene | 95.2 | UQ | 34.3 | 95.2 | 190 |
| 55-63-0 | Nitroglycerin | 95.2 | UQ | 70.5 | 95.2 | 190 |
| 78-11-5 | Pentaerythritol Tetranitrate | 143 | UQ | 116 | 143 | 190 |
| 121-82-4 | RDX | 95.2 | UQ | 17.1 | 95.2 | 190 |
| 479-45-8 | Tetryl <i>R.I</i> | 95.2 | UQ | 39.0 | 95.2 | 190 |

FORM I ORG-1

1D
ORGANICS ANALYSIS DATA SHEET

| | | | |
|------------------|------------------------------------|--------------------|-----------------------------------|
| Report No: | <u>218081812</u> | Client Sample ID: | <u>WIL02DA01A (RE)</u> |
| Collect Date: | <u>08/16/18</u> Time: <u>0930</u> | GCAL Sample ID: | <u>21808181216</u> |
| Matrix: | <u>Solid</u> % Moisture: <u>NA</u> | Instrument ID: | <u>HPLC3</u> |
| Sample Amt: | <u>10.1</u> g | Lab File ID: | <u>2180912\A75</u> |
| Injection Vol.: | <u>1.0</u> (µL) | GC Column: | <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: | <u>40000</u> (µL) | Dilution Factor: | <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: | <u>09/05/18</u> | Analysis Date: | <u>09/13/18</u> Time: <u>0917</u> |
| Prep Batch: | <u>643342</u> | Analytical Batch: | <u>643776</u> |
| Prep Method: | <u>8330B</u> | Analytical Method: | <u>EPA 8330B</u> |

CONCENTRATION UNITS: ug/kg

| CAS | ANALYTE | RESULT | Q | DL | LOD | LOQ |
|------------|------------------------------|--------|----|------|------|-----|
| 99-35-4 | 1,3,5-Trinitrobenzene | 99.0 | UQ | 41.6 | 99.0 | 198 |
| 99-65-0 | 1,3-Dinitrobenzene | 99.0 | U | 76.2 | 99.0 | 198 |
| 118-96-7 | 2,4,6-Trinitrotoluene | 99.0 | U | 50.5 | 99.0 | 198 |
| 121-14-2 | 2,4-Dinitrotoluene | 99.0 | U | 98.0 | 99.0 | 198 |
| 606-20-2 | 2,6-Dinitrotoluene | 99.0 | U | 60.4 | 99.0 | 198 |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | 99.0 | U | 97.0 | 99.0 | 198 |
| 88-72-2 | 2-Nitrotoluene | 99.0 | U | 63.4 | 99.0 | 198 |
| 618-87-1 | 3,5-Dinitroaniline | 99.0 | U | 82.2 | 99.0 | 198 |
| 99-08-1 | 3-Nitrotoluene | 149 | UQ | 124 | 149 | 198 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | 99.0 | U | 76.2 | 99.0 | 198 |
| 99-99-0 | 4-Nitrotoluene | 99.0 | UQ | 76.2 | 99.0 | 198 |
| 2691-41-0 | HMX | 99.0 | U | 25.7 | 99.0 | 198 |
| 98-95-3 | Nitrobenzene | 99.0 | U | 35.6 | 99.0 | 198 |
| 55-63-0 | Nitroglycerin | 99.0 | U | 73.3 | 99.0 | 198 |
| 78-11-5 | Pentaerythritol Tetranitrate | 149 | UQ | 121 | 149 | 198 |
| 121-82-4 | RDX | 99.0 | U | 17.8 | 99.0 | 198 |
| 479-45-8 | Tetryl | 99.0 | UQ | 40.6 | 99.0 | 198 |

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1D
ORGANICS ANALYSIS DATA SHEET

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|------------------|------------------------------------|--------------------|-----------------------------------|
| Report No: | <u>218081812</u> | Client Sample ID: | <u>WIL02DA01B</u> |
| Collect Date: | <u>08/16/18</u> Time: <u>0935</u> | GCAL Sample ID: | <u>21808181210</u> |
| Matrix: | <u>Solid</u> % Moisture: <u>NA</u> | Instrument ID: | <u>HPLC3</u> |
| Sample Amt: | <u>10</u> g | Lab File ID: | <u>2180912VA68</u> |
| Injection Vol.: | <u>1.0</u> (µL) | GC Column: | <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: | <u>40000</u> (µL) | Dilution Factor: | <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: | <u>08/25/18</u> | Analysis Date: | <u>09/13/18</u> Time: <u>0659</u> |
| Prep Batch: | <u>642698</u> | Analytical Batch: | <u>643776</u> |
| Prep Method: | <u>8330B</u> | Analytical Method: | <u>EPA 8330B</u> |

CONCENTRATION UNITS: ug/kg

| CAS | ANALYTE | RESULT | Q | DL | LOD | LOQ |
|------------|------------------------------|--------|----|------|-----|-----|
| 99-35-4 | 1,3,5-Trinitrobenzene | 100 | UQ | 42.0 | 100 | 200 |
| 99-65-0 | 1,3-Dinitrobenzene | 100 | UQ | 77.0 | 100 | 200 |
| 118-96-7 | 2,4,6-Trinitrotoluene | 100 | UQ | 51.0 | 100 | 200 |
| 121-14-2 | 2,4-Dinitrotoluene | 100 | UQ | 99.0 | 100 | 200 |
| 606-20-2 | 2,6-Dinitrotoluene | 100 | UQ | 61.0 | 100 | 200 |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | 100 | UQ | 98.0 | 100 | 200 |
| 88-72-2 | 2-Nitrotoluene | 100 | UQ | 64.0 | 100 | 200 |
| 618-87-1 | 3,5-Dinitroaniline | 100 | UQ | 83.0 | 100 | 200 |
| 99-08-1 | 3-Nitrotoluene | 150 | UQ | 125 | 150 | 200 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | 100 | UQ | 77.0 | 100 | 200 |
| 99-99-0 | 4-Nitrotoluene | 100 | UQ | 77.0 | 100 | 200 |
| 2691-41-0 | HMX | 100 | UQ | 26.0 | 100 | 200 |
| 98-95-3 | Nitrobenzene | 100 | UQ | 36.0 | 100 | 200 |
| 55-63-0 | Nitroglycerin | 100 | UQ | 74.0 | 100 | 200 |
| 78-11-5 | Pentaerythritol Tetranitrate | 150 | UQ | 122 | 150 | 200 |
| 121-82-4 | RDx | 100 | UQ | 18.0 | 100 | 200 |
| 479-45-8 | Tetryl | 100 | UQ | 41.0 | 100 | 200 |

FORM I ORG-1

1D
ORGANICS ANALYSIS DATA SHEET

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|------------------|------------------------------------|--------------------|-----------------------------------|
| Report No: | <u>218081812</u> | Client Sample ID: | <u>WIL02DA01B (RE)</u> |
| Collect Date: | <u>08/16/18</u> Time: <u>0935</u> | GCAL Sample ID: | <u>21808181217</u> |
| Matrix: | <u>Solid</u> % Moisture: <u>NA</u> | Instrument ID: | <u>HPLC3</u> |
| Sample Amt: | <u>10.1</u> g | Lab File ID: | <u>2180912VA76</u> |
| Injection Vol.: | <u>1.0</u> (µL) | GC Column: | <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: | <u>40000</u> (µL) | Dilution Factor: | <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: | <u>09/05/18</u> | Analysis Date: | <u>09/13/18</u> Time: <u>0937</u> |
| Prep Batch: | <u>643342</u> | Analytical Batch: | <u>643776</u> |
| Prep Method: | <u>8330B</u> | Analytical Method: | <u>EPA 8330B</u> |

CONCENTRATION UNITS: ug/kg

| CAS | ANALYTE | RESULT | Q | DL | LOD | LOQ |
|------------|------------------------------|--------|----|------|------|-----|
| 99-35-4 | 1,3,5-Trinitrobenzene | 99.0 | UQ | 41.6 | 99.0 | 198 |
| 99-65-0 | 1,3-Dinitrobenzene | 99.0 | U | 76.2 | 99.0 | 198 |
| 118-96-7 | 2,4,6-Trinitrotoluene | 99.0 | U | 50.5 | 99.0 | 198 |
| 121-14-2 | 2,4-Dinitrotoluene | 99.0 | U | 98.0 | 99.0 | 198 |
| 606-20-2 | 2,6-Dinitrotoluene | 99.0 | U | 60.4 | 99.0 | 198 |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | 99.0 | U | 97.0 | 99.0 | 198 |
| 88-72-2 | 2-Nitrotoluene | 99.0 | U | 63.4 | 99.0 | 198 |
| 618-87-1 | 3,5-Dinitroaniline | 99.0 | U | 82.2 | 99.0 | 198 |
| 99-08-1 | 3-Nitrotoluene | 149 | UQ | 124 | 149 | 198 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | 99.0 | U | 76.2 | 99.0 | 198 |
| 99-99-0 | 4-Nitrotoluene | 99.0 | UQ | 76.2 | 99.0 | 198 |
| 2691-41-0 | HMX | 99.0 | U | 25.7 | 99.0 | 198 |
| 98-95-3 | Nitrobenzene | 99.0 | U | 35.6 | 99.0 | 198 |
| 55-63-0 | Nitroglycerin | 99.0 | U | 73.3 | 99.0 | 198 |
| 78-11-5 | Pentaerythritol Tetranitrate | 149 | UQ | 121 | 149 | 198 |
| 121-82-4 | RDX | 99.0 | U | 17.8 | 99.0 | 198 |
| 479-45-8 | Tetryl | 99.0 | UQ | 40.6 | 99.0 | 198 |

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1D
ORGANICS ANALYSIS DATA SHEET

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|------------------|------------------------------------|--------------------|-----------------------------------|
| Report No: | <u>218081814</u> | Client Sample ID: | <u>WIL02IS01</u> |
| Collect Date: | <u>08/15/18</u> Time: <u>1200</u> | GCAL Sample ID: | <u>21808181401</u> |
| Matrix: | <u>Solid</u> % Moisture: <u>NA</u> | Instrument ID: | <u>HPLC3</u> |
| Sample Amt: | <u>10.1</u> g | Lab File ID: | <u>2180830VA22</u> |
| Injection Vol.: | <u>1.0</u> (µL) | GC Column: | <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: | <u>40000</u> (µL) | Dilution Factor: | <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: | <u>08/24/18</u> | Analysis Date: | <u>08/30/18</u> Time: <u>1741</u> |
| Prep Batch: | <u>642680</u> | Analytical Batch: | <u>643050</u> |
| Prep Method: | <u>8330B</u> | Analytical Method: | <u>EPA 8330B</u> |

CONCENTRATION UNITS: ug/kg

| CAS | ANALYTE | RESULT | Q | DL | LOD | LOQ |
|------------|------------------------------|--------|----|------|------|-----|
| 99-35-4 | 1,3,5-Trinitrobenzene | 99.0 | U | 41.6 | 99.0 | 198 |
| 99-65-0 | 1,3-Dinitrobenzene | 99.0 | U | 76.2 | 99.0 | 198 |
| 118-96-7 | 2,4,6-Trinitrotoluene | 99.0 | U | 50.5 | 99.0 | 198 |
| 121-14-2 | 2,4-Dinitrotoluene | 99.0 | U | 98.0 | 99.0 | 198 |
| 606-20-2 | 2,6-Dinitrotoluene | 99.0 | U | 60.4 | 99.0 | 198 |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | 99.0 | U | 97.0 | 99.0 | 198 |
| 88-72-2 | 2-Nitrotoluene | 99.0 | U | 63.4 | 99.0 | 198 |
| 618-87-1 | 3,5-Dinitroaniline | 99.0 | U | 82.2 | 99.0 | 198 |
| 99-08-1 | 3-Nitrotoluene | 149 | U | 124 | 149 | 198 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | 99.0 | U | 76.2 | 99.0 | 198 |
| 99-99-0 | 4-Nitrotoluene | 99.0 | UQ | 76.2 | 99.0 | 198 |
| 2691-41-0 | HMX | 99.0 | U | 25.7 | 99.0 | 198 |
| 98-95-3 | Nitrobenzene | 99.0 | U | 35.6 | 99.0 | 198 |
| 55-63-0 | Nitroglycerin | 99.0 | U | 73.3 | 99.0 | 198 |
| 78-11-5 | Pentaerythritol Tetranitrate | 149 | U | 121 | 149 | 198 |
| 121-82-4 | RDX | 99.0 | U | 17.8 | 99.0 | 198 |
| 479-45-8 | Tetryl | 99.0 | U | 40.6 | 99.0 | 198 |

FORM I ORG-1

1D
ORGANICS ANALYSIS DATA SHEET

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|------------------|------------------------------------|--------------------|-----------------------------------|
| Report No: | <u>218081814</u> | Client Sample ID: | <u>WIL02IS02</u> |
| Collect Date: | <u>08/15/18</u> Time: <u>1210</u> | GCAL Sample ID: | <u>21808181404</u> |
| Matrix: | <u>Solid</u> % Moisture: <u>NA</u> | Instrument ID: | <u>HPLC3</u> |
| Sample Amt: | <u>10.5</u> g | Lab File ID: | <u>2180830VA25</u> |
| Injection Vol.: | <u>1.0</u> (µL) | GC Column: | <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: | <u>40000</u> (µL) | Dilution Factor: | <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: | <u>08/24/18</u> | Analysis Date: | <u>08/30/18</u> Time: <u>1840</u> |
| Prep Batch: | <u>642680</u> | Analytical Batch: | <u>643050</u> |
| Prep Method: | <u>8330B</u> | Analytical Method: | <u>EPA 8330B</u> |

CONCENTRATION UNITS: ug/kg

| CAS | ANALYTE | RESULT | Q | DL | LOD | LOQ |
|------------|------------------------------|--------|----|------|------|-----|
| 99-35-4 | 1,3,5-Trinitrobenzene | 95.2 | U | 40.0 | 95.2 | 190 |
| 99-65-0 | 1,3-Dinitrobenzene | 95.2 | U | 73.3 | 95.2 | 190 |
| 118-96-7 | 2,4,6-Trinitrotoluene | 95.2 | U | 48.6 | 95.2 | 190 |
| 121-14-2 | 2,4-Dinitrotoluene | 95.2 | U | 94.3 | 95.2 | 190 |
| 606-20-2 | 2,6-Dinitrotoluene | 95.2 | U | 58.1 | 95.2 | 190 |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | 95.2 | U | 93.3 | 95.2 | 190 |
| 88-72-2 | 2-Nitrotoluene | 95.2 | U | 61.0 | 95.2 | 190 |
| 618-87-1 | 3,5-Dinitroaniline | 95.2 | U | 79.0 | 95.2 | 190 |
| 99-08-1 | 3-Nitrotoluene | 143 | U | 119 | 143 | 190 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | 95.2 | U | 73.3 | 95.2 | 190 |
| 99-99-0 | 4-Nitrotoluene | 95.2 | UQ | 73.3 | 95.2 | 190 |
| 2691-41-0 | HMX | 95.2 | U | 24.8 | 95.2 | 190 |
| 98-95-3 | Nitrobenzene | 95.2 | U | 34.3 | 95.2 | 190 |
| 55-63-0 | Nitroglycerin | 95.2 | U | 70.5 | 95.2 | 190 |
| 78-11-5 | Pentaerythritol Tetranitrate | 143 | U | 116 | 143 | 190 |
| 121-82-4 | RDX | 95.2 | U | 17.1 | 95.2 | 190 |
| 479-45-8 | Tetryl | 95.2 | U | 39.0 | 95.2 | 190 |

FORM I ORG-1

1D
ORGANICS ANALYSIS DATA SHEET

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| Report No: <u>218081814</u> | Client Sample ID: <u>WIL02IS03</u> |
| Collect Date: <u>08/15/18</u> Time: <u>1220</u> | GCAL Sample ID: <u>21808181407</u> |
| Matrix: <u>Solid</u> % Moisture: <u>NA</u> | Instrument ID: <u>HPLC3</u> |
| Sample Amt: <u>10</u> g | Lab File ID: <u>2180830A26</u> |
| Injection Vol.: <u>1.0</u> (µL) | GC Column: <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: <u>40000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: <u>08/24/18</u> | Analysis Date: <u>08/30/18</u> Time: <u>1900</u> |
| Prep Batch: <u>642680</u> | Analytical Batch: <u>643050</u> |
| Prep Method: <u>8330B</u> | Analytical Method: <u>EPA 8330B</u> |

CONCENTRATION UNITS: ug/kg

| CAS | ANALYTE | RESULT | Q | DL | LOD | LOQ |
|------------|------------------------------|--------|----|------|-----|-----|
| 99-35-4 | 1,3,5-Trinitrobenzene | 100 | U | 42.0 | 100 | 200 |
| 99-65-0 | 1,3-Dinitrobenzene | 100 | U | 77.0 | 100 | 200 |
| 118-96-7 | 2,4,6-Trinitrotoluene | 100 | U | 51.0 | 100 | 200 |
| 121-14-2 | 2,4-Dinitrotoluene | 100 | U | 99.0 | 100 | 200 |
| 606-20-2 | 2,6-Dinitrotoluene | 100 | U | 61.0 | 100 | 200 |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | 100 | U | 98.0 | 100 | 200 |
| 88-72-2 | 2-Nitrotoluene | 100 | U | 64.0 | 100 | 200 |
| 618-87-1 | 3,5-Dinitroaniline | 100 | U | 83.0 | 100 | 200 |
| 99-08-1 | 3-Nitrotoluene | 150 | U | 125 | 150 | 200 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | 100 | U | 77.0 | 100 | 200 |
| 99-99-0 | 4-Nitrotoluene | 100 | UQ | 77.0 | 100 | 200 |
| 2691-41-0 | HMX | 100 | U | 26.0 | 100 | 200 |
| 98-95-3 | Nitrobenzene | 100 | U | 36.0 | 100 | 200 |
| 55-63-0 | Nitroglycerin | 100 | U | 74.0 | 100 | 200 |
| 78-11-5 | Pentaerythritol Tetranitrate | 150 | U | 122 | 150 | 200 |
| 121-82-4 | RDX | 100 | U | 18.0 | 100 | 200 |
| 479-45-8 | Tetryl | 100 | U | 41.0 | 100 | 200 |

FORM I ORG-1

I
INORGANIC ANALYSIS DATA SHEET

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|---------------|-------------------------------------|--------------------|------------------------------------|
| Report No: | <u>218081812</u> | Client Sample ID: | <u>WIL02DA02A</u> |
| Collect Date: | <u>08/16/18</u> Time: <u>1020</u> | GCAL Sample ID: | <u>21808181206</u> |
| Matrix: | <u>Solid</u> % Solids: <u>69.73</u> | Instrument ID: | <u>ICPMS2</u> |
| Sample Amt: | <u>1.25</u> g | Lab File ID: | <u>2180820B_MS2.b\121277SMPL.d</u> |
| Prep Vol.: | <u>50</u> (mL) | Dilution Factor: | <u>10</u> Analyst: <u>LWZ</u> |
| Prep Date: | <u>08/20/18</u> | Analysis Date: | <u>08/20/18</u> Time: <u>1802</u> |
| Prep Batch: | <u>642222</u> | Analytical Batch: | <u>642309</u> |
| Prep Method: | <u>3050B</u> | Analytical Method: | <u>EPA 6020B</u> |

| ANALYTE | RESULT | UNITS | Q | DL | LOD | LOQ |
|----------|------------------|-------|---|------|------|-------|
| Antimony | 574 <i>US,m</i> | ug/kg | U | 287 | 574 | 1150 |
| Copper | 35400 | ug/kg | | 143 | 287 | 574 |
| Lead | 18900 | ug/kg | | 143 | 287 | 574 |
| Zinc | 101000 <i>JS</i> | ug/kg | | 2870 | 5740 | 11500 |

I
INORGANIC ANALYSIS DATA SHEET

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|---------------|-------------------------------------|--------------------|------------------------------------|
| Report No: | <u>218081812</u> | Client Sample ID: | <u>WIL02DA01A</u> |
| Collect Date: | <u>08/16/18</u> Time: <u>0930</u> | GCAL Sample ID: | <u>21808181209</u> |
| Matrix: | <u>Solid</u> % Solids: <u>70.44</u> | Instrument ID: | <u>ICPMS2</u> |
| Sample Amt: | <u>1.45</u> g | Lab File ID: | <u>2180820B_MS2.b\121282SMPL.d</u> |
| Prep Vol.: | <u>50</u> (mL) | Dilution Factor: | <u>10</u> Analyst: <u>LWZ</u> |
| Prep Date: | <u>08/20/18</u> | Analysis Date: | <u>08/20/18</u> Time: <u>1820</u> |
| Prep Batch: | <u>642222</u> | Analytical Batch: | <u>642309</u> |
| Prep Method: | <u>3050B</u> | Analytical Method: | <u>EPA 6020B</u> |

| ANALYTE | RESULT | UNITS | Q | DL | LOD | LOQ |
|----------|-------------------|-------|---|------|------|------|
| Antimony | 490 <i>U.S.M</i> | ug/kg | U | 245 | 490 | 979 |
| Copper | 39300 | ug/kg | | 122 | 245 | 490 |
| Lead | 17600 | ug/kg | | 122 | 245 | 490 |
| Zinc | 98900 <i>S, 2</i> | ug/kg | | 2450 | 4900 | 9790 |

FORM I - IN

I
INORGANIC ANALYSIS DATA SHEET

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|---------------|-------------------------------------|--------------------|------------------------------------|
| Report No: | <u>218081812</u> | Client Sample ID: | <u>WIL02DA01B</u> |
| Collect Date: | <u>08/16/18</u> Time: <u>0935</u> | GCAL Sample ID: | <u>21808181210</u> |
| Matrix: | <u>Solid</u> % Solids: <u>70.26</u> | Instrument ID: | <u>ICPMS2</u> |
| Sample Amt: | <u>1.33</u> g | Lab File ID: | <u>2180820B_MS2.b\121283SMPL.d</u> |
| Prep Vol.: | <u>50</u> (mL) | Dilution Factor: | <u>10</u> Analyst: <u>LWZ</u> |
| Prep Date: | <u>08/20/18</u> | Analysis Date: | <u>08/20/18</u> Time: <u>1824</u> |
| Prep Batch: | <u>642222</u> | Analytical Batch: | <u>642309</u> |
| Prep Method: | <u>3050B</u> | Analytical Method: | <u>EPA 6020B</u> |

| ANALYTE | RESULT | UNITS | Q | DL | LOD | LOQ |
|----------|------------------|-------|---|------|------|-------|
| Antimony | 535 <i>USm</i> | ug/kg | U | 268 | 535 | 1070 |
| Copper | 38400 | ug/kg | | 134 | 268 | 535 |
| Lead | 17300 | ug/kg | | 134 | 268 | 535 |
| Zinc | 98500 <i>S.S</i> | ug/kg | | 2680 | 5350 | 10700 |

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INORGANIC ANALYSIS DATA SHEET

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|---|--|
| Report No: <u>218081813</u> | Client Sample ID: <u>WIL03IS01</u> |
| Collect Date: <u>08/14/18</u> Time: <u>1600</u> | GCAL Sample ID: <u>21808181301</u> |
| Matrix: <u>Solid</u> % Solids: <u>88.78</u> | Instrument ID: <u>ICPMS2</u> |
| Sample Amt: <u>1.35</u> g | Lab File ID: <u>2180828B_MS2.b\1841SMPL_2180828A_MS2.D</u> |
| Prep Vol.: <u>50</u> (mL) | Dilution Factor: <u>10</u> Analyst: <u>LWZ</u> |
| Prep Date: <u>08/25/18</u> | Analysis Date: <u>08/28/18</u> Time: <u>1305</u> |
| Prep Batch: <u>642531</u> | Analytical Batch: <u>642829</u> |
| Prep Method: <u>EPA 3050B \ ISM</u> | Analytical Method: <u>EPA 6020B</u> |

| ANALYTE | RESULT | UNITS | Q | DL | LOD | LOQ |
|----------|-----------------|-------|---|------|------|------|
| Antimony | 417 <i>US.m</i> | ug/kg | U | 209 | 417 | 834 |
| Copper | 27300 | ug/kg | | 104 | 209 | 417 |
| Lead | 14700 | ug/kg | | 104 | 209 | 417 |
| Zinc | 72900 | ug/kg | | 2090 | 4170 | 8340 |

FORM I - IN

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INORGANIC ANALYSIS DATA SHEET

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|---|--|--|
| Report No: <u>218081813</u> | Client Sample ID: <u>WIL01IS02</u> | |
| Collect Date: <u>08/14/18</u> Time: <u>1350</u> | GCAL Sample ID: <u>21808181302</u> | |
| Matrix: <u>Solid</u> % Solids: <u>93.58</u> | Instrument ID: <u>ICPMS1</u> | |
| Sample Amt: <u>1.25</u> g | Lab File ID: <u>2180823A_MS1.b\038SMPL.d</u> | |
| Prep Vol.: <u>50</u> (mL) | Dilution Factor: <u>10</u> Analyst: <u>LWZ</u> | |
| Prep Date: <u>08/22/18</u> | Analysis Date: <u>08/23/18</u> Time: <u>1209</u> | |
| Prep Batch: <u>642442</u> | Analytical Batch: <u>642536</u> | |
| Prep Method: <u>EPA 3050B \ ISM</u> | Analytical Method: <u>EPA 6020B</u> | |

| ANALYTE | RESULT | UNITS | Q | DL | LOD | LOQ |
|----------|------------------|-------|---|------|------|------|
| Antimony | 427 <i>VSM</i> | ug/kg | U | 214 | 427 | 855 |
| Copper | 21000 <i>J+M</i> | ug/kg | | 107 | 214 | 427 |
| Lead | 63500 | ug/kg | | 107 | 214 | 427 |
| Zinc | 61200 | ug/kg | | 2140 | 4270 | 8550 |

FORM I - IN

I
INORGANIC ANALYSIS DATA SHEET

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|---------------|-----------------------------------|--------------------|-----------------------------------|
| Report No: | <u>218081813</u> | Client Sample ID: | <u>WIL03IS00</u> |
| Collect Date: | <u>08/16/18</u> Time: <u>0820</u> | GCAL Sample ID: | <u>21808181303</u> |
| Matrix: | <u>Water</u> % Solids: <u>NA</u> | Instrument ID: | <u>ICPMS1</u> |
| Sample Amt: | <u>50</u> mL | Lab File ID: | <u>2180821A_MS1.b\023SMPL.d</u> |
| Prep Vol.: | <u>50</u> (mL) | Dilution Factor: | <u>1</u> Analyst: <u>LWZ</u> |
| Prep Date: | <u>08/20/18</u> | Analysis Date: | <u>08/21/18</u> Time: <u>1151</u> |
| Prep Batch: | <u>642277</u> | Analytical Batch: | <u>642381</u> |
| Prep Method: | <u>3010A</u> | Analytical Method: | <u>EPA 6020B</u> |

| ANALYTE | RESULT | UNITS | Q | DL | LOD | LOQ |
|----------|--------|-------|---|------|------|------|
| Antimony | 0.74 | ug/L | J | 0.50 | 1.00 | 2.00 |
| Copper | 0.26 | ug/L | J | 0.25 | 0.50 | 1.00 |
| Lead | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Zinc | 10.0 | ug/L | U | 5.00 | 10.0 | 20.0 |

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|---|--|
| Report No: <u>218081813</u> | Client Sample ID: <u>WIL04IS03</u> |
| Collect Date: <u>08/15/18</u> Time: <u>1610</u> | GCAL Sample ID: <u>21808181304</u> |
| Matrix: <u>Solid</u> % Solids: <u>97.47</u> | Instrument ID: <u>ICPMS1</u> |
| Sample Amt: <u>1.33</u> g | Lab File ID: <u>2180823A_MS1.b\043SMPL.d</u> |
| Prep Vol.: <u>50</u> (mL) | Dilution Factor: <u>10</u> Analyst: <u>LWZ</u> |
| Prep Date: <u>08/22/18</u> | Analysis Date: <u>08/23/18</u> Time: <u>1231</u> |
| Prep Batch: <u>642442</u> | Analytical Batch: <u>642536</u> |
| Prep Method: <u>EPA 3050B \ ISM</u> | Analytical Method: <u>EPA 6020B</u> |

| ANALYTE | RESULT | UNITS | Q | DL | LOD | LOQ |
|----------|-------------------|-------|---|------|------|------|
| Antimony | 386 <i>U.S.M</i> | ug/kg | U | 193 | 386 | 771 |
| Copper | 15400 <i>Sr.M</i> | ug/kg | | 96.4 | 193 | 386 |
| Lead | 7260 | ug/kg | | 96.4 | 193 | 386 |
| Zinc | 56500 | ug/kg | | 1930 | 3860 | 7710 |

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|---|--|--|
| Report No: <u>218081813</u> | Client Sample ID: <u>WIL04IS01</u> | |
| Collect Date: <u>08/15/18</u> Time: <u>1600</u> | GCAL Sample ID: <u>21808181305</u> | |
| Matrix: <u>Solid</u> % Solids: <u>97.68</u> | Instrument ID: <u>ICPMS1</u> | |
| Sample Amt: <u>1.35</u> g | Lab File ID: <u>2180823A_MS1.b\044SMPL.d</u> | |
| Prep Vol.: <u>50</u> (mL) | Dilution Factor: <u>10</u> Analyst: <u>LWZ</u> | |
| Prep Date: <u>08/22/18</u> | Analysis Date: <u>08/23/18</u> Time: <u>1235</u> | |
| Prep Batch: <u>642442</u> | Analytical Batch: <u>642536</u> | |
| Prep Method: <u>EPA 3050B \ ISM</u> | Analytical Method: <u>EPA 6020B</u> | |

| ANALYTE | RESULT | UNITS | Q | DL | LOD | LOQ |
|----------|-----------------|-------|---|------|------|------|
| Antimony | 379 <i>VJM</i> | ug/kg | U | 190 | 379 | 758 |
| Copper | 14900 <i>Jm</i> | ug/kg | | 94.8 | 190 | 379 |
| Lead | 7110 | ug/kg | | 94.8 | 190 | 379 |
| Zinc | 56400 | ug/kg | | 1900 | 3790 | 7580 |

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|---------------|-------------------------------------|--------------------|-----------------------------------|
| Report No: | <u>218081813</u> | Client Sample ID: | <u>WIL04IS02</u> |
| Collect Date: | <u>08/15/18</u> Time: <u>1605</u> | GCAL Sample ID: | <u>21808181306</u> |
| Matrix: | <u>Solid</u> % Solids: <u>97.99</u> | Instrument ID: | <u>ICPMS1</u> |
| Sample Amt: | <u>1.32</u> g | Lab File ID: | <u>2180823A_MS1.b\045SMPL.d</u> |
| Prep Vol.: | <u>50</u> (mL) | Dilution Factor: | <u>10</u> Analyst: <u>LWZ</u> |
| Prep Date: | <u>08/22/18</u> | Analysis Date: | <u>08/23/18</u> Time: <u>1240</u> |
| Prep Batch: | <u>642442</u> | Analytical Batch: | <u>642536</u> |
| Prep Method: | <u>EPA 3050B \ ISM</u> | Analytical Method: | <u>EPA 6020B</u> |

| ANALYTE | RESULT | UNITS | Q | DL | LOD | LOQ |
|----------|------------------|-------|---|------|------|------|
| Antimony | 387 <i>U.S.M</i> | ug/kg | U | 193 | 387 | 773 |
| Copper | 14400 <i>S.M</i> | ug/kg | | 96.6 | 193 | 387 |
| Lead | 6660 | ug/kg | | 96.6 | 193 | 387 |
| Zinc | 51100 | ug/kg | | 1930 | 3870 | 7730 |

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| | | |
|---|--|--|
| Report No: <u>218081813</u> | Client Sample ID: <u>WIL03IS03</u> | |
| Collect Date: <u>08/14/18</u> Time: <u>1610</u> | GCAL Sample ID: <u>21808181307</u> | |
| Matrix: <u>Solid</u> % Solids: <u>70.33</u> | Instrument ID: <u>ICPMS2</u> | |
| Sample Amt: <u>1.26</u> g | Lab File ID: <u>2180828C_MS2.b\001SMPL.d</u> | |
| Prep Vol.: <u>50</u> (mL) | Dilution Factor: <u>10</u> Analyst: <u>LWZ</u> | |
| Prep Date: <u>08/25/18</u> | Analysis Date: <u>08/28/18</u> Time: <u>1355</u> | |
| Prep Batch: <u>642531</u> | Analytical Batch: <u>642829</u> | |
| Prep Method: <u>EPA 3050B \ ISM</u> | Analytical Method: <u>EPA 6020B</u> | |

| ANALYTE | RESULT | UNITS | Q | DL | LOD | LOQ |
|----------|------------------|-------|---|------|------|-------|
| Antimony | 564 <i>U.S.m</i> | ug/kg | U | 282 | 564 | 1130 |
| Copper | 35700 | ug/kg | | 141 | 282 | 564 |
| Lead | 22700 | ug/kg | | 141 | 282 | 564 |
| Zinc | 81700 | ug/kg | | 2820 | 5640 | 11300 |

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| | | |
|---|--|--|
| Report No: <u>218081813</u> | Client Sample ID: <u>WIL03IS02</u> | |
| Collect Date: <u>08/14/18</u> Time: <u>1605</u> | GCAL Sample ID: <u>21808181308</u> | |
| Matrix: <u>Solid</u> % Solids: <u>92.27</u> | Instrument ID: <u>ICPMS2</u> | |
| Sample Amt: <u>1.28</u> g | Lab File ID: <u>2180828B_MS2.b\1842SMPL_2180828A_MS2.D</u> | |
| Prep Vol.: <u>50</u> (mL) | Dilution Factor: <u>10</u> Analyst: <u>LWZ</u> | |
| Prep Date: <u>08/25/18</u> | Analysis Date: <u>08/28/18</u> Time: <u>1308</u> | |
| Prep Batch: <u>642531</u> | Analytical Batch: <u>642829</u> | |
| Prep Method: <u>EPA 3050B \ ISM</u> | Analytical Method: <u>EPA 6020B</u> | |

| ANALYTE | RESULT | UNITS | Q | DL | LOD | LOQ |
|----------|----------------|-------|---|------|------|------|
| Antimony | 423 <i>USM</i> | ug/kg | U | 212 | 423 | 847 |
| Copper | 23100 | ug/kg | | 106 | 212 | 423 |
| Lead | 14300 | ug/kg | | 106 | 212 | 423 |
| Zinc | 66900 | ug/kg | | 2120 | 4230 | 8470 |

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| | | | |
|---------------|-------------------------------------|--------------------|---|
| Report No: | <u>218081814</u> | Client Sample ID: | <u>WIL02IS01</u> |
| Collect Date: | <u>08/15/18</u> Time: <u>1200</u> | GCAL Sample ID: | <u>21808181401</u> |
| Matrix: | <u>Solid</u> % Solids: <u>85.34</u> | Instrument ID: | <u>ICPMS2</u> |
| Sample Amt: | <u>1.25</u> g | Lab File ID: | <u>2180828B_MS2.b\1843SMPL_2180828A_MS2.D</u> |
| Prep Vol.: | <u>50</u> (mL) | Dilution Factor: | <u>10</u> Analyst: <u>LWZ</u> |
| Prep Date: | <u>08/25/18</u> | Analysis Date: | <u>08/28/18</u> Time: <u>1312</u> |
| Prep Batch: | <u>642531</u> | Analytical Batch: | <u>642829</u> |
| Prep Method: | <u>EPA 3050B \ ISM</u> | Analytical Method: | <u>EPA 6020B</u> |

| ANALYTE | RESULT | UNITS | Q | DL | LOD | LOQ |
|----------|----------------|-------|---|------|------|------|
| Antimony | 469 <i>USm</i> | ug/kg | U | 234 | 469 | 937 |
| Copper | 38400 | ug/kg | | 117 | 234 | 469 |
| Lead | 15900 | ug/kg | | 117 | 234 | 469 |
| Zinc | 88500 | ug/kg | | 2340 | 4690 | 9370 |

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| | | | |
|---------------|-------------------------------------|--------------------|---|
| Report No: | <u>218081814</u> | Client Sample ID: | <u>WIL02IS02</u> |
| Collect Date: | <u>08/15/18</u> Time: <u>1210</u> | GCAL Sample ID: | <u>21808181404</u> |
| Matrix: | <u>Solid</u> % Solids: <u>84.25</u> | Instrument ID: | <u>ICPMS2</u> |
| Sample Amt: | <u>1.31</u> g | Lab File ID: | <u>2180828B_MS2.b\1848SMPL_2180828A_MS2.D</u> |
| Prep Vol.: | <u>50</u> (mL) | Dilution Factor: | <u>10</u> Analyst: <u>LWZ</u> |
| Prep Date: | <u>08/25/18</u> | Analysis Date: | <u>08/28/18</u> Time: <u>1329</u> |
| Prep Batch: | <u>642531</u> | Analytical Batch: | <u>642829</u> |
| Prep Method: | <u>EPA 3050B \ ISM</u> | Analytical Method: | <u>EPA 6020B</u> |

| ANALYTE | RESULT | UNITS | Q | DL | LOD | LOQ |
|----------|-----------------|-------|---|------|------|------|
| Antimony | 453 <i>US,m</i> | ug/kg | U | 227 | 453 | 906 |
| Copper | 33900 | ug/kg | | 113 | 227 | 453 |
| Lead | 15100 | ug/kg | | 113 | 227 | 453 |
| Zinc | 77400 | ug/kg | | 2270 | 4530 | 9060 |

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INORGANIC ANALYSIS DATA SHEET

| | | | |
|---------------|-------------------------------------|--------------------|---|
| Report No: | <u>218081814</u> | Client Sample ID: | <u>WIL01IS03</u> |
| Collect Date: | <u>08/14/18</u> Time: <u>1400</u> | GCAL Sample ID: | <u>21808181405</u> |
| Matrix: | <u>Solid</u> % Solids: <u>93.16</u> | Instrument ID: | <u>ICPMS2</u> |
| Sample Amt: | <u>1.4</u> g | Lab File ID: | <u>2180828B_MS2.b\1849SMPL_2180828A_MS2.D</u> |
| Prep Vol.: | <u>50</u> (mL) | Dilution Factor: | <u>10</u> Analyst: <u>LWZ</u> |
| Prep Date: | <u>08/25/18</u> | Analysis Date: | <u>08/28/18</u> Time: <u>1333</u> |
| Prep Batch: | <u>642531</u> | Analytical Batch: | <u>642829</u> |
| Prep Method: | <u>EPA 3050B \ ISM</u> | Analytical Method: | <u>EPA 6020B</u> |

| ANALYTE | RESULT | UNITS | Q | DL | LOD | LOQ |
|----------|-----------------|-------|---|------|------|------|
| Antimony | 383 <i>US,m</i> | ug/kg | U | 192 | 383 | 767 |
| Copper | 24300 | ug/kg | | 95.8 | 192 | 383 |
| Lead | 69100 | ug/kg | | 95.8 | 192 | 383 |
| Zinc | 64500 | ug/kg | | 1920 | 3830 | 7670 |

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INORGANIC ANALYSIS DATA SHEET

| | | | |
|---------------|-------------------------------------|--------------------|---|
| Report No: | <u>218081814</u> | Client Sample ID: | <u>WIL01IS01</u> |
| Collect Date: | <u>08/14/18</u> Time: <u>1340</u> | GCAL Sample ID: | <u>21808181406</u> |
| Matrix: | <u>Solid</u> % Solids: <u>94.08</u> | Instrument ID: | <u>ICPMS2</u> |
| Sample Amt: | <u>1.26</u> g | Lab File ID: | <u>2180828B_MS2.b\1852SMPL_2180828A_MS2.D</u> |
| Prep Vol.: | <u>50</u> (mL) | Dilution Factor: | <u>10</u> Analyst: <u>LWZ</u> |
| Prep Date: | <u>08/25/18</u> | Analysis Date: | <u>08/28/18</u> Time: <u>1344</u> |
| Prep Batch: | <u>642531</u> | Analytical Batch: | <u>642829</u> |
| Prep Method: | <u>EPA 3050B \ ISM</u> | Analytical Method: | <u>EPA 6020B</u> |

| ANALYTE | RESULT | UNITS | Q | DL | LOD | LOQ |
|----------|-----------------|-------|---|------|------|------|
| Antimony | 422 <i>US,m</i> | ug/kg | U | 211 | 422 | 844 |
| Copper | 23800 | ug/kg | | 105 | 211 | 422 |
| Lead | 46500 | ug/kg | | 105 | 211 | 422 |
| Zinc | 67300 | ug/kg | | 2110 | 4220 | 8440 |

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| | | | |
|---------------|-------------------------------------|--------------------|---|
| Report No: | <u>218081814</u> | Client Sample ID: | <u>WIL02IS03</u> |
| Collect Date: | <u>08/15/18</u> Time: <u>1220</u> | GCAL Sample ID: | <u>21808181407</u> |
| Matrix: | <u>Solid</u> % Solids: <u>79.69</u> | Instrument ID: | <u>ICPMS2</u> |
| Sample Amt: | <u>1.34</u> g | Lab File ID: | <u>2180828B_MS2.b\1853SMPL_2180828A_MS2.D</u> |
| Prep Vol.: | <u>50</u> (mL) | Dilution Factor: | <u>10</u> Analyst: <u>LWZ</u> |
| Prep Date: | <u>08/25/18</u> | Analysis Date: | <u>08/28/18</u> Time: <u>1347</u> |
| Prep Batch: | <u>642531</u> | Analytical Batch: | <u>642829</u> |
| Prep Method: | <u>EPA 3050B \ ISM</u> | Analytical Method: | <u>EPA 6020B</u> |

| ANALYTE | RESULT | UNITS | Q | DL | LOD | LOQ |
|----------|------------------|-------|---|------|------|------|
| Antimony | 468 <i>U.S.m</i> | ug/kg | U | 234 | 468 | 936 |
| Copper | 35600 | ug/kg | | 117 | 234 | 468 |
| Lead | 15700 | ug/kg | | 117 | 234 | 468 |
| Zinc | 81400 | ug/kg | | 2340 | 4680 | 9360 |

FORM I - IN

DATA VALIDATION WORKSHEET

Reviewer: Naoum Tavantzis
Date: 10/24/2018
DV Level: II III IV

Explosives

Project Name: Williston LTA
Project Number: 60520956
Laboratory: GCAL
SDG No.: 218081812+13+14
Test Name: Explosives
Method No.: 8330B

Review Document:

- USEPA CLP National Functional Guidelines for Superfund Inorganic Methods Data Review
- SW-846/DoD QSM v5.1
- Project QAPP/SAP

1.0 Laboratory Deliverables

| | | Yes | No | NA |
|-----|--|-----|----|----|
| 1.1 | Do Chain-of-Custody forms list all samples that were analyzed? | X | | |
| 1.2 | Are all Chain-of-Custody forms signed, indicating sample chain-of-custody was maintained? | X | | |
| 1.3 | Do sample preservation, collection and storage condition meet method requirement? | X | | |
| | If samples were received with the cooler temperature exceeding 10 ⁰ C, then flag J(+)/UJ(-). | | | |
| 1.4 | Were the sediment samples dried and sieved appropriately? | X | | |
| 1.5 | Do the traffic Reports, chain-of-custody, and lab narrative indicate any problems with sample receipt, condition of samples, analytical problems or special circumstances affecting the quality of the data? | | X | |

Notes:

2.0 Holding Times

| | | Yes | No | NA |
|-----|--|-----|----|----|
| 2.1 | Have any technical holding times, determined from date of sampling to date of analysis, been exceeded? If yes, J(+)/UJ(-). Extraction: (aqueous) 7 days and (soil) 14 days; and Analysis: 40 days. | X | | |
| 2.2 | Have any technical holding time grossly (twice the holding time) been exceeded? If yes, J(+)/UJ(-) OR R(-) . | | X | |

Notes: The field sample results associated with low LCS percent recoveries were re-extracted after the holding time of 14 days had expired.

3.0 Blanks (Laboratory and Field)

| | | Yes | No | NA |
|-----|---|-----|----|----|
| 3.1 | Were method blanks (MB) prepared at the appropriate frequency (one per 20 samples, per batch per matrix?) | X | | |
| 3.2 | Do any method blanks have positive results? Action: If yes: | | | |
| | Blank Result: ≤LOD >LOD | | X | |
| | Sample Result: ≥[LOD] <[LOD] <LOD ≥LOD, <[Blank] ≥[LOD], ≥[Blank] | | | |
| 3.3 | Do any field equipment blanks/trip blanks have positive results? If yes, use same rules above. | | X | |

Notes:

4.0 Initial and Continuing Calibration

| | | Yes | No | NA |
|--|---|-----|----|----|
| 4.1 | Are the Quartly LOD establishment forms provided? Is the signal/noise ratio of the LODV greater than 3:1? | X | | |
| 4.2 | If not, is the determination repeated at a higher concentration? | X | | |
| 4.3 | Are at least five standards included in the calibration curve? If no, flag "R". | X | | |
| 4.4 | Was a second source calibration verification analyzed for each calibration curve? If no, flag "R". | X | | |
| 4.5 | Were continuing calibration standards analyzed every 12 hours or ten samples and at the end of the sequence? If no, flag "R". | X | | |
| 4.6 | Are all calibration standard %RSD (<15%ICAL, r>0.995), second source (±20%) or %D(≤20%) within the control limits? | X | | |
| For initial calibration: %RSD > 15% J(+) If RRF<0.05 J(+)/R(-) | | | | |
| For second source: %D>20%, J(+)/R(-). | | | | |
| For continuing calibration: Positive Bias - %D >+ 20%, J(+), only. Negative Bias - %D>-20%, but<50% J(+)/UJ(-) | | | | |

Notes:

5.0 Laboratory Control Sample (LCS)

| | | Yes | No | NA |
|-----|--|-----|----|----|
| 5.1 | Is the LCS/LCSD recovery form present? | X | | |
| 5.2 | Were LCS/LCSDs analyzed at the required frequency (one per 20 samples per batch) for each matrix? | | X | |
| 5.3 | Are there any %Rs for LCS/LCSD recoveries outside the QC limits of 80-120%? Action: If Yes, for %R > UCL, J(+) only; for %R < LCL, J(+)/R(-); | X | | |
| 5.4 | Are there any RPD >20%? Action: If Yes, J(+) only. | X | | |

Notes: Several LCS recoveries less than the lower QC limits in batch 642698, and recoveries higher than the QC limits in batch 643342. Both batches displayed zero percent recoveries for tetryl. In addition, the LCS/LCSD displayed several relative percent difference (RPD) anomalies greater than the laboratory QC limit of 20%.

6.0 Surrogate Recovery

| | | Yes | No | NA |
|-----|--|-----|----|----|
| 6.1 | Are surrogate recoveries within acceptance criteria for all samples and method blanks? | X | | |
| 6.2 | If No in Section 6.1, are these sample(s) or method blank(s) reanalyzed? Action: If any surrogate in the fraction is out of specification, there should be a reanalysis to confirm that the non-compliance is because of sample matrix effects rather than laboratory deficiencies. | | | X |
| 6.3 | If No in Section 6.2, is any sample dilution factor greater than 10? (recoveries may be diluted out.) Action: for a %R > UCL, J(+) only; for a %R < LCL, but > 10%, J(+)/UJ(-); for any %R < 10%, J(+)/R(-). | | | X |

Notes:

7.0 Matrix Spike/Matrix Spike Duplicate (MS/MSD)

| | | Yes | No | NA |
|-----|--|-----|----|----|
| 7.1 | Is the matrix spike/matrix spike duplicate recovery form present? | X | | |
| 7.2 | Were matrix spikes analyzed at required frequency (one per 20 samples per batch) for each matrix? | X | | |
| 7.3 | Are there any %R for matrix spike and matrix spike duplicate recoveries outside the QC limits? | X | | |
| 7.4 | Are there any RPDs for matrix spike and matrix spike duplicate recoveries outside the QC limits? | X | | |
| | Action: %R or % RPD > UCL J(+); %R < LCL: J(+)/UJ(-) [or possible R(-) for <20%] to parent sample results for the specific outlier analyte(s) or class of analytes using professional judgement. | | | |

Notes: Matrix spike performed on WIL02DA02A in QC batch 643342 displayed a recovery less than the lower QC limit for HMX; the MS pair performed on WIL02DA02A displayed several RPD anomalies.

8.0 Field Duplicate /Triplicate

| | | Yes | No | NA |
|-----|---|-----|----|----|
| 8.1 | Evaluate field duplicate results For sample results > 5 x RL, a control limit of ≤30% RPD/RSD will be used. For sample results < 5 x RL, a control limit of 2 x RL will be used. | X | | |

Notes:

9.0 Compound Identification and Detection Limit Verification

| | | Yes | No | NA |
|-----|---|-----|----|----|
| 9.1 | Are any target compounds detected in the field samples? If Yes, are all positive identifications confirmed in second column? Apply J flag if RPD >40% between first and second columns. | | | X |
| 9.2 | Do detection limits meet those required by the project QAPP and were they properly adjusted for dilution factors and moisture (including adjustment of wet weight aliquot)? | X | | |

Notes:

10.0 Data Completeness

| | | Yes | No | NA |
|--------|---|-----|----|----|
| 10.1 | Is % completeness within the control limits? (Control limit 95% _{aq} and 90% _{so}) | X | | |
| 10.1.1 | Number of samples: <u>6</u> | | | |
| 10.1.2 | Number of target compounds in each analysis: <u>17</u> | | | |
| 10.1.3 | Number of results rejected or not reported: <u>3</u> | | | |
| | % Completeness = $(10.1.1 \times 10.1.2 - 10.1.3) \times 100 / (10.1.1 \times 10.1.2)$ | | | |
| | % Completeness = <u>97.1%</u> | | | |

| Vial | Sample | Dilution | File Number |
|---------|-------------|----------|-------------|
| P1-A-02 | 1207*2500 | 1 | A000003.D |
| P1-A-02 | 1206*1250 | 1 | A000004.D |
| P1-A-02 | 1205*1000 | 1 | A000005.D |
| P1-A-03 | 1204*500 | 1 | A000006.D |
| P1-A-03 | 1203*250 | 1 | A000007.D |
| P1-A-03 | 1202*125 | 1 | A000008.D |
| P1-A-03 | 1201*50 | 1 | A000009.D |
| P1-A-04 | 1600*1000 | 1 | A000010.D |
| P1-E-07 | 21807123417 | 1 | A000011.D |
| P1-E-08 | 1831319 | 1 | A000012.D |
| P1-E-09 | 1831318 | 1 | A000013.D |
| P1-F-02 | 21807123418 | 1 | A000014.D |
| P1-F-03 | 21807123419 | 1 | A000015.D |
| P1-F-04 | 21807123420 | 1 | A000016.D |
| P1-F-05 | 21807123421 | 1 | A000017.D |
| P1-F-06 | 21807123422 | 1 | A000018.D |
| P1-F-07 | 1831320 | 1 | A000019.D |
| P1-F-08 | 1831314 | 1 | A000020.D |
| P1-F-09 | 1831315 | 1 | A000021.D |
| P1-A-01 | 1400*1000 | 1 | A000022.D |
| P2-A-01 | 1831831 | 1 | A000023.D |
| P2-A-02 | 1831832 | 1 | A000024.D |
| P2-A-03 | 1831833 | 1 | A000025.D |
| P2-A-04 | 21807132408 | 1 | A000026.D |
| P2-A-05 | 21807132410 | 1 | A000027.D |
| P2-A-06 | 21807132411 | 1 | A000028.D |
| P2-A-07 | 21807132412 | 1 | A000029.D |
| P2-A-08 | 21807132413 | 1 | A000030.D |
| P2-A-09 | 1831834 | 1 | A000031.D |
| P2-B-01 | 1831835 | 1 | A000032.D |
| P2-B-02 | 21807132414 | 1 | A000033.D |
| P2-B-03 | 21807132415 | 1 | A000034.D |
| P2-B-04 | 21807132416 | 1 | A000035.D |
| P1-A-01 | 1400*1000 | 1 | A000036.D |
| P2-B-05 | 21807132417 | 1 | A000037.D |
| P2-B-06 | 1831836 | 1 | A000038.D |
| P2-B-07 | 1831837 | 1 | A000039.D |
| P2-B-08 | 21807132418 | 5 | A000040.D |
| P2-B-09 | 21807132419 | 1 | A000041.D |
| P2-C-01 | 21807132421 | 5 | A000042.D |
| P2-C-02 | 1831838 | 1 | A000043.D |
| P2-C-03 | 1831832 | 1 | A000044.D |
| P2-C-04 | 1831833 | 1 | A000045.D |
| P1-A-01 | 1400*1000 | 1 | A000046.D |

Analyst MEG
Method ARCI
CCV 005-45-6
RunDate 7/24/2018

ORGANICS INITIAL CALIBRATION DATA

| | | | | | | | | | | | | | | | |
|----------------|-----------|---------|------|--------------------|-----------|-------------------------|--------------------------|--------------------------|--------------------------|---------------------------|---------------------------|-----|---|----------|--|
| Report No: | 218081812 | ID | (mm) | Instrument ID: | HPLC3 | GCALID - FileID - Conc | | | | | | | | | |
| GC Column: | | 1201 | 1202 | 1203 | 1204 | 1205 | 1206 | 1207 | 1208 | 1209 | RF/b/A | m/B | C | FIT TYPE | |
| Calib. Date 1: | 07/24/18 | Time 1: | 1355 | Analyst: | MEG | 1201 ~ 2180724VA09 ~ 50 | 1202 ~ 2180724VA08 ~ 125 | 1203 ~ 2180724VA07 ~ 250 | 1204 ~ 2180724VA06 ~ 500 | 1205 ~ 2180724VA05 ~ 1000 | 1206 ~ 2180724VA04 ~ 1250 | | | | |
| Calib. Date 2: | 07/24/18 | Time 2: | 1552 | Analytical Batch: | 640653 | | | | | | | | | | |
| | | | | Analytical Method: | EPA 8330B | | | | | | | | | | |

| ANALYTE | 1201 | 1202 | 1203 | 1204 | 1205 | 1206 | 1207 | 1208 | 1209 | RF/b/A | m/B | C | FIT TYPE |
|------------------------------|------|------|------|------|------|------|------|------|------|--------|-----|---|----------|
| 1,2-Dinitrobenzene | 16.7 | 16.6 | 16.5 | 17.4 | 18.6 | 18.5 | 17.5 | | | 17.8 | | | 11.78 A |
| 1,3,5-Trinitrobenzene | 10.9 | 11.5 | 11.4 | 11.1 | 12.4 | 12.4 | 11.9 | | | 11.9 | | | 11.66 A |
| 1,3-Dinitrobenzene | 11.5 | 12.0 | 11.4 | 10.7 | 12.8 | 12.7 | 12.5 | | | 12.2 | | | 12.12 A |
| 2,4,6-Trinitrotoluene | 17.2 | 16.3 | 16.4 | 16.5 | 18.2 | 18.1 | 17.9 | | | 17.6 | | | 10.85 A |
| 2,4-Dinitrotoluene | 12.1 | 12.0 | 12.2 | 13.6 | 14.1 | 14.0 | 13.6 | | | 13.5 | | | 12.38 A |
| 2,6-Dinitrotoluene | 22.7 | 25.0 | 25.1 | 21.1 | 26.4 | 26.8 | 27.1 | | | 25.3 | | | 13.72 A |
| 2-Amino-4,6-dinitrotoluene | 19.0 | 20.9 | 20.4 | 18.9 | 23.5 | 23.4 | 23.2 | | | 21.9 | | | 14.61 A |
| 2-Nitrotoluene | 13.3 | 13.1 | 13.5 | 14.0 | 15.1 | 15.1 | 15.3 | | | 14.6 | | | 11.90 A |
| 3,5-Dinitroaniline | 11.1 | 10.7 | 10.7 | 11.2 | 11.9 | 11.8 | 11.8 | | | 11.6 | | | 10.28 A |
| 3-Nitrotoluene | 11.7 | 11.5 | 11.4 | 12.0 | 13.0 | 12.7 | 13.1 | | | 12.5 | | | 10.60 A |
| 4-Amino-2,6-dinitrotoluene | 28.6 | 32.3 | 30.0 | 35.0 | 33.3 | 33.1 | 33.0 | | | 32.9 | | | 11.12 A |
| 4-Nitrotoluene | 21.1 | 23.5 | 24.3 | 22.2 | 23.0 | 23.7 | 23.8 | | | 23.6 | | | 10.02 A |
| HMX | 36.9 | 37.6 | 39.7 | 40.7 | 44.2 | 43.6 | 43.9 | | | 42.1 | | | 12.44 A |
| Nitrobenzene | 18.3 | 19.7 | 19.6 | 19.8 | 21.6 | 21.6 | 21.5 | | | 20.8 | | | 11.95 A |
| Nitroglycerin | 29.6 | 29.3 | 30.7 | 31.0 | 34.1 | 34.1 | 34.4 | | | 32.7 | | | 12.34 A |
| Pentaerythritol Tetranitrate | 26.2 | 27.6 | 27.5 | 28.4 | 30.3 | 30.6 | 32.0 | | | 29.7 | | | 11.85 A |
| RDX | 29.5 | 30.1 | 31.0 | 32.0 | 34.6 | 34.6 | 35.0 | | | 33.3 | | | 12.36 A |
| Tetryl | 21.6 | 23.2 | 23.6 | 23.3 | 25.6 | 25.8 | 25.5 | | | 24.4 | | | 12.87 A |

FIT = %RSD For Average Curve And Calibration Coefficient For Linear And Quadratic
 Curve Types: A - Averged, L - Linear Regression, W - Weighted Linear, Q - Quadratic
 For curve types L and Q, the RRF and RSP (Response) are shown on separate lines to allow for evaluation against minimum RRF

FORM V I SVOA

ORGANICS INITIAL CALIBRATION VERIFICATION

Report No: 218081812 Instrument ID: HPLC3
 Analysis Date: 07/24/18 1612 Lab File ID: 2180724\A10
 Analytical Method: EPA 8330B Analytical Batch: 640653

| <i>ANALYTE</i> | <i>UNITS</i> | <i>TRUE</i> | <i>FOUND</i> | <i>% REC</i> | <i>LCL</i> | <i>UCL</i> | <i>Q</i> |
|------------------------------|--------------|-------------|--------------|--------------|------------|------------|----------|
| 1,3,5-Trinitrobenzene | ug/L | 1000 | 1060 | 106 | 80 | 120 | |
| 1,3-Dinitrobenzene | ug/L | 1000 | 1050 | 105 | 80 | 120 | |
| 2,4,6-Trinitrotoluene | ug/L | 1000 | 1060 | 106 | 80 | 120 | |
| 2,4-Dinitrotoluene | ug/L | 1000 | 1070 | 107 | 80 | 120 | |
| 2,6-Dinitrotoluene | ug/L | 1000 | 1050 | 105 | 80 | 120 | |
| 2-Amino-4,6-dinitrotoluene | ug/L | 1000 | 1030 | 103 | 80 | 120 | |
| 2-Nitrotoluene | ug/L | 1000 | 1060 | 106 | 80 | 120 | |
| 3,5-Dinitroaniline | ug/L | 1000 | 1060 | 106 | 80 | 120 | |
| 3-Nitrotoluene | ug/L | 1000 | 1060 | 106 | 80 | 120 | |
| 4-Amino-2,6-dinitrotoluene | ug/L | 1000 | 1100 | 110 | 80 | 120 | |
| 4-Nitrotoluene | ug/L | 1000 | 1100 | 110 | 80 | 120 | |
| HMX | ug/L | 1000 | 1070 | 107 | 80 | 120 | |
| Nitrobenzene | ug/L | 1000 | 1080 | 108 | 80 | 120 | |
| Nitroglycerin | ug/L | 1000 | 1050 | 105 | 80 | 120 | |
| Pentaerythritol Tetranitrate | ug/L | 1000 | 1030 | 103 | 80 | 120 | |
| RDX | ug/L | 1000 | 1050 | 105 | 80 | 120 | |
| Tetryl | ug/L | 1000 | 1010 | 101 | 80 | 120 | |

FORM 6I - ORG

| Vial | Sample | Dilution | File Number |
|---------|-------------|----------|-------------|
| Vial 2 | 1400*1000 | 1 | A000003.D |
| P2-A-01 | 1843189 | 1 | A000004.D |
| P2-A-02 | 1843190 | 1 | A000005.D |
| P2-A-03 | 1843191 | 1 | A000006.D |
| P2-A-04 | 21808241101 | 1 | A000007.D |
| P2-A-05 | 1843193 | 1 | A000008.D |
| P2-A-06 | 1843192 | 1 | A000009.D |
| P2-A-07 | 1843194 | 1 | A000010.D |
| P2-A-08 | 1843195 | 1 | A000011.D |
| P2-A-09 | 1843196 | 1 | A000012.D |
| Vial 2 | 1400*1000 | 1 | A000013.D |
| P2-B-01 | 21808172601 | 1 | A000014.D |
| P2-B-02 | 21808172602 | 1 | A000015.D |
| P2-B-03 | 21808172603 | 1 | A000016.D |
| P2-B-04 | 21808172604 | 1 | A000017.D |
| Vial 2 | 1400*1000 | 1 | A000018.D |
| P2-B-06 | 1843101 | 1 | A000019.D |
| P2-B-07 | 1843102 | 1 | A000020.D |
| P2-B-08 | 1843103 | 1 | A000021.D |
| P2-B-09 | 21808181401 | 1 | A000022.D |
| P2-C-01 | 21808181402 | 1 | A000023.D |
| P2-C-02 | 21808181403 | 1 | A000024.D |
| P2-C-03 | 21808181404 | 1 | A000025.D |
| P2-C-04 | 21808181407 | 1 | A000026.D |
| P2-C-05 | 1843105 | 1 | A000027.D |
| P2-C-06 | 1843104 | 1 | A000028.D |
| P2-C-07 | 1843106 | 1 | A000029.D |
| P2-C-08 | 1843214 | 1 | A000030.D |
| P2-C-09 | 1843215 | 1 | A000031.D |
| Vial 2 | 1400*1000 | 1 | A000032.D |

Analyst MEG
Method ARCI
CCV 005-49-4
RunDate 8/30/2018

7E
ORGANICS CONTINUING CALIBRATION CHECK

| | | | |
|----------------------|-------------------------------------|--------------------|-------------------|
| Report No: | <u>218081812</u> | CCAL ID: | <u>1400</u> |
| GC Column: | <u>ARC18</u> ID <u>3</u> (mm) | Instrument ID: | <u>HPLC3</u> |
| Injection Vol.: | <u>1.0</u> (µL) | Lab File ID: | <u>2180830A03</u> |
| Init. Calib. Date 1: | <u>07/24/18</u> Time 1: <u>1355</u> | Analyst: | <u>MEG</u> |
| Init. Calib. Date 2: | <u>07/24/18</u> Time 2: <u>1552</u> | Analytical Batch: | <u>643050</u> |
| Analysis Date: | <u>08/30/18</u> Time: <u>1116</u> | Analytical Method: | <u>EPA 8330B</u> |

| ANALYTE | TRUE | CONC | RRF | RRF CCV | Min RRF | %D/%Drift | Max %D/ %Drift | TYPE | Q |
|------------------------------|------|-------|------|---------|---------|-----------|-------------------|------|---|
| 1,2-Dinitrobenzene | 1000 | 1156 | 17.8 | 15.2 | NA | -14.6 | 20 | A | |
| 1,3,5-Trinitrobenzene | 1000 | 1037 | 11.9 | 11.4 | NA | -4.2 | 20 | A | |
| 1,3-Dinitrobenzene | 1000 | 995.7 | 12.2 | 12.1 | NA | -8 | 20 | A | |
| 2,4,6-Trinitrotoluene | 1000 | 1061 | 17.6 | 16.5 | NA | -6.3 | 20 | A | |
| 2,4-Dinitrotoluene | 1000 | 986.4 | 13.5 | 13.5 | NA | 0 | 20 | A | |
| 2,6-Dinitrotoluene | 1000 | 976.3 | 25.3 | 25.6 | NA | 1.2 | 20 | A | |
| 2-Amino-4,6-dinitrotoluene | 1000 | 960.0 | 21.9 | 22.4 | NA | 2.3 | 20 | A | |
| 2-Nitrotoluene | 1000 | 866.8 | 14.6 | 16.6 | NA | 13.7 | 20 | A | |
| 3,5-Dinitroaniline | 1000 | 970.9 | 11.6 | 11.9 | NA | 2.6 | 20 | A | |
| 3-Nitrotoluene | 1000 | 890.7 | 12.5 | 13.9 | NA | 11.2 | 20 | A | |
| 4-Amino-2,6-dinitrotoluene | 1000 | 1033 | 32.9 | 31.6 | NA | -4 | 20 | A | |
| 4-Nitrotoluene | 1000 | 921.0 | 23.6 | 25.4 | NA | 7.6 | 20 | A | |
| HMX | 1000 | 958.3 | 42.1 | 43.4 | NA | 3.1 | 20 | A | |
| Nitrobenzene | 1000 | 858.6 | 20.8 | 23.9 | NA | 14.9 | 20 | A | |
| Nitroglycerin | 1000 | 989.0 | 32.7 | 32.7 | NA | 0 | 20 | A | |
| Pentaerythritol Tetranitrate | 1000 | 1021 | 29.7 | 28.8 | NA | -3 | 20 | A | |
| RDX | 1000 | 961.8 | 33.3 | 34.2 | NA | 2.7 | 20 | A | |
| Tetryl | 1000 | 972.3 | 24.4 | 24.8 | NA | 1.6 | 20 | A | |

FORM V II SV

ORGANICS CONTINUING CALIBRATION CHECK

Report No: 218081812 CCAL ID: 1400
 GC Column: ARC18 ID 3 (mm) Instrument ID: HPLC3
 Injection Vol.: 1.0 (µL) Lab File ID: 2180830A13
 Init. Calib. Date 1: 07/24/18 Time 1: 1355 Analyst: MEG
 Init. Calib. Date 2: 07/24/18 Time 2: 1552 Analytical Batch: 643050
 Analysis Date: 08/30/18 Time: 1443 Analytical Method: EPA 8330B

| ANALYTE | TRUE | CONC | RRF | RRF CCV | Min RRF | %D/%Drift | Max %D/ %Drift | TYPE | Q |
|------------------------------|------|-------|------|---------|---------|-----------|-------------------|------|---|
| 1,2-Dinitrobenzene | 1000 | 1108 | 17.8 | 15.9 | NA | -10.7 | 20 | A | |
| 1,3,5-Trinitrobenzene | 1000 | 992.1 | 11.9 | 11.9 | NA | 0 | 20 | A | |
| 1,3-Dinitrobenzene | 1000 | 952.0 | 12.2 | 12.7 | NA | 4.1 | 20 | A | |
| 2,4,6-Trinitrotoluene | 1000 | 965.8 | 17.6 | 18.1 | NA | 2.8 | 20 | A | |
| 2,4-Dinitrotoluene | 1000 | 937.7 | 13.5 | 14.2 | NA | 5.2 | 20 | A | |
| 2,6-Dinitrotoluene | 1000 | 933.9 | 25.3 | 26.7 | NA | 5.5 | 20 | A | |
| 2-Amino-4,6-dinitrotoluene | 1000 | 906.6 | 21.9 | 23.7 | NA | 8.2 | 20 | A | |
| 2-Nitrotoluene | 1000 | 935.7 | 14.6 | 15.4 | NA | 5.5 | 20 | A | |
| 3,5-Dinitroaniline | 1000 | 929.8 | 11.6 | 12.4 | NA | 6.9 | 20 | A | |
| 3-Nitrotoluene | 1000 | 953.7 | 12.5 | 13.0 | NA | 4 | 20 | A | |
| 4-Amino-2,6-dinitrotoluene | 1000 | 993.7 | 32.9 | 32.8 | NA | -3 | 20 | A | |
| 4-Nitrotoluene | 1000 | 991.7 | 23.6 | 23.6 | NA | 0 | 20 | A | |
| HMX | 1000 | 916.9 | 42.1 | 45.3 | NA | 7.6 | 20 | A | |
| Nitrobenzene | 1000 | 928.8 | 20.8 | 22.1 | NA | 6.3 | 20 | A | |
| Nitroglycerin | 1000 | 945.8 | 32.7 | 34.2 | NA | 4.6 | 20 | A | |
| Pentaerythritol Tetranitrate | 1000 | 1008 | 29.7 | 29.2 | NA | -1.7 | 20 | A | |
| RDX | 1000 | 924.8 | 33.3 | 35.6 | NA | 6.9 | 20 | A | |
| Tetryl | 1000 | 1021 | 24.4 | 23.6 | NA | -3.3 | 20 | A | |

7E
ORGANICS CONTINUING CALIBRATION CHECK

| | | | |
|----------------------|-------------------------------------|--------------------|--------------------|
| Report No: | <u>218081814</u> | CCAL ID: | <u>1400</u> |
| GC Column: | <u>ARC18</u> ID <u>3</u> (mm) | Instrument ID: | <u>HPLC3</u> |
| Injection Vol.: | <u>1.0</u> (µL) | Lab File ID: | <u>2180830VA18</u> |
| Init. Calib. Date 1: | <u>07/24/18</u> Time 1: <u>1355</u> | Analyst: | <u>MEG</u> |
| Init. Calib. Date 2: | <u>07/24/18</u> Time 2: <u>1552</u> | Analytical Batch: | <u>643050</u> |
| Analysis Date: | <u>08/30/18</u> Time: <u>1622</u> | Analytical Method: | <u>EPA 8330B</u> |

| ANALYTE | TRUE | CONC | RRF | RRF CCV | Min RRF | %D/%Drift | Max %D/ %Drift | TYPE | Q |
|------------------------------|------|-------|------|---------|---------|-----------|-------------------|------|---|
| 1,2-Dinitrobenzene | 1000 | 1112 | 17.8 | 15.8 | NA | -11.2 | 20 | A | |
| 1,3,5-Trinitrobenzene | 1000 | 998.5 | 11.9 | 11.8 | NA | -8 | 20 | A | |
| 1,3-Dinitrobenzene | 1000 | 957.2 | 12.2 | 12.6 | NA | 3.3 | 20 | A | |
| 2,4,6-Trinitrotoluene | 1000 | 983.1 | 17.6 | 17.8 | NA | 1.1 | 20 | A | |
| 2,4-Dinitrotoluene | 1000 | 938.4 | 13.5 | 14.2 | NA | 5.2 | 20 | A | |
| 2,6-Dinitrotoluene | 1000 | 942.0 | 25.3 | 26.5 | NA | 4.7 | 20 | A | |
| 2-Amino-4,6-dinitrotoluene | 1000 | 904.0 | 21.9 | 23.8 | NA | 8.7 | 20 | A | |
| 2-Nitrotoluene | 1000 | 961.0 | 14.6 | 15.0 | NA | 2.7 | 20 | A | |
| 3,5-Dinitroaniline | 1000 | 919.0 | 11.6 | 12.5 | NA | 7.8 | 20 | A | |
| 3-Nitrotoluene | 1000 | 958.6 | 12.5 | 12.9 | NA | 3.2 | 20 | A | |
| 4-Amino-2,6-dinitrotoluene | 1000 | 969.2 | 32.9 | 33.6 | NA | 2.1 | 20 | A | |
| 4-Nitrotoluene | 1000 | 1014 | 23.6 | 23.1 | NA | -2.1 | 20 | A | |
| HMX | 1000 | 916.4 | 42.1 | 45.3 | NA | 7.6 | 20 | A | |
| Nitrobenzene | 1000 | 937.6 | 20.8 | 21.9 | NA | 5.3 | 20 | A | |
| Nitroglycerin | 1000 | 941.9 | 32.7 | 34.3 | NA | 4.9 | 20 | A | |
| Pentaerythritol Tetranitrate | 1000 | 990.1 | 29.7 | 29.7 | NA | 0 | 20 | A | |
| RDX | 1000 | 901.8 | 33.3 | 36.5 | NA | 9.6 | 20 | A | |
| Tetryl | 1000 | 990.2 | 24.4 | 24.4 | NA | 0 | 20 | A | |

FORM V II SV

7E
ORGANICS CONTINUING CALIBRATION CHECK

| | | | |
|----------------------|-------------------------------------|--------------------|--------------------|
| Report No: | <u>218081814</u> | CCAL ID: | <u>1400</u> |
| GC Column: | <u>ARC18</u> ID <u>3</u> (mm) | Instrument ID: | <u>HPLC3</u> |
| Injection Vol.: | <u>1.0</u> (µL) | Lab File ID: | <u>2180830VA32</u> |
| Init. Calib. Date 1: | <u>07/24/18</u> Time 1: <u>1355</u> | Analyst: | <u>MEG</u> |
| Init. Calib. Date 2: | <u>07/24/18</u> Time 2: <u>1552</u> | Analytical Batch: | <u>643050</u> |
| Analysis Date: | <u>08/30/18</u> Time: <u>2058</u> | Analytical Method: | <u>EPA 8330B</u> |

| ANALYTE | TRUE | CONC | RRF | RRF CCV | Min RRF | %D/%Drift | Max %D/ | | TYPE | Q |
|------------------------------|------|-------|------|---------|---------|-----------|---------|--|------|---|
| | | | | | | | %Drift | | | |
| 1,2-Dinitrobenzene | 1000 | 1141 | 17.8 | 15.4 | NA | -13.5 | 20 | | A | |
| 1,3,5-Trinitrobenzene | 1000 | 1003 | 11.9 | 11.8 | NA | -.8 | 20 | | A | |
| 1,3-Dinitrobenzene | 1000 | 979.0 | 12.2 | 12.3 | NA | .8 | 20 | | A | |
| 2,4,6-Trinitrotoluene | 1000 | 976.1 | 17.6 | 17.9 | NA | 1.7 | 20 | | A | |
| 2,4-Dinitrotoluene | 1000 | 892.0 | 13.5 | 14.9 | NA | 10.4 | 20 | | A | |
| 2,6-Dinitrotoluene | 1000 | 1012 | 25.3 | 24.6 | NA | -2.8 | 20 | | A | |
| 2-Amino-4,6-dinitrotoluene | 1000 | 919.1 | 21.9 | 23.4 | NA | 6.8 | 20 | | A | |
| 2-Nitrotoluene | 1000 | 937.8 | 14.6 | 15.4 | NA | 5.5 | 20 | | A | |
| 3,5-Dinitroaniline | 1000 | 925.7 | 11.6 | 12.4 | NA | 6.9 | 20 | | A | |
| 3-Nitrotoluene | 1000 | 925.8 | 12.5 | 13.4 | NA | 7.2 | 20 | | A | |
| 4-Amino-2,6-dinitrotoluene | 1000 | 969.1 | 32.9 | 33.6 | NA | 2.1 | 20 | | A | |
| 4-Nitrotoluene | 1000 | 963.8 | 23.6 | 24.3 | NA | 3 | 20 | | A | |
| HMX | 1000 | 919.8 | 42.1 | 45.2 | NA | 7.4 | 20 | | A | |
| Nitrobenzene | 1000 | 914.0 | 20.8 | 22.5 | NA | 8.2 | 20 | | A | |
| Nitroglycerin | 1000 | 939.2 | 32.7 | 34.4 | NA | 5.2 | 20 | | A | |
| Pentaerythritol Tetranitrate | 1000 | 958.4 | 29.7 | 30.7 | NA | 3.4 | 20 | | A | |
| RDX | 1000 | 923.1 | 33.3 | 35.6 | NA | 6.9 | 20 | | A | |
| Tetryl | 1000 | 981.3 | 24.4 | 24.6 | NA | .8 | 20 | | A | |

FORM V II SV

| Vial | Sample | Dilution | File Number |
|---------|-------------|----------|-------------|
| Vial 2 | 1400*1000 | 1 | A000004.D |
| P2-A-01 | 1845522 | 1 | A000010.D |
| P2-A-02 | 1845523 | 1 | A000011.D |
| P2-A-03 | 1845524 | 1 | A000012.D |
| P2-A-04 | 21808311301 | 1 | A000013.D |
| P2-A-05 | 1845525 | 1 | A000014.D |
| P2-A-06 | 1845526 | 1 | A000015.D |
| P2-A-07 | 1845527 | 1 | A000016.D |
| P2-A-08 | 1845523 | 1 | A000017.D |
| P2-A-09 | 1845524 | 1 | A000021.D |
| P2-B-09 | 21808220401 | 1 | A000022.D |
| Vial 2 | 1400*1000 | 1 | A000023.D |
| P2-B-01 | 21808241301 | 1 | A000024.D |
| P2-B-02 | 21808241302 | 1 | A000025.D |
| P2-B-04 | 21808242301 | 1 | A000026.D |
| P2-B-05 | 1843500 | 1 | A000027.D |
| P2-B-06 | 1843499 | 1 | A000028.D |
| P2-B-07 | 21808242302 | 1 | A000029.D |
| P2-B-08 | 21808242303 | 1 | A000030.D |
| Vial 2 | 1400*1000 | 1 | A000031.D |
| P2-C-01 | 21808240301 | 1 | A000032.D |
| P2-C-02 | 21808240302 | 1 | A000033.D |
| P2-C-03 | 21808240303 | 1 | A000034.D |
| P2-C-04 | 21808240304 | 1 | A000035.D |
| P2-C-05 | 21808240305 | 1 | A000036.D |
| P2-C-06 | 21808240306 | 1 | A000037.D |
| P2-C-07 | 21808240307 | 1 | A000038.D |
| P2-C-08 | 21808240308 | 1 | A000039.D |
| P2-D-01 | 21808172301 | 1 | A000040.D |
| Vial 2 | 1400*1000 | 1 | A000041.D |
| P2-D-02 | 21808172302 | 1 | A000042.D |
| P2-D-03 | 21808172303 | 1 | A000043.D |
| P2-D-04 | 21808172304 | 1 | A000044.D |
| P2-D-05 | 21808172305 | 1 | A000045.D |
| P2-D-06 | 21808172306 | 1 | A000046.D |
| P2-D-07 | 21808172307 | 1 | A000047.D |
| P2-D-08 | 21808172308 | 1 | A000048.D |
| P2-D-09 | 21808172309 | 1 | A000049.D |
| P2-E-01 | 21808172310 | 1 | A000050.D |
| P2-E-02 | 21808172311 | 1 | A000051.D |
| Vial 2 | 1400*1000 | 1 | A000052.D |
| P2-E-03 | 21808172312 | 1 | A000053.D |
| P2-E-04 | 21808172313 | 1 | A000054.D |
| P2-E-05 | 21808172314 | 1 | A000055.D |
| P2-E-06 | 21808172315 | 1 | A000056.D |
| P2-E-07 | 21808172316 | 1 | A000057.D |

| | | | |
|---------|-------------|---|-----------|
| P2-E-08 | 21808172317 | 1 | A000058.D |
| P2-E-09 | 21808172318 | 1 | A000059.D |
| P2-F-01 | 21808172319 | 1 | A000060.D |
| P2-F-02 | 21808172320 | 1 | A000061.D |
| P2-F-03 | 21808172321 | 1 | A000062.D |
| Vial 2 | 1400*1000 | 1 | A000063.D |
| P1-A-01 | 21808181206 | 1 | A000064.D |
| P1-A-02 | 21808181207 | 1 | A000065.D |
| P1-A-03 | 21808181208 | 1 | A000066.D |
| P1-A-04 | 21808181209 | 1 | A000067.D |
| P1-A-05 | 21808181210 | 1 | A000068.D |
| P1-A-06 | 1846370 | 1 | A000069.D |
| P1-A-07 | 1846371 | 1 | A000070.D |
| P1-A-08 | 1846372 | 1 | A000071.D |
| P1-A-09 | 21808181213 | 1 | A000072.D |
| P1-B-01 | 21808181214 | 1 | A000073.D |
| P1-B-02 | 21808181215 | 1 | A000074.D |
| P1-B-03 | 21808181216 | 1 | A000075.D |
| P1-B-04 | 21808181217 | 1 | A000076.D |
| P1-B-05 | 1846373 | 1 | A000077.D |
| P1-B-06 | 1846374 | 1 | A000078.D |
| P1-B-07 | 1846375 | 1 | A000079.D |
| P1-B-08 | 1846376 | 1 | A000080.D |
| P1-B-09 | 1846377 | 1 | A000081.D |
| Vial 2 | 1400*1000 | 1 | A000082.D |
| P2-F-06 | 1845471 | 1 | A000083.D |
| P2-F-07 | 1845472 | 1 | A000084.D |
| P2-F-08 | 1845473 | 1 | A000085.D |
| P2-F-09 | 21808302601 | 1 | A000086.D |
| Vial 2 | 1400*1000 | 1 | A000087.D |

Analyst MEG
Method ARCI
CCV 005-49-9
RunDate 9/12/2018

ORGANICS CONTINUING CALIBRATION CHECK

Report No: 218081812 CCAL ID: 1400
 GC Column: ARC18 ID 3 (mm) Instrument ID: HPLC3
 Injection Vol.: 1.0 (µL) Lab File ID: 2180912\A63
 Init. Calib. Date 1: 07/24/18 Time 1: 1355 Analyst: MEG
 Init. Calib. Date 2: 07/24/18 Time 2: 1552 Analytical Batch: 643776
 Analysis Date: 09/13/18 Time: 0520 Analytical Method: EPA 8330B

| ANALYTE | TRUE | CONC | RRF | RRF CCV | Min RRF | %D/%Drift | Max %D/ | | TYPE | Q |
|------------------------------|------|-------|------|---------|---------|-----------|---------|--|------|---|
| | | | | | | | %Drift | | | |
| 1,2-Dinitrobenzene | 1000 | 1168 | 17.8 | 15.1 | NA | -15.2 | 20 | | A | |
| 1,3,5-Trinitrobenzene | 1000 | 1128 | 11.9 | 10.5 | NA | -11.8 | 20 | | A | |
| 1,3-Dinitrobenzene | 1000 | 1083 | 12.2 | 11.2 | NA | -8.2 | 20 | | A | |
| 2,4,6-Trinitrotoluene | 1000 | 1150 | 17.6 | 15.2 | NA | -13.6 | 20 | | A | |
| 2,4-Dinitrotoluene | 1000 | 992.5 | 13.5 | 13.4 | NA | -.7 | 20 | | A | |
| 2,6-Dinitrotoluene | 1000 | 1059 | 25.3 | 23.6 | NA | -6.7 | 20 | | A | |
| 2-Amino-4,6-dinitrotoluene | 1000 | 980.6 | 21.9 | 21.9 | NA | 0 | 20 | | A | |
| 2-Nitrotoluene | 1000 | 954.5 | 14.6 | 15.1 | NA | 3.4 | 20 | | A | |
| 3,5-Dinitroaniline | 1000 | 1017 | 11.6 | 11.3 | NA | -2.6 | 20 | | A | |
| 3-Nitrotoluene | 1000 | 899.2 | 12.5 | 13.8 | NA | 10.4 | 20 | | A | |
| 4-Amino-2,6-dinitrotoluene | 1000 | 1117 | 32.9 | 29.2 | NA | -11.2 | 20 | | A | |
| 4-Nitrotoluene | 1000 | 882.8 | 23.6 | 26.5 | NA | 12.3 | 20 | | A | |
| HMX | 1000 | 1035 | 42.1 | 40.1 | NA | -4.8 | 20 | | A | |
| Nitrobenzene | 1000 | 939.4 | 20.8 | 21.9 | NA | 5.3 | 20 | | A | |
| Nitroglycerin | 1000 | 1071 | 32.7 | 30.2 | NA | -7.6 | 20 | | A | |
| Pentaerythritol Tetranitrate | 1000 | 1132 | 29.7 | 26.0 | NA | -12.5 | 20 | | A | |
| RDX | 1000 | 1024 | 33.3 | 32.1 | NA | -3.6 | 20 | | A | |
| Tetryl | 1000 | 935.7 | 24.4 | 25.8 | NA | 5.7 | 20 | | A | |

FORM V II SV

ORGANICS CONTINUING CALIBRATION CHECK

Report No: 218081812 CCAL ID: 1400
 GC Column: ARC18 ID 3 (mm) Instrument ID: HPLC3
 Injection Vol.: 1.0 (µL) Lab File ID: 2180912\A82
 Init. Calib. Date 1: 07/24/18 Time 1: 1355 Analyst: MEG
 Init. Calib. Date 2: 07/24/18 Time 2: 1552 Analytical Batch: 643776
 Analysis Date: 09/13/18 Time: 1135 Analytical Method: EPA 8330B

| ANALYTE | TRUE | CONC | RRF | RRF CCV | Min RRF | %D/%Drift | Max %D/ %Drift | TYPE | Q |
|------------------------------|------|-------|------|---------|---------|-----------|-------------------|------|---|
| 1,2-Dinitrobenzene | 1000 | 1130 | 17.8 | 15.6 | NA | -12.4 | 20 | A | |
| 1,3,5-Trinitrobenzene | 1000 | 1083 | 11.9 | 10.9 | NA | -8.4 | 20 | A | |
| 1,3-Dinitrobenzene | 1000 | 1081 | 12.2 | 11.2 | NA | -8.2 | 20 | A | |
| 2,4,6-Trinitrotoluene | 1000 | 1108 | 17.6 | 15.8 | NA | -10.2 | 20 | A | |
| 2,4-Dinitrotoluene | 1000 | 921.8 | 13.5 | 14.4 | NA | 6.7 | 20 | A | |
| 2,6-Dinitrotoluene | 1000 | 1037 | 25.3 | 24.1 | NA | -4.7 | 20 | A | |
| 2-Amino-4,6-dinitrotoluene | 1000 | 987.5 | 21.9 | 21.8 | NA | -.5 | 20 | A | |
| 2-Nitrotoluene | 1000 | 969.2 | 14.6 | 14.9 | NA | 2.1 | 20 | A | |
| 3,5-Dinitroaniline | 1000 | 1037 | 11.6 | 11.1 | NA | -4.3 | 20 | A | |
| 3-Nitrotoluene | 1000 | 827.4 | 12.5 | 15.0 | NA | 20 | 20 | A | |
| 4-Amino-2,6-dinitrotoluene | 1000 | 1139 | 32.9 | 28.6 | NA | -13.1 | 20 | A | |
| 4-Nitrotoluene | 1000 | 1033 | 23.6 | 22.7 | NA | -3.8 | 20 | A | |
| HMX | 1000 | 1031 | 42.1 | 40.3 | NA | -4.3 | 20 | A | |
| Nitrobenzene | 1000 | 1078 | 20.8 | 19.1 | NA | -8.2 | 20 | A | |
| Nitroglycerin | 1000 | 1083 | 32.7 | 29.9 | NA | -8.6 | 20 | A | |
| Pentaerythritol Tetranitrate | 1000 | 1100 | 29.7 | 26.7 | NA | -10.1 | 20 | A | |
| RDX | 1000 | 1012 | 33.3 | 32.5 | NA | -2.4 | 20 | A | |
| Tetryl | 1000 | 1115 | 24.4 | 21.7 | NA | -11.1 | 20 | A | |

EPA 8330B REPLICATE SUMMARY

| | | | |
|--------------------|----------------------------|-------------------|------------------------|
| Report No: | <u>218081812</u> | Parent Sample ID: | <u>WIL02DA01B (RE)</u> |
| Prep Method: | <u>EPA 8330B</u> | Parent GCAL ID: | <u>21808181217</u> |
| Prep Date: | <u>9/5/2018 6:51:00 PM</u> | Prep Batch: | <u>643342</u> |
| Analytical Method: | <u>EPA 8330B</u> | | |

| <i>ANALYTE</i> | <i>CAS</i> | <i>UNITS</i> | <i>PARENT RESULT</i> | <i>REP #1 RESULT (1846373)</i> | <i>REP #2 RESULT (1846374)</i> | <i>%RSD</i> | <i>#</i> |
|------------------------------|------------|--------------|--------------------------|--|--|-------------|----------|
| 1,3,5-Trinitrobenzene | 99-35-4 | ug/Kg | 0 | 0 | 0 | 0 | |
| 1,3-Dinitrobenzene | 99-65-0 | ug/Kg | 0 | 0 | 0 | 0 | |
| 2,4,6-Trinitrotoluene | 118-96-7 | ug/Kg | 0 | 0 | 0 | 0 | |
| 2,4-Dinitrotoluene | 121-14-2 | ug/Kg | 0 | 0 | 0 | 0 | |
| 2,6-Dinitrotoluene | 606-20-2 | ug/Kg | 0 | 0 | 0 | 0 | |
| 2-Amino-4,6-dinitrotoluene | 35572-78-2 | ug/Kg | 0 | 0 | 0 | 0 | |
| 2-Nitrotoluene | 88-72-2 | ug/Kg | 0 | 0 | 0 | 0 | |
| 3,5-Dinitroaniline | 618-87-1 | ug/Kg | 0 | 0 | 0 | 0 | |
| 3-Nitrotoluene | 99-08-1 | ug/Kg | 0 | 0 | 0 | 0 | |
| 4-Amino-2,6-dinitrotoluene | 19406-51-0 | ug/Kg | 0 | 0 | 0 | 0 | |
| 4-Nitrotoluene | 99-99-0 | ug/Kg | 0 | 0 | 0 | 0 | |
| HMX | 2691-41-0 | ug/Kg | 0 | 0 | 0 | 0 | |
| Nitrobenzene | 98-95-3 | ug/Kg | 0 | 0 | 0 | 0 | |
| Nitroglycerin | 55-63-0 | ug/Kg | 0 | 0 | 0 | 0 | |
| Pentaerythritol Tetranitrate | 78-11-5 | ug/Kg | 0 | 0 | 0 | 0 | |
| RDX | 121-82-4 | ug/Kg | 0 | 0 | 0 | 0 | |
| Tetryl | 479-45-8 | ug/Kg | 0 | 0 | 0 | 0 | |

* - RSD greater than 20%

2F
ORGANIC SURROGATE RECOVERY

Report No: 218081812

Analytical Method: EPA 8330B

| <i>Client Sample ID</i> | <i>GCAL Sample ID</i> | <i>SMC1 #</i> | <i>SMC2 #</i> | <i>SMC3 #</i> | <i>SMC4 #</i> | <i>TOT OUT</i> |
|---------------------------------|-----------------------|---------------|---------------|---------------|---------------|----------------|
| WIL02DA02A | 21808181206 | 106 | | | | 0 |
| WIL02DA02A MS | 21808181207 | 111 | | | | 0 |
| WIL02DA02A MSD | 21808181208 | 105 | | | | 0 |
| WIL02DA01A | 21808181209 | 104 | | | | 0 |
| WIL02DA01B | 21808181210 | 106 | | | | 0 |
| WIL02DA02A (RE) | 21808181213 | 103 | | | | 0 |
| WIL02DA02A MS (RE) | 21808181214 | 98 | | | | 0 |
| WIL02DA02A MSD (RE) | 21808181215 | 112 | | | | 0 |
| WIL02DA01A (RE) | 21808181216 | 97 | | | | 0 |
| WIL02DA01B (RE) | 21808181217 | 96 | | | | 0 |
| MB1843189 | 1843189 | 71 | | | | 0 |
| LCS1843190 | 1843190 | 64 | | | | 0 |
| LCSD1843191 | 1843191 | 45 | * | | | 1 |
| GRBLK for HBN 642698 [EXTO/549] | 1843194 | 83 | | | | 0 |
| MB1846370 | 1846370 | 114 | | | | 0 |
| LCS1846371 | 1846371 | 96 | | | | 0 |
| LCSD1846372 | 1846372 | 100 | | | | 0 |
| WIL02DA01B (RE)DUP | 1846373 | 107 | | | | 0 |
| WIL02DA01B (RE)DUP | 1846374 | 108 | | | | 0 |
| GRBLK for HBN 643342 [EXTO/550] | 1846375 | 107 | | | | 0 |

QC LIMITS

| | | | | |
|---------|--------------------|----|-----|---|
| SMC 1 : | 1,2-Dinitrobenzene | 50 | 150 | # Column to be used to flag recovery limits |
| SMC 2 : | | | | * Value outside of QC limits |
| SMC 3 : | | | | D Surrogate diluted out |
| SMC 4 : | | | | |

FORM II ORG-2

4C
ORGANIC METHOD BLANK SUMMARY

| | |
|------------------------------------|--|
| Report No: <u>218081812</u> | Method Blank ID: <u>1843189</u> |
| Matrix: <u>Solid</u> | Instrument ID: <u>HPLC3</u> |
| Sample Amt: <u>10</u> g | Lab File ID: <u>2180830\A04</u> |
| Injection Vol.: <u>1.0</u> (µL) | GC Column: <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: <u>40000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: <u>08/25/18</u> | Analysis Date: <u>08/30/18</u> Time: <u>1145</u> |
| Prep Batch: <u>642698</u> | Analytical Batch: <u>643050</u> |
| Prep Method: <u>8330B</u> | Analytical Method: <u>EPA 8330B</u> |

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD

| CLIENT SAMPLE ID | GCAL SAMPLE ID | DATE ANALYZED | TIME ANALYZED | INSTRUMENT ID |
|------------------------------------|-------------------|------------------|------------------|------------------|
| 1. LCS1843190 | 1843190 | 08/30/18 | 1205 | HPLC3 |
| 2. LCSD1843191 | 1843191 | 08/30/18 | 1225 | HPLC3 |
| 3. GRBLK for HBN 642698 [EXTO/549] | 1843194 | 08/30/18 | 1344 | HPLC3 |
| 4. WIL02DA02A | 21808181206 | 09/13/18 | 0540 | HPLC3 |
| 5. WIL02DA02A MS | 21808181207 | 09/13/18 | 0559 | HPLC3 |
| 6. WIL02DA02A MSD | 21808181208 | 09/13/18 | 0619 | HPLC3 |
| 7. WIL02DA01A | 21808181209 | 09/13/18 | 0639 | HPLC3 |
| 8. WIL02DA01B | 21808181210 | 09/13/18 | 0659 | HPLC3 |

FORM IV ORGANIC



8330B Explosives Prep Solid



| | | | | | | | |
|--------------|------------|-----------------|---------------------|---------------|----------------------|-------|---------------|
| ANALYST/TECH | <i>DUP</i> | START DATE/TIME | <i>8/25/18 1649</i> | END DATE/TIME | <i>8/26/18 15:02</i> | BATCH | 642698 |
|--------------|------------|-----------------|---------------------|---------------|----------------------|-------|---------------|

| # | CLIENT | TYPE | CLIENT ID | GCAL ID | INITIAL WGT (g) | FINAL VOL (mL) | COMMENT | STANDARDS/ REAGENTS |
|----|--------|-------|----------------------------|-------------|-----------------|----------------|---------|---|
| 1 | QC | MB | MB 1843189 | 1843189 | <i>10.0</i> | <i>40.0</i> | | 8330 Surrogate 4/10ug/ml / Volume 1.0ml |
| 2 | QC | LCS | SRM 1843190 ** Use SRM ** | 1843190 | <i>10.0</i> | <i>40.0</i> | | <i>703-27-9</i> |
| 3 | QC | LCSD | SRMD 1843191 ** Use SRM ** | 1843191 | <i>10.0</i> | <i>40.0</i> | | 8330 Spike 4/10ug/ml / Volume 1.0ml |
| 4 | 4838 | SAMP | WIL02DB02A | 21808181204 | <i>10.1</i> | <i>40.0</i> | | <i>703-23-1</i> |
| 5 | 4838 | SAMP | WIL02DB01A | 21808181205 | <i>10.0</i> | <i>40.0</i> | | Solid Reference Material |
| 6 | 4838 | SAMP | WIL02DA02A | 21808181206 | <i>10.3</i> | <i>40.0</i> | | <i>2107150</i> |
| 7 | 4838 | MS | WIL02DA02A MS | 21808181207 | <i>10.0</i> | <i>40.0</i> | | HPLC Water |
| 8 | 4838 | MSD | WIL02DA02A MSD | 21808181208 | <i>10.5</i> | <i>40.0</i> | | |
| 9 | 4838 | SAMP | WIL02DA01A | 21808181209 | <i>10.5</i> | <i>40.0</i> | | Acetonitrile |
| 10 | 4838 | SAMP | WIL02DA01B | 21808181210 | <i>10.0</i> | <i>40.0</i> | | <i>2127054</i> |
| 11 | 4612 | SAMP | ADA-20X38-A03-SP07... | 21808241101 | <i>10.1</i> | <i>40.0</i> | | Sand |
| 12 | QC | DUP | DUP 1843193 | 1843193 | <i>10.1</i> | <i>40.0</i> | | |
| 13 | QC | DUP | DUP 1843192 | 1843192 | <i>10.5</i> | <i>40.0</i> | | |
| 14 | QC | GRBLK | GRBLK 1843194 | 1843194 | <i>10.0</i> | <i>40.0</i> | | |
| 15 | QC | LCS | SRM 1843195 ** Use SRM ** | 1843195 | <i>10.0</i> | <i>40.0</i> | | |
| 16 | QC | LCSD | SRMD 1843196 ** Use SRM ** | 1843196 | <i>10.0</i> | <i>40.0</i> | | |
| 17 | | | | | | | | |
| 18 | | | | | | | | |
| 19 | | | | | | | | |
| 20 | | | | | | | | |
| 21 | | | | | | | | |
| 22 | | | | | | | | |
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| 24 | | | | | | | | |
| 25 | | | | | | | | |
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| 27 | | | | | | | | |
| 28 | | | | | | | | |
| 29 | | | | | | | | |
| 30 | | | | | | | | |

EQUIPMENT CONDITIONS

| | | | |
|------------|-----------|----------------------|--|
| BALANCE ID | <i>D3</i> | GRINDER ID - PUCK 01 | |
|------------|-----------|----------------------|--|

NOTES

| |
|--|
| Matrix-Solid. Batch Batch Rule 8330B_S_EX. |
|--|

1D
ORGANICS ANALYSIS DATA SHEET

| | | | |
|------------------|------------------------------------|--------------------|-----------------------------------|
| Report No: | <u>218081812</u> | Client Sample ID: | <u>MB1843189</u> |
| Collect Date: | <u>NA</u> Time: <u>NA</u> | GCAL Sample ID: | <u>1843189</u> |
| Matrix: | <u>Solid</u> % Moisture: <u>NA</u> | Instrument ID: | <u>HPLC3</u> |
| Sample Amt: | <u>10</u> g | Lab File ID: | <u>2180830VA04</u> |
| Injection Vol.: | <u>1.0</u> (µL) | GC Column: | <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: | <u>40000</u> (µL) | Dilution Factor: | <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: | <u>08/25/18</u> | Analysis Date: | <u>08/30/18</u> Time: <u>1145</u> |
| Prep Batch: | <u>642698</u> | Analytical Batch: | <u>643050</u> |
| Prep Method: | <u>8330B</u> | Analytical Method: | <u>EPA 8330B</u> |

CONCENTRATION UNITS: ug/kg

| CAS | ANALYTE | RESULT | Q | DL | LOD | LOQ |
|------------|------------------------------|--------|---|------|-----|-----|
| 99-35-4 | 1,3,5-Trinitrobenzene | 100 | U | 42.0 | 100 | 200 |
| 99-65-0 | 1,3-Dinitrobenzene | 100 | U | 77.0 | 100 | 200 |
| 118-96-7 | 2,4,6-Trinitrotoluene | 100 | U | 51.0 | 100 | 200 |
| 121-14-2 | 2,4-Dinitrotoluene | 100 | U | 99.0 | 100 | 200 |
| 606-20-2 | 2,6-Dinitrotoluene | 100 | U | 61.0 | 100 | 200 |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | 100 | U | 98.0 | 100 | 200 |
| 88-72-2 | 2-Nitrotoluene | 100 | U | 64.0 | 100 | 200 |
| 618-87-1 | 3,5-Dinitroaniline | 100 | U | 83.0 | 100 | 200 |
| 99-08-1 | 3-Nitrotoluene | 150 | U | 125 | 150 | 200 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | 100 | U | 77.0 | 100 | 200 |
| 99-99-0 | 4-Nitrotoluene | 100 | U | 77.0 | 100 | 200 |
| 2691-41-0 | HMX | 100 | U | 26.0 | 100 | 200 |
| 98-95-3 | Nitrobenzene | 100 | U | 36.0 | 100 | 200 |
| 55-63-0 | Nitroglycerin | 100 | U | 74.0 | 100 | 200 |
| 78-11-5 | Pentaerythritol Tetranitrate | 150 | U | 122 | 150 | 200 |
| 121-82-4 | RDX | 100 | U | 18.0 | 100 | 200 |
| 479-45-8 | Tetryl | 100 | U | 41.0 | 100 | 200 |

FORM I ORG-1

1D
ORGANICS ANALYSIS DATA SHEET

| | | | |
|------------------|------------------------------------|--------------------|--|
| Report No: | <u>218081812</u> | Client Sample ID: | <u>GRBLK for HBN 642698 [EXTO/549]</u> |
| Collect Date: | <u>NA</u> Time: <u>NA</u> | GCAL Sample ID: | <u>1843194</u> |
| Matrix: | <u>Solid</u> % Moisture: <u>NA</u> | Instrument ID: | <u>HPLC3</u> |
| Sample Amt: | <u>10</u> g | Lab File ID: | <u>2180830VA10</u> |
| Injection Vol.: | <u>1.0</u> (µL) | GC Column: | <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: | <u>40000</u> (µL) | Dilution Factor: | <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: | <u>08/25/18</u> | Analysis Date: | <u>08/30/18</u> Time: <u>1344</u> |
| Prep Batch: | <u>642698</u> | Analytical Batch: | <u>643050</u> |
| Prep Method: | <u>8330B</u> | Analytical Method: | <u>EPA 8330B</u> |

CONCENTRATION UNITS: ug/kg

| CAS | ANALYTE | RESULT | Q | DL | LOD | LOQ |
|------------|------------------------------|--------|---|------|-----|-----|
| 99-35-4 | 1,3,5-Trinitrobenzene | 100 | U | 42.0 | 100 | 200 |
| 99-65-0 | 1,3-Dinitrobenzene | 100 | U | 77.0 | 100 | 200 |
| 118-96-7 | 2,4,6-Trinitrotoluene | 100 | U | 51.0 | 100 | 200 |
| 121-14-2 | 2,4-Dinitrotoluene | 100 | U | 99.0 | 100 | 200 |
| 606-20-2 | 2,6-Dinitrotoluene | 100 | U | 61.0 | 100 | 200 |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | 100 | U | 98.0 | 100 | 200 |
| 88-72-2 | 2-Nitrotoluene | 100 | U | 64.0 | 100 | 200 |
| 618-87-1 | 3,5-Dinitroaniline | 100 | U | 83.0 | 100 | 200 |
| 99-08-1 | 3-Nitrotoluene | 150 | U | 125 | 150 | 200 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | 100 | U | 77.0 | 100 | 200 |
| 99-99-0 | 4-Nitrotoluene | 100 | U | 77.0 | 100 | 200 |
| 2691-41-0 | HMX | 100 | U | 26.0 | 100 | 200 |
| 98-95-3 | Nitrobenzene | 100 | U | 36.0 | 100 | 200 |
| 55-63-0 | Nitroglycerin | 100 | U | 74.0 | 100 | 200 |
| 78-11-5 | Pentaerythritol Tetranitrate | 150 | U | 122 | 150 | 200 |
| 121-82-4 | RDX | 100 | U | 18.0 | 100 | 200 |
| 479-45-8 | Tetryl | 100 | U | 41.0 | 100 | 200 |

FORM I ORG-1

3F
SOIL ORGANICS LCS/LCSD RECOVERY

Report No: 218081812
 Prep Method: 8330B
 Analytical Method: EPA 8330B

Prep Batch: 642698
 Analytical Batch: 643050

GCAL QC ID: 1843190

| ANALYTE | UNITS | SPIKE ADDED | SAMPLE RESULT | LCS RESULT | LCS % REC | # | QC LIMITS |
|------------------------------|-------|----------------|------------------|---------------|--------------|---|-----------|
| 1,3,5-Trinitrobenzene | ug/kg | 2050 | 0 | 971 | 47 | * | 80 - 116 |
| 1,3-Dinitrobenzene | ug/kg | 1520 | 0 | 755 | 50 | * | 73 - 119 |
| 2,4,6-Trinitrotoluene | ug/kg | 1400 | 0 | 584 | 42 | * | 71 - 120 |
| 2,4-Dinitrotoluene | ug/kg | 1610 | 0 | 424 | 26 | * | 75 - 121 |
| 2,6-Dinitrotoluene | ug/kg | 1200 | 0 | 1250 | 105 | | 79 - 117 |
| 2-Amino-4,6-dinitrotoluene | ug/kg | 1260 | 0 | 469 | 37 | * | 71 - 123 |
| 2-Nitrotoluene | ug/kg | 4080 | 0 | 1810 | 44 | * | 84 - 120 |
| 3,5-Dinitroaniline | ug/kg | 3910 | 0 | 1530 | 39 | * | 86 - 118 |
| 3-Nitrotoluene | ug/kg | 1750 | 0 | 739 | 42 | * | 67 - 129 |
| 4-Amino-2,6-dinitrotoluene | ug/kg | 2600 | 0 | 987 | 38 | * | 64 - 127 |
| 4-Nitrotoluene | ug/kg | 652 | 0 | 305 | 47 | * | 71 - 124 |
| HMX | ug/kg | 3500 | 0 | 1350 | 39 | * | 74 - 124 |
| Nitrobenzene | ug/kg | 3390 | 0 | 1320 | 39 | * | 80 - 128 |
| Nitroglycerin | ug/kg | 4600 | 0 | 1710 | 37 | * | 73 - 124 |
| Pentaerythritol Tetranitrate | ug/kg | 3100 | 0 | 957 | 31 | * | 72 - 128 |
| RDX | ug/kg | 1200 | 0 | 506 | 42 | * | 67 - 129 |
| Tetryl | ug/kg | 100 | 0 | 0 | 0 | * | 68 - 135 |

RPD : 3 out of 17 outside limits

Column to be used to flag recovery and RPD values with an asterisk

Spike Recovery: 33 out of 34 outside limits

* Values outside of QC limits

FORM III ORG-1

3F
SOIL ORGANICS LCS/LCSD RECOVERY

Report No: 218081812
 Prep Method: 8330B
 Analytical Method: EPA 8330B

Prep Batch: 642698
 Analytical Batch: 643050

GCAL QC ID: 1843191

| ANALYTE | UNITS | SPIKE ADDED | LCSD RESULT | LCSD % REC | # | % RPD | # | QC LIMITS | |
|------------------------------|-------|----------------|----------------|---------------|---|----------|---|-----------|--------|
| | | | | | | | | REC | RPD |
| 1,3,5-Trinitrobenzene | ug/kg | 2050 | 842 | 41 | * | 14 | | 80 - 116 | 0 - 20 |
| 1,3-Dinitrobenzene | ug/kg | 1520 | 657 | 43 | * | 14 | | 73 - 119 | 0 - 20 |
| 2,4,6-Trinitrotoluene | ug/kg | 1400 | 542 | 39 | * | 8 | | 71 - 120 | 0 - 20 |
| 2,4-Dinitrotoluene | ug/kg | 1610 | 519 | 32 | * | 20 | | 75 - 121 | 0 - 20 |
| 2,6-Dinitrotoluene | ug/kg | 1200 | 673 | 56 | * | 60 | * | 79 - 117 | 0 - 20 |
| 2-Amino-4,6-dinitrotoluene | ug/kg | 1260 | 430 | 34 | * | 9 | | 71 - 123 | 0 - 20 |
| 2-Nitrotoluene | ug/kg | 4080 | 1450 | 35 | * | 22 | * | 84 - 120 | 0 - 20 |
| 3,5-Dinitroaniline | ug/kg | 3910 | 1340 | 34 | * | 14 | | 86 - 118 | 0 - 20 |
| 3-Nitrotoluene | ug/kg | 1750 | 607 | 35 | * | 20 | | 67 - 129 | 0 - 20 |
| 4-Amino-2,6-dinitrotoluene | ug/kg | 2600 | 886 | 34 | * | 11 | | 64 - 127 | 0 - 20 |
| 4-Nitrotoluene | ug/kg | 652 | 392 | 60 | * | 25 | * | 71 - 124 | 0 - 20 |
| HMX | ug/kg | 3500 | 1190 | 34 | * | 13 | | 74 - 124 | 0 - 20 |
| Nitrobenzene | ug/kg | 3390 | 1140 | 34 | * | 15 | | 80 - 128 | 0 - 20 |
| Nitroglycerin | ug/kg | 4600 | 1460 | 32 | * | 16 | | 73 - 124 | 0 - 20 |
| Pentaerythritol Tetranitrate | ug/kg | 3100 | 915 | 30 | * | 4 | | 72 - 128 | 0 - 20 |
| RDX | ug/kg | 1200 | 499 | 42 | * | 1 | | 67 - 129 | 0 - 20 |
| Tetryl | ug/kg | 100 | 0 | 0 | * | 0 | | 68 - 135 | 0 - 20 |

RPD : 3 out of 17 outside limits

Column to be used to flag recovery and RPD values with an asterisk

Spike Recovery: 33 out of 34 outside limits

* Values outside of QC limits

FORM III ORG-1

3F
SOIL ORGANICS MS/MSD RECOVERY

Report No: 218081812
 Prep Method: 8330B
 Analytical Method: EPA 8330B

Parent Sample ID: WIL02DA02A
 Prep Batch: 642698
 Analytical Batch: 643776

GCAL QC ID: 21808181207

| ANALYTE | UNITS | SPIKE ADDED | SAMPLE RESULT | MS RESULT | MS % REC | # | QC LIMITS |
|------------------------------|-------|----------------|------------------|--------------|-------------|---|-----------|
| 1,3,5-Trinitrobenzene | ug/kg | 1000 | 0 | 1090 | 109 | | 80 - 116 |
| 1,3-Dinitrobenzene | ug/kg | 1000 | 0 | 1090 | 109 | | 73 - 119 |
| 2,4,6-Trinitrotoluene | ug/kg | 1000 | 0 | 1060 | 106 | | 71 - 120 |
| 2,4-Dinitrotoluene | ug/kg | 1000 | 0 | 1060 | 106 | | 75 - 121 |
| 2,6-Dinitrotoluene | ug/kg | 1000 | 0 | 1100 | 110 | | 79 - 117 |
| 2-Amino-4,6-dinitrotoluene | ug/kg | 1000 | 0 | 997 | 100 | | 71 - 123 |
| 2-Nitrotoluene | ug/kg | 1000 | 0 | 1130 | 113 | | 84 - 120 |
| 3,5-Dinitroaniline | ug/kg | 1000 | 0 | 1040 | 104 | | 86 - 118 |
| 3-Nitrotoluene | ug/kg | 1000 | 0 | 1040 | 104 | | 67 - 129 |
| 4-Amino-2,6-dinitrotoluene | ug/kg | 1000 | 0 | 1050 | 105 | | 64 - 127 |
| 4-Nitrotoluene | ug/kg | 1000 | 0 | 1120 | 112 | | 71 - 124 |
| HMX | ug/kg | 1000 | 0 | 986 | 99 | | 74 - 124 |
| Nitrobenzene | ug/kg | 1000 | 0 | 1010 | 101 | | 80 - 128 |
| Nitroglycerin | ug/kg | 1000 | 0 | 977 | 98 | | 73 - 124 |
| Pentaerythritol Tetranitrate | ug/kg | 1000 | 0 | 987 | 99 | | 72 - 128 |
| RDX | ug/kg | 1000 | 0 | 917 | 92 | | 67 - 129 |
| Tetryl | ug/kg | 1000 | 0 | 1030 | 103 | | 68 - 135 |

RPD : 3 out of 17 outside limits

Column to be used to flag recovery and RPD values with an asterisk

Spike Recovery: 0 out of 34 outside limits

* Values outside of QC limits

FORM III ORG-1

3F
SOIL ORGANICS MS/MSD RECOVERY

Report No: 218081812
 Prep Method: 8330B
 Analytical Method: EPA 8330B

Parent Sample ID: WIL02DA02A
 Prep Batch: 642698
 Analytical Batch: 643776

GCAL QC ID: 21808181208

| ANALYTE | UNITS | SPIKE ADDED | MSD RESULT | MSD % REC | # | % RPD | # | QC LIMITS | |
|------------------------------|-------|----------------|---------------|--------------|---|----------|---|-----------|--------|
| | | | | | | | | REC | RPD |
| 1,3,5-Trinitrobenzene | ug/kg | 952 | 1080 | 114 | | .6 | | 80 - 116 | 0 - 20 |
| 1,3-Dinitrobenzene | ug/kg | 952 | 1070 | 112 | | 2 | | 73 - 119 | 0 - 20 |
| 2,4,6-Trinitrotoluene | ug/kg | 952 | 957 | 100 | | 10 | | 71 - 120 | 0 - 20 |
| 2,4-Dinitrotoluene | ug/kg | 952 | 1040 | 109 | | 2 | | 75 - 121 | 0 - 20 |
| 2,6-Dinitrotoluene | ug/kg | 952 | 989 | 104 | | 10 | | 79 - 117 | 0 - 20 |
| 2-Amino-4,6-dinitrotoluene | ug/kg | 952 | 885 | 93 | | 12 | | 71 - 123 | 0 - 20 |
| 2-Nitrotoluene | ug/kg | 952 | 812 | 85 | | 32 | * | 84 - 120 | 0 - 20 |
| 3,5-Dinitroaniline | ug/kg | 952 | 1100 | 116 | | 5 | | 86 - 118 | 0 - 20 |
| 3-Nitrotoluene | ug/kg | 952 | 901 | 95 | | 14 | | 67 - 129 | 0 - 20 |
| 4-Amino-2,6-dinitrotoluene | ug/kg | 952 | 1040 | 109 | | .8 | | 64 - 127 | 0 - 20 |
| 4-Nitrotoluene | ug/kg | 952 | 846 | 89 | | 28 | * | 71 - 124 | 0 - 20 |
| HMX | ug/kg | 952 | 955 | 100 | | 3 | | 74 - 124 | 0 - 20 |
| Nitrobenzene | ug/kg | 952 | 1010 | 106 | | .5 | | 80 - 128 | 0 - 20 |
| Nitroglycerin | ug/kg | 952 | 956 | 100 | | 2 | | 73 - 124 | 0 - 20 |
| Pentaerythritol Tetranitrate | ug/kg | 952 | 1040 | 109 | | 5 | | 72 - 128 | 0 - 20 |
| RDX | ug/kg | 952 | 1010 | 106 | | 10 | | 67 - 129 | 0 - 20 |
| Tetryl | ug/kg | 952 | 820 | 86 | | 23 | * | 68 - 135 | 0 - 20 |

RPD : 3 out of 17 outside limits

Column to be used to flag recovery and RPD values with an asterisk

Spike Recovery: 0 out of 34 outside limits

* Values outside of QC limits

FORM III ORG-1

4C
ORGANIC METHOD BLANK SUMMARY

| | | | |
|------------------|-------------------|--------------------|-----------------------------------|
| Report No: | <u>218081812</u> | Method Blank ID: | <u>1846370</u> |
| Matrix: | <u>Solid</u> | Instrument ID: | <u>HPLC3</u> |
| Sample Amt: | <u>10</u> g | Lab File ID: | <u>2180912\A69</u> |
| Injection Vol.: | <u>1.0</u> (µL) | GC Column: | <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: | <u>40000</u> (µL) | Dilution Factor: | <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: | <u>09/05/18</u> | Analysis Date: | <u>09/13/18</u> Time: <u>0719</u> |
| Prep Batch: | <u>643342</u> | Analytical Batch: | <u>643776</u> |
| Prep Method: | <u>8330B</u> | Analytical Method: | <u>EPA 8330B</u> |

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD

| | <i>GCAL</i> | <i>DATE</i> | <i>TIME</i> | <i>INSTRUMENT</i> |
|-------------------------------------|------------------|-----------------|-----------------|-------------------|
| <i>CLIENT SAMPLE ID</i> | <i>SAMPLE ID</i> | <i>ANALYZED</i> | <i>ANALYZED</i> | <i>ID</i> |
| 1. LCS1846371 | 1846371 | 09/13/18 | 0738 | HPLC3 |
| 2. LCSD1846372 | 1846372 | 09/13/18 | 0758 | HPLC3 |
| 3. WIL02DA02A (RE) | 21808181213 | 09/13/18 | 0818 | HPLC3 |
| 4. WIL02DA02A MS (RE) | 21808181214 | 09/13/18 | 0838 | HPLC3 |
| 5. WIL02DA02A MSD (RE) | 21808181215 | 09/13/18 | 0857 | HPLC3 |
| 6. WIL02DA01A (RE) | 21808181216 | 09/13/18 | 0917 | HPLC3 |
| 7. WIL02DA01B (RE) | 21808181217 | 09/13/18 | 0937 | HPLC3 |
| 8. WIL02DA01B (RE)DUP | 1846373 | 09/13/18 | 0957 | HPLC3 |
| 9. WIL02DA01B (RE)DUP | 1846374 | 09/13/18 | 1016 | HPLC3 |
| 10. GRBLK for HBN 643342 [EXTO/550] | 1846375 | 09/13/18 | 1036 | HPLC3 |

FORM IV ORGANIC



8330B Explosives Prep Solid



| | | | |
|----------------------------|---------------------------------------|-------------------------------------|------------------------|
| ANALYST/TECH <i>JWA</i> | START DATE/TIME <i>9/5/18 1851</i> | END DATE/TIME <i>9/6/18 1510</i> | BATCH 643342 |
|----------------------------|---------------------------------------|-------------------------------------|------------------------|

| # | CLIENT | TYPE | CLIENT ID | GCAL ID | INITIAL WGT (g) | FINAL VOL (ml) | COMMENT | STANDARDS/ REAGENTS |
|----|--------|-------|----------------------------|-------------|-----------------|----------------|---------|---|
| 1 | QC | MB | MB 1846370 | 1846370 | <i>10.0</i> | <i>40</i> | | 8330 Surrogate 4/10ug/ml / Volume 1.0ml |
| 2 | QC | LCS | SRM 1846371 ** Use SRM ** | 1846371 | <i>10.0</i> | <i>40</i> | | <i>703-27-9</i> |
| 3 | QC | LCSD | SRMD 1846372 ** Use SRM ** | 1846372 | <i>10.0</i> | <i>40</i> | | 8330 Spike 4/10ug/ml / Volume 1.0ml |
| 4 | 4838 | SAMP | WIL02DB02A (RE) | 21808181211 | <i>10.1</i> | <i>40</i> | | <i>703-23-1</i> |
| 5 | 4838 | SAMP | WIL02DB01A (RE) | 21808181212 | <i>10.5</i> | <i>40</i> | | Solid Reference Material |
| 6 | 4838 | SAMP | WIL02DA02A (RE) | 21808181213 | <i>10.1</i> | <i>40</i> | | <i>2181150</i> |
| 7 | 4838 | MS | WIL02DA02A MS (RE) | 21808181214 | <i>10.3</i> | <i>40</i> | | HPLC Water |
| 8 | 4838 | MSD | WIL02DA02A MSD (RE) | 21808181215 | <i>10.5</i> | <i>40</i> | | <i>2181989</i> |
| 9 | 4838 | SAMP | WIL02DA01A (RE) | 21808181216 | <i>10.1</i> | <i>40</i> | | Acetonitrile |
| 10 | 4838 | SAMP | WIL02DA01B (RE) | 21808181217 | <i>10.1</i> | <i>40</i> | | <i>2124002</i> |
| 11 | QC | DUP | DUP 1846374 | 1846374 | <i>10.0</i> | <i>40</i> | | Sand |
| 12 | QC | DUP | DUP 1846373 | 1846373 | <i>10.0</i> | <i>40</i> | | |
| 13 | QC | GRBLK | GRBLK 1846375 | 1846375 | <i>10.0</i> | <i>40</i> | | |
| 14 | | | <i>LCS 1846376</i> | | <i>10.0</i> | <i>40</i> | | |
| 15 | | | <i>LCSD 1846377</i> | | <i>10.0</i> | <i>40</i> | | |
| 16 | | | | | | | | |
| 17 | | | | | | | | |
| 18 | | | | | | | | |
| 19 | | | | | | | | |
| 20 | | | | | | | | |
| 21 | | | | | | | | |
| 22 | | | | | | | | |
| 23 | | | | | | | | |
| 24 | | | | | | | | |
| 25 | | | | | | | | |
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| 27 | | | | | | | | |
| 28 | | | | | | | | |
| 29 | | | | | | | | |
| 30 | | | | | | | | |

EQUIPMENT/CONDITIONS

| | |
|------------|----------------------|
| BALANCE ID | GRINDER ID - PUCK 01 |
|------------|----------------------|

NOTES

Matrix-Solid. Batch Batch Rule 8330B_S_EX.

1D
ORGANICS ANALYSIS DATA SHEET

| | | | |
|------------------|------------------------------------|--------------------|-----------------------------------|
| Report No: | <u>218081812</u> | Client Sample ID: | <u>MB1846370</u> |
| Collect Date: | <u>NA</u> Time: <u>NA</u> | GCAL Sample ID: | <u>1846370</u> |
| Matrix: | <u>Solid</u> % Moisture: <u>NA</u> | Instrument ID: | <u>HPLC3</u> |
| Sample Amt: | <u>10</u> g | Lab File ID: | <u>2180912VA69</u> |
| Injection Vol.: | <u>1.0</u> (µL) | GC Column: | <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: | <u>40000</u> (µL) | Dilution Factor: | <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: | <u>09/05/18</u> | Analysis Date: | <u>09/13/18</u> Time: <u>0719</u> |
| Prep Batch: | <u>643342</u> | Analytical Batch: | <u>643776</u> |
| Prep Method: | <u>8330B</u> | Analytical Method: | <u>EPA 8330B</u> |

CONCENTRATION UNITS: ug/kg

| CAS | ANALYTE | RESULT | Q | DL | LOD | LOQ |
|------------|------------------------------|--------|---|------|-----|-----|
| 99-35-4 | 1,3,5-Trinitrobenzene | 100 | U | 42.0 | 100 | 200 |
| 99-65-0 | 1,3-Dinitrobenzene | 100 | U | 77.0 | 100 | 200 |
| 118-96-7 | 2,4,6-Trinitrotoluene | 100 | U | 51.0 | 100 | 200 |
| 121-14-2 | 2,4-Dinitrotoluene | 100 | U | 99.0 | 100 | 200 |
| 606-20-2 | 2,6-Dinitrotoluene | 100 | U | 61.0 | 100 | 200 |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | 100 | U | 98.0 | 100 | 200 |
| 88-72-2 | 2-Nitrotoluene | 100 | U | 64.0 | 100 | 200 |
| 618-87-1 | 3,5-Dinitroaniline | 100 | U | 83.0 | 100 | 200 |
| 99-08-1 | 3-Nitrotoluene | 150 | U | 125 | 150 | 200 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | 100 | U | 77.0 | 100 | 200 |
| 99-99-0 | 4-Nitrotoluene | 100 | U | 77.0 | 100 | 200 |
| 2691-41-0 | HMX | 100 | U | 26.0 | 100 | 200 |
| 98-95-3 | Nitrobenzene | 100 | U | 36.0 | 100 | 200 |
| 55-63-0 | Nitroglycerin | 100 | U | 74.0 | 100 | 200 |
| 78-11-5 | Pentaerythritol Tetranitrate | 150 | U | 122 | 150 | 200 |
| 121-82-4 | RDX | 100 | U | 18.0 | 100 | 200 |
| 479-45-8 | Tetryl | 100 | U | 41.0 | 100 | 200 |

FORM I ORG-1

1D
ORGANICS ANALYSIS DATA SHEET

| | | | |
|------------------|------------------------------------|--------------------|--|
| Report No: | <u>218081812</u> | Client Sample ID: | <u>GRBLK for HBN 643342 [EXTO/550]</u> |
| Collect Date: | <u>NA</u> Time: <u>NA</u> | GCAL Sample ID: | <u>1846375</u> |
| Matrix: | <u>Solid</u> % Moisture: <u>NA</u> | Instrument ID: | <u>HPLC3</u> |
| Sample Amt: | <u>10</u> g | Lab File ID: | <u>2180912VA79</u> |
| Injection Vol.: | <u>1.0</u> (µL) | GC Column: | <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: | <u>40000</u> (µL) | Dilution Factor: | <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: | <u>09/05/18</u> | Analysis Date: | <u>09/13/18</u> Time: <u>1036</u> |
| Prep Batch: | <u>643342</u> | Analytical Batch: | <u>643776</u> |
| Prep Method: | <u>8330B</u> | Analytical Method: | <u>EPA 8330B</u> |

CONCENTRATION UNITS: ug/kg

| CAS | ANALYTE | RESULT | Q | DL | LOD | LOQ |
|------------|------------------------------|--------|---|------|-----|-----|
| 99-35-4 | 1,3,5-Trinitrobenzene | 100 | U | 42.0 | 100 | 200 |
| 99-65-0 | 1,3-Dinitrobenzene | 100 | U | 77.0 | 100 | 200 |
| 118-96-7 | 2,4,6-Trinitrotoluene | 100 | U | 51.0 | 100 | 200 |
| 121-14-2 | 2,4-Dinitrotoluene | 100 | U | 99.0 | 100 | 200 |
| 606-20-2 | 2,6-Dinitrotoluene | 100 | U | 61.0 | 100 | 200 |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | 100 | U | 98.0 | 100 | 200 |
| 88-72-2 | 2-Nitrotoluene | 100 | U | 64.0 | 100 | 200 |
| 618-87-1 | 3,5-Dinitroaniline | 100 | U | 83.0 | 100 | 200 |
| 99-08-1 | 3-Nitrotoluene | 150 | U | 125 | 150 | 200 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | 100 | U | 77.0 | 100 | 200 |
| 99-99-0 | 4-Nitrotoluene | 100 | U | 77.0 | 100 | 200 |
| 2691-41-0 | HMX | 100 | U | 26.0 | 100 | 200 |
| 98-95-3 | Nitrobenzene | 100 | U | 36.0 | 100 | 200 |
| 55-63-0 | Nitroglycerin | 100 | U | 74.0 | 100 | 200 |
| 78-11-5 | Pentaerythritol Tetranitrate | 150 | U | 122 | 150 | 200 |
| 121-82-4 | RDX | 100 | U | 18.0 | 100 | 200 |
| 479-45-8 | Tetryl | 100 | U | 41.0 | 100 | 200 |

FORM I ORG-1

3F
SOIL ORGANICS LCS/LCSD RECOVERY

Report No: 218081812
 Prep Method: 8330B
 Analytical Method: EPA 8330B

Prep Batch: 643342
 Analytical Batch: 643776

GCAL QC ID: 1846371

| ANALYTE | UNITS | SPIKE ADDED | SAMPLE RESULT | LCS RESULT | LCS % REC | # | QC LIMITS |
|------------------------------|-------|----------------|------------------|---------------|--------------|---|-----------|
| 1,3,5-Trinitrobenzene | ug/kg | 1750 | 0 | 2110 | 121 | * | 80 - 116 |
| 1,3-Dinitrobenzene | ug/kg | 1280 | 0 | 1620 | 126 | * | 73 - 119 |
| 2,4,6-Trinitrotoluene | ug/kg | 1120 | 0 | 1240 | 111 | | 71 - 120 |
| 2,4-Dinitrotoluene | ug/kg | 1210 | 0 | 1300 | 107 | | 75 - 121 |
| 2,6-Dinitrotoluene | ug/kg | 1310 | 0 | 1370 | 104 | | 79 - 117 |
| 2-Amino-4,6-dinitrotoluene | ug/kg | 1210 | 0 | 1250 | 104 | | 71 - 123 |
| 2-Nitrotoluene | ug/kg | 3160 | 0 | 3540 | 112 | | 84 - 120 |
| 3,5-Dinitroaniline | ug/kg | 2690 | 0 | 3020 | 112 | | 86 - 118 |
| 3-Nitrotoluene | ug/kg | 1360 | 0 | 1540 | 113 | | 67 - 129 |
| 4-Amino-2,6-dinitrotoluene | ug/kg | 1390 | 0 | 1360 | 98 | | 64 - 127 |
| 4-Nitrotoluene | ug/kg | 688 | 0 | 741 | 108 | | 71 - 124 |
| HMX | ug/kg | 2850 | 0 | 3290 | 116 | | 74 - 124 |
| Nitrobenzene | ug/kg | 2340 | 0 | 2510 | 107 | | 80 - 128 |
| Nitroglycerin | ug/kg | 3870 | 0 | 4540 | 117 | | 73 - 124 |
| Pentaerythritol Tetranitrate | ug/kg | 2650 | 0 | 3070 | 116 | | 72 - 128 |
| RDX | ug/kg | 923 | 0 | 1030 | 111 | | 67 - 129 |
| Tetryl | ug/kg | 100 | 0 | 0 | 0 | * | 68 - 135 |

RPD: 3 out of 17 outside limits

Column to be used to flag recovery and RPD values with an asterisk

Spike Recovery: 8 out of 34 outside limits

* Values outside of QC limits

FORM III ORG-1

3F
SOIL ORGANICS LCS/LCSD RECOVERY

Report No: 218081812
 Prep Method: 8330B
 Analytical Method: EPA 8330B

Prep Batch: 643342
 Analytical Batch: 643776

GCAL QC ID: 1846372

| ANALYTE | UNITS | SPIKE ADDED | LCSD RESULT | LCSD % REC | # | % RPD | # | QC LIMITS | |
|------------------------------|-------|----------------|----------------|---------------|---|----------|---|-----------|--------|
| | | | | | | | | REC | RPD |
| 1,3,5-Trinitrobenzene | ug/kg | 1750 | 2070 | 119 | * | 2 | | 80 - 116 | 0 - 20 |
| 1,3-Dinitrobenzene | ug/kg | 1280 | 1420 | 111 | | 13 | | 73 - 119 | 0 - 20 |
| 2,4,6-Trinitrotoluene | ug/kg | 1120 | 1320 | 118 | | 6 | | 71 - 120 | 0 - 20 |
| 2,4-Dinitrotoluene | ug/kg | 1210 | 1290 | 106 | | 1 | | 75 - 121 | 0 - 20 |
| 2,6-Dinitrotoluene | ug/kg | 1310 | 1300 | 99 | | 5 | | 79 - 117 | 0 - 20 |
| 2-Amino-4,6-dinitrotoluene | ug/kg | 1210 | 1480 | 122 | | 17 | | 71 - 123 | 0 - 20 |
| 2-Nitrotoluene | ug/kg | 3160 | 3380 | 107 | | 5 | | 84 - 120 | 0 - 20 |
| 3,5-Dinitroaniline | ug/kg | 2690 | 2880 | 107 | | 4 | | 86 - 118 | 0 - 20 |
| 3-Nitrotoluene | ug/kg | 1360 | 1970 | 145 | * | 24 | * | 67 - 129 | 0 - 20 |
| 4-Amino-2,6-dinitrotoluene | ug/kg | 1390 | 1330 | 96 | | 2 | | 64 - 127 | 0 - 20 |
| 4-Nitrotoluene | ug/kg | 688 | 2430 | 354 | * | 107 | * | 71 - 124 | 0 - 20 |
| HMX | ug/kg | 2850 | 3370 | 118 | | 2 | | 74 - 124 | 0 - 20 |
| Nitrobenzene | ug/kg | 2340 | 2480 | 106 | | 1 | | 80 - 128 | 0 - 20 |
| Nitroglycerin | ug/kg | 3870 | 4190 | 108 | | 8 | | 73 - 124 | 0 - 20 |
| Pentaerythritol Tetranitrate | ug/kg | 2650 | 4410 | 166 | * | 36 | * | 72 - 128 | 0 - 20 |
| RDX | ug/kg | 923 | 1010 | 110 | | 2 | | 67 - 129 | 0 - 20 |
| Tetryl | ug/kg | 100 | 0 | 0 | * | 0 | | 68 - 135 | 0 - 20 |

RPD : 3 out of 17 outside limits

Column to be used to flag recovery and RPD values with an asterisk

Spike Recovery: 8 out of 34 outside limits

* Values outside of QC limits

FORM III ORG-1

3F
SOIL ORGANICS MS/MSD RECOVERY

Report No: 218081812
 Prep Method: 8330B
 Analytical Method: EPA 8330B

Parent Sample ID: WIL02DA02A (RE)
 Prep Batch: 643342
 Analytical Batch: 643776

GCAL QC ID: 21808181214

| ANALYTE | UNITS | SPIKE ADDED | SAMPLE RESULT | MS RESULT | MS % REC | # | QC LIMITS |
|------------------------------|-------|----------------|------------------|--------------|-------------|---|-----------|
| 1,3,5-Trinitrobenzene | ug/kg | 971 | 0 | 828 | 85 | | 80 - 116 |
| 1,3-Dinitrobenzene | ug/kg | 971 | 0 | 836 | 86 | | 73 - 119 |
| 2,4,6-Trinitrotoluene | ug/kg | 971 | 0 | 767 | 79 | | 71 - 120 |
| 2,4-Dinitrotoluene | ug/kg | 971 | 0 | 884 | 91 | | 75 - 121 |
| 2,6-Dinitrotoluene | ug/kg | 971 | 0 | 803 | 83 | | 79 - 117 |
| 2-Amino-4,6-dinitrotoluene | ug/kg | 971 | 0 | 889 | 92 | | 71 - 123 |
| 2-Nitrotoluene | ug/kg | 971 | 0 | 864 | 89 | | 84 - 120 |
| 3,5-Dinitroaniline | ug/kg | 971 | 0 | 923 | 95 | | 86 - 118 |
| 3-Nitrotoluene | ug/kg | 971 | 0 | 833 | 86 | | 67 - 129 |
| 4-Amino-2,6-dinitrotoluene | ug/kg | 971 | 0 | 1030 | 106 | | 64 - 127 |
| 4-Nitrotoluene | ug/kg | 971 | 0 | 930 | 96 | | 71 - 124 |
| HMX | ug/kg | 971 | 0 | 596 | 61 | * | 74 - 124 |
| Nitrobenzene | ug/kg | 971 | 0 | 918 | 95 | | 80 - 128 |
| Nitroglycerin | ug/kg | 971 | 0 | 848 | 87 | | 73 - 124 |
| Pentaerythritol Tetranitrate | ug/kg | 971 | 0 | 947 | 98 | | 72 - 128 |
| RDX | ug/kg | 971 | 0 | 816 | 84 | | 67 - 129 |
| Tetryl | ug/kg | 971 | 0 | 734 | 76 | | 68 - 135 |

RPD : 2 out of 17 outside limits

Column to be used to flag recovery and RPD values with an asterisk

Spike Recovery: 1 out of 34 outside limits

* Values outside of QC limits

FORM III ORG-1

3F
SOIL ORGANICS MS/MSD RECOVERY

Report No: 218081812
 Prep Method: 8330B
 Analytical Method: EPA 8330B

Parent Sample ID: WIL02DA02A (RE)
 Prep Batch: 643342
 Analytical Batch: 643776

GCAL QC ID: 21808181215

| ANALYTE | UNITS | SPIKE ADDED | MSD RESULT | MSD % REC | # | % RPD | # | QC LIMITS | |
|------------------------------|-------|----------------|---------------|--------------|---|----------|---|-----------|--------|
| | | | | | | | | REC | RPD |
| 1,3,5-Trinitrobenzene | ug/kg | 952 | 956 | 100 | | 14 | | 80 - 116 | 0 - 20 |
| 1,3-Dinitrobenzene | ug/kg | 952 | 884 | 93 | | 6 | | 73 - 119 | 0 - 20 |
| 2,4,6-Trinitrotoluene | ug/kg | 952 | 887 | 93 | | 15 | | 71 - 120 | 0 - 20 |
| 2,4-Dinitrotoluene | ug/kg | 952 | 778 | 82 | | 13 | | 75 - 121 | 0 - 20 |
| 2,6-Dinitrotoluene | ug/kg | 952 | 1090 | 114 | | 30 | * | 79 - 117 | 0 - 20 |
| 2-Amino-4,6-dinitrotoluene | ug/kg | 952 | 818 | 86 | | 8 | | 71 - 123 | 0 - 20 |
| 2-Nitrotoluene | ug/kg | 952 | 827 | 87 | | 4 | | 84 - 120 | 0 - 20 |
| 3,5-Dinitroaniline | ug/kg | 952 | 944 | 99 | | 2 | | 86 - 118 | 0 - 20 |
| 3-Nitrotoluene | ug/kg | 952 | 848 | 89 | | 2 | | 67 - 129 | 0 - 20 |
| 4-Amino-2,6-dinitrotoluene | ug/kg | 952 | 1040 | 109 | | 1 | | 64 - 127 | 0 - 20 |
| 4-Nitrotoluene | ug/kg | 952 | 928 | 97 | | .2 | | 71 - 124 | 0 - 20 |
| HMX | ug/kg | 952 | 794 | 83 | | 29 | * | 74 - 124 | 0 - 20 |
| Nitrobenzene | ug/kg | 952 | 927 | 97 | | 1 | | 80 - 128 | 0 - 20 |
| Nitroglycerin | ug/kg | 952 | 925 | 97 | | 9 | | 73 - 124 | 0 - 20 |
| Pentaerythritol Tetranitrate | ug/kg | 952 | 915 | 96 | | 3 | | 72 - 128 | 0 - 20 |
| RDX | ug/kg | 952 | 808 | 85 | | .9 | | 67 - 129 | 0 - 20 |
| Tetryl | ug/kg | 952 | 769 | 81 | | 5 | | 68 - 135 | 0 - 20 |

RPD : 2 out of 17 outside limits

Column to be used to flag recovery and RPD values with an asterisk

Spike Recovery: 1 out of 34 outside limits

* Values outside of QC limits

FORM III ORG-1

EPA 8330B REPLICATE SUMMARY

| | | | |
|--------------------|-----------------------------|-------------------|--------------------|
| Report No: | <u>218081814</u> | Parent Sample ID: | <u>WIL02IS03</u> |
| Prep Method: | <u>EPA 8330B</u> | Parent GCAL ID: | <u>21808181407</u> |
| Prep Date: | <u>8/24/2018 6:00:00 PM</u> | Prep Batch: | <u>642680</u> |
| Analytical Method: | <u>EPA 8330B</u> | | |

| ANALYTE | CAS | UNITS | PARENT RESULT | REP #1 RESULT (1843104) | REP #2 RESULT (1843105) | %RSD | # |
|------------------------------|------------|-------|------------------|-------------------------------|-------------------------------|------|---|
| 1,3,5-Trinitrobenzene | 99-35-4 | ug/Kg | 0 | 0 | 0 | 0 | |
| 1,3-Dinitrobenzene | 99-65-0 | ug/Kg | 0 | 0 | 0 | 0 | |
| 2,4,6-Trinitrotoluene | 118-96-7 | ug/Kg | 0 | 0 | 0 | 0 | |
| 2,4-Dinitrotoluene | 121-14-2 | ug/Kg | 0 | 0 | 0 | 0 | |
| 2,6-Dinitrotoluene | 606-20-2 | ug/Kg | 0 | 0 | 0 | 0 | |
| 2-Amino-4,6-dinitrotoluene | 35572-78-2 | ug/Kg | 0 | 0 | 0 | 0 | |
| 2-Nitrotoluene | 88-72-2 | ug/Kg | 0 | 0 | 0 | 0 | |
| 3,5-Dinitroaniline | 618-87-1 | ug/Kg | 0 | 0 | 0 | 0 | |
| 3-Nitrotoluene | 99-08-1 | ug/Kg | 0 | 0 | 0 | 0 | |
| 4-Amino-2,6-dinitrotoluene | 19406-51-0 | ug/Kg | 0 | 0 | 0 | 0 | |
| 4-Nitrotoluene | 99-99-0 | ug/Kg | 0 | 0 | 0 | 0 | |
| HMX | 2691-41-0 | ug/Kg | 0 | 0 | 0 | 0 | |
| Nitrobenzene | 98-95-3 | ug/Kg | 0 | 0 | 0 | 0 | |
| Nitroglycerin | 55-63-0 | ug/Kg | 0 | 0 | 0 | 0 | |
| Pentaerythritol Tetranitrate | 78-11-5 | ug/Kg | 0 | 0 | 0 | 0 | |
| RDX | 121-82-4 | ug/Kg | 0 | 0 | 0 | 0 | |
| Tetryl | 479-45-8 | ug/Kg | 0 | 0 | 0 | 0 | |

* - RSD greater than 20%

4C
ORGANIC METHOD BLANK SUMMARY

| | |
|------------------------------------|--|
| Report No: <u>218081814</u> | Method Blank ID: <u>1843101</u> |
| Matrix: <u>Solid</u> | Instrument ID: <u>HPLC3</u> |
| Sample Amt: <u>10</u> g | Lab File ID: <u>2180830A19</u> |
| Injection Vol.: <u>1.0</u> (µL) | GC Column: <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: <u>40000</u> (µL) | Dilution Factor: <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: <u>08/24/18</u> | Analysis Date: <u>08/30/18</u> Time: <u>1641</u> |
| Prep Batch: <u>642680</u> | Analytical Batch: <u>643050</u> |
| Prep Method: <u>8330B</u> | Analytical Method: <u>EPA 8330B</u> |

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD

| <i>CLIENT SAMPLE ID</i> | <i>GCAL SAMPLE ID</i> | <i>DATE ANALYZED</i> | <i>TIME ANALYZED</i> | <i>INSTRUMENT ID</i> |
|------------------------------------|---------------------------|--------------------------|--------------------------|--------------------------|
| 1. WIL02IS01 | 21808181401 | 08/30/18 | 1741 | HPLC3 |
| 2. WIL02IS01 MS | 21808181402 | 08/30/18 | 1800 | HPLC3 |
| 3. WIL02IS01 MSD | 21808181403 | 08/30/18 | 1820 | HPLC3 |
| 4. WIL02IS02 | 21808181404 | 08/30/18 | 1840 | HPLC3 |
| 5. WIL02IS03 | 21808181407 | 08/30/18 | 1900 | HPLC3 |
| 6. WIL02IS03DUP | 1843105 | 08/30/18 | 1919 | HPLC3 |
| 7. WIL02IS03DUP | 1843104 | 08/30/18 | 1939 | HPLC3 |
| 8. GRBLK for HBN 642680 [EXTO/549] | 1843106 | 08/30/18 | 1959 | HPLC3 |
| 9. LCS1843214 | 1843214 | 08/30/18 | 2019 | HPLC3 |
| 10. LCSD1843215 | 1843215 | 08/30/18 | 2039 | HPLC3 |

FORM IV ORGANIC



8330B Explosives Prep Solid



| | | | | | | | |
|--------------|------------|-----------------|---------------------|---------------|---------------------|-------|--------|
| ANALYST/TECH | <i>DNB</i> | START DATE/TIME | <i>8/24/18 1800</i> | END DATE/TIME | <i>8/25/18 1700</i> | BATCH | 642680 |
|--------------|------------|-----------------|---------------------|---------------|---------------------|-------|--------|

| # | CLIENT | TYPE | CLIENT ID | GCAL ID | INITIAL WGT (g) | FINAL VOL (mL) | COMMENT | STANDARDS/ REAGENTS |
|----|--------|-------|----------------------------|-------------|-----------------|----------------|--------------|---|
| 1 | QC | MB | MB 1843101 | 1843101 | <i>10.0</i> | <i>40</i> | | 8330 Surrogate 4/10ug/ml / Volume 1.0ml |
| 2 | QC | LCS | SRM 1843102 ** Use SRM ** | 1843102 | <i>10.0</i> | <i>40</i> | | <i>703-27-8</i> |
| 3 | QC | LCSD | SRMD 1843103 ** Use SRM ** | 1843103 | <i>10.0</i> | <i>40</i> | | 8330 Spike 4/10ug/ml / Volume 1.0ml |
| 4 | 4838 | SAMP | WIL02IS01 | 21808181401 | <i>10.1</i> | <i>40</i> | | <i>703-23-1</i> |
| 5 | 4838 | MS | WIL02IS01 MS | 21808181402 | <i>10.1</i> | <i>40</i> | | Solid Reference Material |
| 6 | 4838 | MSD | WIL02IS01 MSD | 21808181403 | <i>10.0</i> | <i>40</i> | | <i>2127150</i> |
| 7 | 4838 | SAMP | WIL02IS02 | 21808181404 | <i>10.5</i> | <i>40</i> | | HPLC Water |
| 8 | 4838 | SAMP | WIL02IS03 | 21808181407 | <i>10.0</i> | <i>40</i> | | <i>2126989</i> |
| 9 | QC | DUP | DUP 1843105 | 1843105 | <i>10.0</i> | <i>40</i> | | Acetonitrile |
| 10 | QC | DUP | DUP 1843104 | 1843104 | <i>10.0</i> | <i>40</i> | | <i>2127054</i> |
| 11 | QC | GRBLK | GRBLK 1843106 | 1843106 | <i>10.0</i> | <i>40</i> | | Sand |
| 12 | QC | LCS | SRM 1843214 ** Use SRM ** | 1843214 | <i>10.0</i> | <i>40</i> | <i>spike</i> | |
| 13 | QC | LCSD | SRMD 1843215 ** Use SRM ** | 1843215 | <i>10.0</i> | <i>40</i> | <i>spike</i> | |
| 14 | | | | | | | | |
| 15 | | | | | | | | |
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| 30 | | | | | | | | |

EQUIPMENT/CONDITIONS

| | | | |
|------------|-----------|----------------------|--|
| BALANCE ID | <i>03</i> | GRINDER ID - PUCK 01 | |
|------------|-----------|----------------------|--|

NOTES

Matrix-Solid. Batch:Batch Rule 8330B_S_EX.

1D
ORGANICS ANALYSIS DATA SHEET

| | | | |
|------------------|------------------------------------|--------------------|-----------------------------------|
| Report No: | <u>218081814</u> | Client Sample ID: | <u>MB1843101</u> |
| Collect Date: | <u>NA</u> Time: <u>NA</u> | GCAL Sample ID: | <u>1843101</u> |
| Matrix: | <u>Solid</u> % Moisture: <u>NA</u> | Instrument ID: | <u>HPLC3</u> |
| Sample Amt: | <u>10</u> g | Lab File ID: | <u>2180830A19</u> |
| Injection Vol.: | <u>1.0</u> (µL) | GC Column: | <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: | <u>40000</u> (µL) | Dilution Factor: | <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: | <u>08/24/18</u> | Analysis Date: | <u>08/30/18</u> Time: <u>1641</u> |
| Prep Batch: | <u>642680</u> | Analytical Batch: | <u>643050</u> |
| Prep Method: | <u>8330B</u> | Analytical Method: | <u>EPA 8330B</u> |

CONCENTRATION UNITS: ug/kg

| CAS | ANALYTE | RESULT | Q | DL | LOD | LOQ |
|------------|------------------------------|--------|---|------|-----|-----|
| 99-35-4 | 1,3,5-Trinitrobenzene | 100 | U | 42.0 | 100 | 200 |
| 99-65-0 | 1,3-Dinitrobenzene | 100 | U | 77.0 | 100 | 200 |
| 118-96-7 | 2,4,6-Trinitrotoluene | 100 | U | 51.0 | 100 | 200 |
| 121-14-2 | 2,4-Dinitrotoluene | 100 | U | 99.0 | 100 | 200 |
| 606-20-2 | 2,6-Dinitrotoluene | 100 | U | 61.0 | 100 | 200 |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | 100 | U | 98.0 | 100 | 200 |
| 88-72-2 | 2-Nitrotoluene | 100 | U | 64.0 | 100 | 200 |
| 618-87-1 | 3,5-Dinitroaniline | 100 | U | 83.0 | 100 | 200 |
| 99-08-1 | 3-Nitrotoluene | 150 | U | 125 | 150 | 200 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | 100 | U | 77.0 | 100 | 200 |
| 99-99-0 | 4-Nitrotoluene | 100 | U | 77.0 | 100 | 200 |
| 2691-41-0 | HMX | 100 | U | 26.0 | 100 | 200 |
| 98-95-3 | Nitrobenzene | 100 | U | 36.0 | 100 | 200 |
| 55-63-0 | Nitroglycerin | 100 | U | 74.0 | 100 | 200 |
| 78-11-5 | Pentaerythritol Tetranitrate | 150 | U | 122 | 150 | 200 |
| 121-82-4 | RDX | 100 | U | 18.0 | 100 | 200 |
| 479-45-8 | Tetryl | 100 | U | 41.0 | 100 | 200 |

FORM I ORG-1

1D
ORGANICS ANALYSIS DATA SHEET

| | | | |
|------------------|------------------------------------|--------------------|--|
| Report No: | <u>218081814</u> | Client Sample ID: | <u>GRBLK for HBN 642680 [EXTO/549]</u> |
| Collect Date: | <u>NA</u> Time: <u>NA</u> | GCAL Sample ID: | <u>1843106</u> |
| Matrix: | <u>Solid</u> % Moisture: <u>NA</u> | Instrument ID: | <u>HPLC3</u> |
| Sample Amt: | <u>10</u> g | Lab File ID: | <u>2180830A29</u> |
| Injection Vol.: | <u>1.0</u> (µL) | GC Column: | <u>ARC18</u> ID <u>3</u> (mm) |
| Prep Final Vol.: | <u>40000</u> (µL) | Dilution Factor: | <u>1</u> Analyst: <u>MEG</u> |
| Prep Date: | <u>08/24/18</u> | Analysis Date: | <u>08/30/18</u> Time: <u>1959</u> |
| Prep Batch: | <u>642680</u> | Analytical Batch: | <u>643050</u> |
| Prep Method: | <u>8330B</u> | Analytical Method: | <u>EPA 8330B</u> |

CONCENTRATION UNITS: ug/kg

| CAS | ANALYTE | RESULT | Q | DL | LOD | LOQ |
|------------|------------------------------|--------|---|------|-----|-----|
| 99-35-4 | 1,3,5-Trinitrobenzene | 100 | U | 42.0 | 100 | 200 |
| 99-65-0 | 1,3-Dinitrobenzene | 100 | U | 77.0 | 100 | 200 |
| 118-96-7 | 2,4,6-Trinitrotoluene | 100 | U | 51.0 | 100 | 200 |
| 121-14-2 | 2,4-Dinitrotoluene | 100 | U | 99.0 | 100 | 200 |
| 606-20-2 | 2,6-Dinitrotoluene | 100 | U | 61.0 | 100 | 200 |
| 35572-78-2 | 2-Amino-4,6-dinitrotoluene | 100 | U | 98.0 | 100 | 200 |
| 88-72-2 | 2-Nitrotoluene | 100 | U | 64.0 | 100 | 200 |
| 618-87-1 | 3,5-Dinitroaniline | 100 | U | 83.0 | 100 | 200 |
| 99-08-1 | 3-Nitrotoluene | 150 | U | 125 | 150 | 200 |
| 19406-51-0 | 4-Amino-2,6-dinitrotoluene | 100 | U | 77.0 | 100 | 200 |
| 99-99-0 | 4-Nitrotoluene | 100 | U | 77.0 | 100 | 200 |
| 2691-41-0 | HMX | 100 | U | 26.0 | 100 | 200 |
| 98-95-3 | Nitrobenzene | 100 | U | 36.0 | 100 | 200 |
| 55-63-0 | Nitroglycerin | 100 | U | 74.0 | 100 | 200 |
| 78-11-5 | Pentaerythritol Tetranitrate | 150 | U | 122 | 150 | 200 |
| 121-82-4 | RDX | 100 | U | 18.0 | 100 | 200 |
| 479-45-8 | Tetryl | 100 | U | 41.0 | 100 | 200 |

FORM I ORG-1

3F
SOIL ORGANICS LCS/LCSD RECOVERY

Report No: 218081814

Prep Method: 8330B

Analytical Method: EPA 8330B

Prep Batch: 642680

Analytical Batch: 643050

GCAL QC ID: 1843214

| ANALYTE | UNITS | SPIKE ADDED | SAMPLE RESULT | LCS RESULT | LCS % REC | # | QC LIMITS |
|------------------------------|-------|----------------|------------------|---------------|--------------|---|-----------|
| 1,3,5-Trinitrobenzene | ug/kg | 1000 | 0 | 807 | 81 | | 80 - 116 |
| 1,3-Dinitrobenzene | ug/kg | 1000 | 0 | 787 | 79 | | 73 - 119 |
| 2,4,6-Trinitrotoluene | ug/kg | 1000 | 0 | 747 | 75 | | 71 - 120 |
| 2,4-Dinitrotoluene | ug/kg | 1000 | 0 | 862 | 86 | | 75 - 121 |
| 2,6-Dinitrotoluene | ug/kg | 1000 | 0 | 815 | 81 | | 79 - 117 |
| 2-Amino-4,6-dinitrotoluene | ug/kg | 1000 | 0 | 757 | 76 | | 71 - 123 |
| 2-Nitrotoluene | ug/kg | 1000 | 0 | 913 | 91 | | 84 - 120 |
| 3,5-Dinitroaniline | ug/kg | 1000 | 0 | 875 | 87 | | 86 - 118 |
| 3-Nitrotoluene | ug/kg | 1000 | 0 | 708 | 71 | | 67 - 129 |
| 4-Amino-2,6-dinitrotoluene | ug/kg | 1000 | 0 | 739 | 74 | | 64 - 127 |
| 4-Nitrotoluene | ug/kg | 1000 | 0 | 816 | 82 | | 71 - 124 |
| HMX | ug/kg | 1000 | 0 | 833 | 83 | | 74 - 124 |
| Nitrobenzene | ug/kg | 1000 | 0 | 819 | 82 | | 80 - 128 |
| Nitroglycerin | ug/kg | 1000 | 0 | 782 | 78 | | 73 - 124 |
| Pentaerythritol Tetranitrate | ug/kg | 1000 | 0 | 836 | 84 | | 72 - 128 |
| RDX | ug/kg | 1000 | 0 | 822 | 82 | | 67 - 129 |
| Tetryl | ug/kg | 1000 | 0 | 827 | 83 | | 68 - 135 |

RPD : 1 out of 17 outside limits

Column to be used to flag recovery and RPD values with an asterisk

Spike Recovery: 0 out of 34 outside limits

* Values outside of QC limits

FORM III ORG-1

3F
SOIL ORGANICS LCS/LCSD RECOVERY

Report No: 218081814

Prep Method: 8330B

Analytical Method: EPA 8330B

Prep Batch: 642680

Analytical Batch: 643050

GCAL QC ID: 1843215

| ANALYTE | UNITS | SPIKE ADDED | LCSD RESULT | LCSD % REC | # | % RPD | # | QC LIMITS | |
|------------------------------|-------|----------------|----------------|---------------|---|----------|---|-----------|--------|
| | | | | | | | | REC | RPD |
| 1,3,5-Trinitrobenzene | ug/kg | 1000 | 910 | 91 | | 12 | | 80 - 116 | 0 - 20 |
| 1,3-Dinitrobenzene | ug/kg | 1000 | 872 | 87 | | 10 | | 73 - 119 | 0 - 20 |
| 2,4,6-Trinitrotoluene | ug/kg | 1000 | 842 | 84 | | 12 | | 71 - 120 | 0 - 20 |
| 2,4-Dinitrotoluene | ug/kg | 1000 | 882 | 88 | | 2 | | 75 - 121 | 0 - 20 |
| 2,6-Dinitrotoluene | ug/kg | 1000 | 877 | 88 | | 7 | | 79 - 117 | 0 - 20 |
| 2-Amino-4,6-dinitrotoluene | ug/kg | 1000 | 888 | 89 | | 16 | | 71 - 123 | 0 - 20 |
| 2-Nitrotoluene | ug/kg | 1000 | 959 | 96 | | 5 | | 84 - 120 | 0 - 20 |
| 3,5-Dinitroaniline | ug/kg | 1000 | 902 | 90 | | 3 | | 86 - 118 | 0 - 20 |
| 3-Nitrotoluene | ug/kg | 1000 | 845 | 84 | | 18 | | 67 - 129 | 0 - 20 |
| 4-Amino-2,6-dinitrotoluene | ug/kg | 1000 | 872 | 87 | | 16 | | 64 - 127 | 0 - 20 |
| 4-Nitrotoluene | ug/kg | 1000 | 1100 | 110 | | 30 | * | 71 - 124 | 0 - 20 |
| HMX | ug/kg | 1000 | 829 | 83 | | .4 | | 74 - 124 | 0 - 20 |
| Nitrobenzene | ug/kg | 1000 | 904 | 90 | | 10 | | 80 - 128 | 0 - 20 |
| Nitroglycerin | ug/kg | 1000 | 854 | 85 | | 9 | | 73 - 124 | 0 - 20 |
| Pentaerythritol Tetranitrate | ug/kg | 1000 | 839 | 84 | | .3 | | 72 - 128 | 0 - 20 |
| RDX | ug/kg | 1000 | 909 | 91 | | 10 | | 67 - 129 | 0 - 20 |
| Tetryl | ug/kg | 1000 | 916 | 92 | | 10 | | 68 - 135 | 0 - 20 |

RPD : 1 out of 17 outside limits

Column to be used to flag recovery and RPD values with an asterisk

Spike Recovery: 0 out of 34 outside limits

* Values outside of QC limits

FORM III ORG-1

3F
SOIL ORGANICS MS/MSD RECOVERY

Report No: 218081814
 Prep Method: 8330B
 Analytical Method: EPA 8330B

Parent Sample ID: WIL02IS01
 Prep Batch: 642680
 Analytical Batch: 643050

GCAL QC ID: 21808181402

| ANALYTE | UNITS | SPIKE ADDED | SAMPLE RESULT | MS RESULT | MS % REC | # | QC LIMITS |
|------------------------------|-------|----------------|------------------|--------------|-------------|---|-----------|
| 1,3,5-Trinitrobenzene | ug/kg | 990 | 0 | 975 | 98 | | 80 - 116 |
| 1,3-Dinitrobenzene | ug/kg | 990 | 0 | 916 | 92 | | 73 - 119 |
| 2,4,6-Trinitrotoluene | ug/kg | 990 | 0 | 908 | 92 | | 71 - 120 |
| 2,4-Dinitrotoluene | ug/kg | 990 | 0 | 863 | 87 | | 75 - 121 |
| 2,6-Dinitrotoluene | ug/kg | 990 | 0 | 991 | 100 | | 79 - 117 |
| 2-Amino-4,6-dinitrotoluene | ug/kg | 990 | 0 | 895 | 90 | | 71 - 123 |
| 2-Nitrotoluene | ug/kg | 990 | 0 | 909 | 92 | | 84 - 120 |
| 3,5-Dinitroaniline | ug/kg | 990 | 0 | 889 | 90 | | 86 - 118 |
| 3-Nitrotoluene | ug/kg | 990 | 0 | 871 | 88 | | 67 - 129 |
| 4-Amino-2,6-dinitrotoluene | ug/kg | 990 | 0 | 898 | 91 | | 64 - 127 |
| 4-Nitrotoluene | ug/kg | 990 | 0 | 976 | 99 | | 71 - 124 |
| HMX | ug/kg | 990 | 0 | 834 | 84 | | 74 - 124 |
| Nitrobenzene | ug/kg | 990 | 0 | 1030 | 104 | | 80 - 128 |
| Nitroglycerin | ug/kg | 990 | 0 | 904 | 91 | | 73 - 124 |
| Pentaerythritol Tetranitrate | ug/kg | 990 | 0 | 986 | 100 | | 72 - 128 |
| RDX | ug/kg | 990 | 0 | 891 | 90 | | 67 - 129 |
| Tetryl | ug/kg | 990 | 0 | 720 | 73 | | 68 - 135 |

RPD : 0 out of 17 outside limits

Column to be used to flag recovery and RPD values with an asterisk

Spike Recovery: 0 out of 34 outside limits

* Values outside of QC limits

FORM III ORG-1

3F
SOIL ORGANICS MS/MSD RECOVERY

Report No: 218081814
 Prep Method: 8330B
 Analytical Method: EPA 8330B

Parent Sample ID: WIL02IS01
 Prep Batch: 642680
 Analytical Batch: 643050

GCAL QC ID: 21808181403

| ANALYTE | UNITS | SPIKE ADDED | MSD RESULT | MSD % REC | # | % RPD | # | QC LIMITS | |
|------------------------------|-------|----------------|---------------|--------------|---|----------|---|-----------|--------|
| | | | | | | | | REC | RPD |
| 1,3,5-Trinitrobenzene | ug/kg | 1000 | 922 | 92 | | 6 | | 80 - 116 | 0 - 20 |
| 1,3-Dinitrobenzene | ug/kg | 1000 | 887 | 89 | | 3 | | 73 - 119 | 0 - 20 |
| 2,4,6-Trinitrotoluene | ug/kg | 1000 | 903 | 90 | | .6 | | 71 - 120 | 0 - 20 |
| 2,4-Dinitrotoluene | ug/kg | 1000 | 896 | 90 | | 4 | | 75 - 121 | 0 - 20 |
| 2,6-Dinitrotoluene | ug/kg | 1000 | 851 | 85 | | 15 | | 79 - 117 | 0 - 20 |
| 2-Amino-4,6-dinitrotoluene | ug/kg | 1000 | 910 | 91 | | 2 | | 71 - 123 | 0 - 20 |
| 2-Nitrotoluene | ug/kg | 1000 | 972 | 97 | | 7 | | 84 - 120 | 0 - 20 |
| 3,5-Dinitroaniline | ug/kg | 1000 | 872 | 87 | | 2 | | 86 - 118 | 0 - 20 |
| 3-Nitrotoluene | ug/kg | 1000 | 925 | 92 | | 6 | | 67 - 129 | 0 - 20 |
| 4-Amino-2,6-dinitrotoluene | ug/kg | 1000 | 958 | 96 | | 7 | | 64 - 127 | 0 - 20 |
| 4-Nitrotoluene | ug/kg | 1000 | 959 | 96 | | 2 | | 71 - 124 | 0 - 20 |
| HMX | ug/kg | 1000 | 798 | 80 | | 4 | | 74 - 124 | 0 - 20 |
| Nitrobenzene | ug/kg | 1000 | 872 | 87 | | 16 | | 80 - 128 | 0 - 20 |
| Nitroglycerin | ug/kg | 1000 | 879 | 88 | | 3 | | 73 - 124 | 0 - 20 |
| Pentaerythritol Tetranitrate | ug/kg | 1000 | 902 | 90 | | 9 | | 72 - 128 | 0 - 20 |
| RDX | ug/kg | 1000 | 847 | 85 | | 5 | | 67 - 129 | 0 - 20 |
| Tetryl | ug/kg | 1000 | 726 | 73 | | .8 | | 68 - 135 | 0 - 20 |

RPD : 0 out of 17 outside limits

Column to be used to flag recovery and RPD values with an asterisk

Spike Recovery: 0 out of 34 outside limits

* Values outside of QC limits

FORM III ORG-1

**DATA VALIDATION WORKSHEET
INORGANIC - ICPMS (Sb, Cu, Pb, Zn)**

SDG No.: 218081812+13+14

Project No.: 60520956

X USEPA CLP National Functional Guidelines for Superfund Inorganic Methods Data Review

X SW-846/DoD QSM v5.1

X Project QAPP/SAP

Project: Williston LTA

Reviewer: Naoum Tavantzis

Date: October 24, 2018

1.0 Chain of Custody/Sample Condition/Raw Data

| | | Yes | No | NA |
|-----|--|-----|----|----|
| 1.1 | Do Chain-of-Custody forms list all samples which were analyzed? | X | | |
| 1.2 | Are all Chain-of-Custody forms signed, indicating sample chain-of-custody was maintained? | X | | |
| 1.3 | Do the traffic Reports, chain-of-custody, and lab narrative indicate any problems with sample receipt, condition of samples, analytical problems or special circumstances affecting the quality of the data? | | X | |
| 1.4 | Does sample preservation, collection and storage meet method requirement? (For metal: water samples: with Nitric Acid to pH < 2, and soil/sediment samples: 4 °C ± 2 °C). Action: Professional judgement or J-(+)/R(-) | X | | |
| 1.5 | Are the digestion logs present and complete with pH values, sample weights, dilutions, final volumes. % solids (for soil samples), and preparation dates? For any missing or incomplete documentation, contact the laboratory for explanation/resubmittal. | X | | |
| 1.6 | Were the sediment samples dried and sieved appropriately? | X | | |
| 1.7 | Are the measurement read out records legible and complete (properly labeled, and include all samples and QC)? | X | | |

Note:

2.0 Holding Time

| | | Yes | No | NA |
|-----|--|-----|----|----|
| 2.1 | Have any technical holding times of 6 months, determined from date of collection to date of analysis, been exceeded? | | X | |
| | Action: Use professional judgement, then J-(+)/R(-) | | | |

Note:

3.0 Instrument Calibration

| | | Yes | No | NA |
|-------|--|-----|----|----|
| 3.1 | Are the Quartly LOD establishment forms provided for each instrument? | X | | |
| 3.2 | Is the signal/noise ratio less than 3:1? If not, is the determination repeated at a higher concentration? | X | | |
| 3.3 | Are sufficient standards of a blank + one standard & a RL standard OR 3 standards and a blank with one standard at the RL included in the calibration curve? If not, qualify with "R". | X | | |
| 3.3.a | If more than one standard is used, are the correlation coefficients > 0.995? Action: J(+)/UJ(-). | X | | |
| 3.4 | Was an initial calibration check standard (ICV) analyzed immediately after instrument system had been calibrated? Action: If no, all associated data are rejected "R". | X | | |
| 3.5 | Was continuing calibration (CCV) analyzed at a minimum frequency of 10% (every 10 samples) during and at the end of the analytical run? If not, document and flag based on professional judgement. | X | | |
| 3.6 | Are all calibration standard percent recoveries within the control limits of 90%-110%? ICV and CCV: | X | | |
| | < 75% 75%-89% 111%-125% 125%<R<160% >160% | | | |
| | Action: J-(+)/R(-) J-(+)/UJ(-) J+(+) J+(+)(maybe R(+)) R(+) | | | |
| 3.7 | Was the high-level check standard included within 10% of the true value? | X | | |

Note:

4.0 Blanks

| | | Yes | No | NA |
|-----|--|-----|----|----|
| 4.1 | Were method blank (MB) prepared at the appropriate frequency (one / 20 samples, batch, or matrix)? | X | | |
| 4.2 | Were calibration blanks (ICB and CCBs) analyzed immediately after each ICV and CCVs? | X | | |
| 4.3 | Are there reported ICB/CCBs values > DL? If yes, refer to action table. | | X | |
| 4.4 | Are there negative ICB/CCB blank results with the absolute value > LOD? If yes use 5x rule | | X | |
| 4.5 | Are there reported MB values > LOD? If yes use 5x rule | X | | |
| 4.6 | Are there negative method blank results with the absolute value > LOD? If yes use 5x rule | | X | |
| 4.7 | Are there reported field blank or equipment blank > + DL? If yes use 5x rule | X | | |

Note: MB in batch 642222 had a detection for zinc at 518 ug/kg. Equipment blank had detections for antimony (0.74 ug/L) and copper (0.26 ug/L)

5.0 Laboratory Control Sample (LCS)

| | | Yes | No | NA |
|-----|--|-----|----|----|
| 5.1 | Was an LCS prepared and analyzed at the correct frequency (one / 20 samples, batch, or matrix)? Action: If no, J(+)/R(-) any sample not associated with LCS results. | X | | |
| 5.2 | Is any LCS recovery outside the control limits? | | X | |
| | % Recoveries: <40% 40%-LCL% >UCL,<130% >150% | | | |
| | J-+)/R(-) J-+)/UJ(-) J+(+) R(+/-) | | | |

Note:

6.0 ICP Interference Check Sample (ICS)

| | | Yes | No | NA |
|-----|--|-----|----|----|
| 6.1 | Was ICS analyzed at beginning of each ICP run and every 12 hours? | X | | |
| 6.2 | Are the ICS AB recoveries within 80% - 120%? | X | | |
| 6.3 | Are the results for unspiked analytes (in ICS A) <LOQ? | X | | |
| 6.4 | If not, are the associated sample Al, Ca, Fe, and Mg concentrations less than the level in the ICS? | X | | |
| | Action: Not Spiked Analytes Spiked analytes | | | |
| | False negative False Positive < 50% 50%-79% 120% | | | |
| | ≥DL, but <10x [ICS] ≥DL | | | |
| | J-+)/UJ(-) J+(+) R(-) J-+)/UJ(-) J+(+) | | | |

Note:

7.0 Laboratory Duplicates (MD)

| | | Yes | No | NA |
|-----|--|-----|----|----|
| 7.1 | Were Laboratory duplicates prepared and analyzed at the correct frequency (one / 20 samples, batch, or matrix)? If no, J(+), using professional judgement, analytes not associated with duplicate results. | X | | |
| 7.2 | Are all RPDs less than 20%? If no, qualify all associated field samples J(+)/UJ(-) for the batch. | X | | |

Note:

8.0 Matrix/Matrix Spike Duplicate

| | | Yes | No | NA |
|-----|--|-----|----|----|
| 8.1 | Was a spiked sample prepared and analyzed at the correct frequency (one / 20 samples, batch, or matrix)? If not, J(+) with professional judgement. | X | | |
| 8.2 | Are any MS/MSD recovery outside the control limits? | X | | |
| | For all analytes with sample concentration ,4x spike concentration: | | | |
| | Matrix Spike %R <30% 30%-74% >125% | | | |
| | No PDS, Action: J-+)/R(-, reanalyze) J-+)/UJ(-) J+(+) | | | |
| | Matrix Spike %R <30% 30%-74% >125% | | | |
| | PDS%R <75% ≥75% <75% ≥75% <125% ≥125% | | | |
| | Action J-+)/R(-) J(+)/UJ(-) J-+)/UJ(-) J(+)/UJ(-) J(+) J+(+) | | | |
| 8.3 | Are all RPDs less than 20%? If no, qualify all associated field samples J(+) for the batch. | | X | |

Note: Matrix spikes on WIL02DA02A and WIL01IS02 - low recoveries for antimony in MS/MSD; WIL01IS02 high recovery for copper in MS.

9.0 ICP/AA Serial Dilutions (Not for Mercury Analysis)

| | | Yes | No | NA |
|-----|---|-----|----|----|
| 9.1 | Were serial dilutions performed? | X | | |
| 9.2 | Was a five-fold dilution performed? | X | | |
| 9.3 | Were results agree within 10% for [sample] > 50 X DL in the original sample? If no, J(+)/UJ(-) | | X | |
| 9.4 | Where dilution test fails or [sample]<50xLOD, was a post-digestion spike performed? Note any recoveries greater than ±25%D in DV report, but no data qualifying action is required. | X | | |
| 9.5 | When both dilution test and post-digestion spikes do not pass, was a method of standard additions used to quantitate the reported sample concentration? If not, ask lab to comment. | | | X |

Note: WIL02DA02A in analytical sequence 642309 displayed a percent difference greater than 10% for zinc at 10.9%.

10.0 Field Duplicate Samples

| | | Yes | No | NA |
|------|---|-----|----|----|
| 10.1 | Were any field duplicates submitted for metal analysis? | X | | |
| | For sample results > 5 x RL, a control limit of ≤35% RPD will be used. For sample results < 5 x RL, a control limit of 2 x RL will be used. | | | |
| 10.2 | Are all analyte duplicate results within control limits? If not, J(+) the parent samples. | X | | |

Note:

11.0 Result Verification/ Internal Standards/ Tune

| | | Yes | No | NA |
|------|--|-----|----|----|
| 11.1 | Are all DLs/RLs equal to or less than the reporting limits specified? | X | | |
| 11.2 | Were all results and detection limits for solid-matrix samples reported on a dry-weight basis? | X | | |
| 11.3 | Were all dilutions reflected in the positive results and detection limits? | X | | |
| 11.4 | Is there an Internal Standard associated with all analytes? R(+/-) | X | | |
| 11.5 | Were the Internal Standard recoveries within control limits? If not, J(+)/UJ(-) | X | | |
| 11.6 | Was a tune performed? If not, R(+/-) all associated samples. | X | | |
| 11.7 | Were the tunes run at a minimum of four times with RSD < 5% for analytes in solution? Action: J(+)/UJ(-) | X | | |
| 11.8 | Were the tune mass calibrations < 0.1 amu from the true value? Action: J(+)/UJ(-) | X | | |
| 11.9 | Was the resolution check peak width < 0.9 amu at 10% peak height? Action: J(+)/UJ(-) | X | | |

Note:

12.0 Completeness Calculation

| | | Yes | No | NA |
|------|---|-----|----|----|
| 12.1 | Is % completeness within the control limits? (Control limit 90%) | X | | |
| 12.2 | Number of samples: 16 | | | |
| 12.3 | Number of target compounds in each analysis: 4 | | | |
| 12.4 | Number of results rejected and not reported: 0 | | | |
| | % Completeness = (12.1.1 x 12.1.2 - 12.1.3) x 100/(12.1.1 x 12.1.2) | | | |
| | % Completeness = 100% | | | |

XIV
ANALYSIS RUN LOG

Report No: 218081813 Analytical Batch: 642309 Start Date: 08/20/18
 Instrument ID: ICPMS2 Analytical Method: EPA 6020B End Date: 08/20/18

| CLIENT SAMPLE ID | GCAL SAMPLE ID | PF | D/F | TIME | Analyte Symbols | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|------------------|----------------|----|-----|------|-----------------|----|----|----|----|---|----|----|----|----|----|----|----|----|----|----|----|----|----|---|----|----|----|----|----|----|----|----|---|----|----|--|
| | | | | | Al | Sb | As | Ba | Be | B | Cd | Ca | Cr | Co | Cu | Fe | Pb | Li | Mg | Mn | Hg | Mo | Ni | K | Se | Si | Ag | Na | Sr | Ti | Sn | Ti | V | Zn | Zr | |
| ITUNE | 1150 | * | 1 | 1252 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| ICALB | 1300 | * | 1 | 1305 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| ICAL2 | 1302 | * | 1 | 1309 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| ICAL4 | 1304 | * | 1 | 1313 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| ICAL5 | 1305 | * | 1 | 1317 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| ICAL6 | 1306 | * | 1 | 1320 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| ICV | 1600 | * | 1 | 1324 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| ICB | 1700 | * | 1 | 1342 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| LLCCV | 1803 | * | 1 | 1353 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| ICSA | 2000 | * | 1 | 1356 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| ICSAB | 2100 | * | 1 | 1400 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| LDR | 2500 | * | 1 | 1403 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| CCV | 1800 | * | 1 | 1734 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| CCB | 1900 | * | 1 | 1738 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| MB1841052 | 1841052 | * | 1 | 1827 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| LCS1841053 | 1841053 | * | 1 | 1831 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| CCV | 1800 | * | 1 | 1838 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| CCB | 1900 | * | 1 | 1841 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |

FORM XIV - IN

XIV
ANALYSIS RUN LOG

Report No: 218081812 Analytical Batch: 642309 Start Date: 08/20/18
 Instrument ID: ICPMS2 Analytical Method: EPA 6020B End Date: 08/20/18

| CLIENT SAMPLE ID | GCAL SAMPLE ID | PF | D/F | TIME | Analyte Symbols | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|------------------|----------------|----|-----|------|-----------------|----|----|----|----|---|----|----|----|----|----|----|----|----|----|----|----|----|----|---|----|----|----|----|----|----|----|----|
| | | | | | Al | Sb | As | Ba | Be | B | Cd | Ca | Cr | Co | Cu | Fe | Pb | Li | Mg | Mn | Hg | Mo | Ni | K | Se | Si | Ag | Na | Sr | Ti | Sh | Ti |
| ITUNE | 1150 | * | 1 | 1252 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| ICALB | 1300 | * | 1 | 1305 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| ICAL2 | 1302 | * | 1 | 1309 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| ICAL4 | 1304 | * | 1 | 1313 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| ICAL5 | 1305 | * | 1 | 1317 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| ICAL6 | 1306 | * | 1 | 1320 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| ICV | 1600 | * | 1 | 1324 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| ICB | 1700 | * | 1 | 1342 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| LLCCV | 1803 | * | 1 | 1353 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| ICSA | 2000 | * | 1 | 1356 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| ICSAB | 2100 | * | 1 | 1400 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| LDR | 2500 | * | 1 | 1403 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| CCV | 1800 | * | 1 | 1734 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| CCB | 1900 | * | 1 | 1738 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| WIL02DA02A | 21808181206 | * | 10 | 1802 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| WIL02DA02A MS | 21808181207 | * | 10 | 1806 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| WIL02DA02A MSD | 21808181208 | * | 10 | 1810 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| WIL02DA02APDS | 1841231 | * | 10 | 1813 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| WIL02DA02ASD | 1841232 | * | 50 | 1817 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| WIL02DA01A | 21808181209 | * | 10 | 1820 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| WIL02DA01B | 21808181210 | * | 10 | 1824 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| CCV | 1800 | * | 1 | 1838 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |
| CCB | 1900 | * | 1 | 1841 | | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | X |

FORM XIV - IN

VIII
ICP-MS TUNE

Report No: 218081812 GCAL QC ID: 1150
 Instrument ID: ICPMS2 Lab File ID: 2180820B_MS2.b\QCTune\2180820B_MS2-QCTu
 Analyst: AWG Analytical Batch: 642309
 Analysis Date: 08/20/18 Time: 1253 Analytical Method: EPA 6020B

| <i>ELEMENT - MASS</i> | <i>AVG MEASURED MASS (amu)</i> | <i>PEAK WIDTH AT 5% PEAK HEIGHT (amu)</i> | <i>%RSD</i> |
|-----------------------|--|---|-------------|
| Be-9 | 9.05 | .7887 | .7645 |
| Mg-24 | 24 | .7914 | .9917 |
| Mg-25 | 25 | .7928 | 1.2727 |
| Mg-26 | 26 | .8258 | .4753 |
| Co-59 | 58.95 | .7864 | .6309 |
| In-115 | 115 | .7729 | 1.2881 |
| Pb-206 | 206 | .8291 | 2.0969 |
| Pb-207 | 206.95 | .7931 | 1.9959 |
| Pb-208 | 207.95 | .8279 | 2.4434 |

II
INITIAL CALIBRATION VERIFICATION (ICV) STANDARD

| | |
|--|--|
| Report No: <u>218081812</u> | GCAL QC ID: <u>1600</u> |
| Instrument ID: <u>ICPMS2</u> | Lab File ID: <u>2180820B_MS2.b\009_ICV.d</u> |
| Analyst: <u>LWZ</u> | Analytical Batch: <u>642309</u> |
| Analysis Date: <u>08/20/18</u> Time: <u>1324</u> | Analytical Method: <u>EPA 6020B</u> |

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|----------------|-------------|--------------|------------------|----------|--------------|
| Antimony | 50.0 | 54.6 | 109 | | ug/L |
| Copper | 50.0 | 52.0 | 104 | | ug/L |
| Lead | 50.0 | 49.3 | 99 | | ug/L |
| Zinc | 1000 | 1010 | 101 | | ug/L |

CONTROL LIMITS 90-110%

FORM II - IN

III
INITIAL CALIBRATION BLANK

Report No: 218081812 Blank ID: 1700
Instrument ID: ICPMS2 Lab File ID: 2180820B_MS2.b\1211_ICB.d
Analyst: LWZ Analytical Batch: 642309
Analysis Date: 08/20/18 Time: 1342 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>RESULT</i> | <i>UNITS</i> | <i>Q</i> | <i>DL</i> | <i>LOD</i> | <i>LOQ</i> |
|----------------|---------------|--------------|----------|-----------|------------|------------|
| Antimony | 1.00 | ug/L | U | 0.50 | 1.00 | 2.00 |
| Copper | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Lead | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Zinc | 10.0 | ug/L | U | 5.00 | 10.0 | 20.0 |

FORM III - IN

II
 LOW LEVEL CONTINUING CALIBRATION VERIFICATION (LLCCV) STANDARD

| | |
|--|---|
| Report No: <u>218081812</u> | GCAL QC ID: <u>1803</u> |
| Instrument ID: <u>ICPMS2</u> | Lab File ID: <u>2180820B_MS2.b\121211CCV1.d</u> |
| Analyst: <u>LWZ</u> | Analytical Batch: <u>642309</u> |
| Analysis Date: <u>08/20/18</u> Time: <u>1353</u> | Analytical Method: <u>EPA 6020B</u> |

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|----------------|-------------|--------------|------------------|----------|--------------|
| Antimony | 2.00 | 1.90 | 95 | | ug/L |
| Copper | 1.00 | 1.06 | 106 | | ug/L |
| Lead | 1.00 | 0.990 | 99 | | ug/L |
| Zinc | 20.0 | 21.2 | 106 | | ug/L |

CONTROL LIMITS 80-120%

FORM II - IN

IV
ICPMS INTERFERENCE CHECKS

| | |
|--|---|
| Report No: <u>218081812</u> | ICSA \ AB ID: <u>2000 \ 2100</u> |
| Instrument ID: <u>ICPMS2</u> | Analytical Batch: <u>642309</u> |
| Analyst: <u>LWZ</u> | Analytical Method: <u>EPA 6020B</u> |
| Lab File ID ICSA1: <u>2180820B_MS2.b\121212\ICSA.d</u> | Lab File ID ICSAB1: <u>2180820B_MS2.b\121213\ICSB.d</u> |
| Lab File ID ICSA2: _____ | Lab File ID ICSAB2: _____ |

Concentration Units: ug/L

| Analyzed (A/AB): | | | 08/20/18 1356 | 08/20/18 1400 | | | | |
|------------------|--------|---------|---------------|---------------|------|-------|--------|----|
| ANALYTE | TRUE A | TRUE AB | ICSA1 | ICSAB1 | %R | ICSA2 | ICSAB2 | %R |
| Aluminum | 1000 | 1000 | 1010 | 999 | 100 | | | |
| Antimony | 0 | 0 | 0.039 | 0.031 | | | | |
| Arsenic | 0 | 10.0 | 0.0030 | 10.3 | 103 | | | |
| Barium | 0 | 0 | -0.0040 | 0.016 | | | | |
| Beryllium | 0 | 0 | -0.0010 | -0.0010 | | | | |
| Boron | 0 | 20.0 | 0.17 | 20.2 | 101 | | | |
| Cadmium | 0 | 10.0 | 0.017 | 10.1 | 101 | | | |
| Calcium | 3000 | 3000 | 2950 | 2980 | 99 | | | |
| Chromium | 0 | 20.0 | 0.088 | 20.8 | 104 | | | |
| Cobalt | 0 | 20.0 | 0.0040 | 21.0 | 105 | | | |
| Copper | 0 | 20.0 | 0.18 | 21.5 | 108 | | | |
| Iron | 2500 | 2500 | 2570 | 2510 | 100 | | | |
| Lead | 0 | 0 | 0.0040 | -0.0020 | | | | |
| Lithium | 0 | 20.0 | -0.069 | 23.9 | 120 | | | |
| Magnesium | 1000 | 1000 | 1020 | 1020 | 102 | | | |
| Manganese | 0 | 20.0 | 0.035 | 20.6 | 103 | | | |
| Molybdenum | 20.0 | 20.0 | 19.5 | 19.3 | 96 | | | |
| Nickel | 0 | 20.0 | 0.032 | 21.3 | 106 | | | |
| Potassium | 1000 | 1000 | 1010 | 1010 | 101 | | | |
| Selenium | 0 | 10.0 | -0.040 | 10.1 | 101 | | | |
| Silicon | 0 | 1000 | -4500 | -3500 | -351 | | | |
| Silver | 0 | 5.00 | 0.0060 | 5.15 | 103 | | | |
| Sodium | 2500 | 2500 | 2540 | 2540 | 102 | | | |
| Strontium | 0 | 10.0 | 0.10 | 10.1 | 101 | | | |
| Thallium | 0 | 0 | 0.0 | -0.0010 | | | | |
| Tin | 0 | 10.0 | 0.040 | 7.34 | 73 | | | |
| Titanium | 20.0 | 20.0 | 19.4 | 19.3 | 96 | | | |
| Vanadium | 0 | 20.0 | -0.015 | 17.6 | 88 | | | |
| Zinc | 0 | 20.0 | 0.029 | 21.1 | 106 | | | |
| Zirconium | 0 | 20.0 | 0.017 | 19.1 | 96 | | | |

FORM IV - IN

II
LINEAR DYNAMIC RANGE (LDR) STANDARD

Report No: 218081812 GCAL QC ID: 2500
Instrument ID: ICPMS2 Lab File ID: 2180820B_MS2.b\121214_QC1.d
Analyst: LWZ Analytical Batch: 642309
Analysis Date: 08/20/18 Time: 1403 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|----------------|-------------|--------------|------------------|----------|--------------|
| Antimony | 1000 | 924 | 92 | | ug/L |
| Copper | 1000 | 929 | 93 | | ug/L |
| Lead | 1000 | 915 | 92 | | ug/L |
| Zinc | 20000 | 18100 | 91 | | ug/L |

CONTROL LIMITS 90-110%

FORM II - IN

II
CONTINUING CALIBRATION VERIFICATION (CCV) STANDARD

| | |
|--|---|
| Report No: <u>218081812</u> | GCAL QC ID: <u>1800</u> |
| Instrument ID: <u>ICPMS2</u> | Lab File ID: <u>2180820B_MS2.b\121269_CCV.d</u> |
| Analyst: <u>LWZ</u> | Analytical Batch: <u>642309</u> |
| Analysis Date: <u>08/20/18</u> Time: <u>1734</u> | Analytical Method: <u>EPA 6020B</u> |

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|-----------------------|--------------------|---------------------|-------------------------|-----------------|---------------------|
| Antimony | 20.0 | 19.9 | 99 | | ug/L |
| Copper | 10.0 | 10.2 | 102 | | ug/L |
| Lead | 10.0 | 9.84 | 98 | | ug/L |
| Zinc | 200 | 201 | 100 | | ug/L |

CONTROL LIMITS 90-110%

III
CONTINUING CALIBRATION BLANK

Report No: 218081812 Blank ID: 1900
Instrument ID: ICPMS2 Lab File ID: 2180820B_MS2.b\121270_CCB.d
Analyst: LWZ Analytical Batch: 642309
Analysis Date: 08/20/18 Time: 1738 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>RESULT</i> | <i>UNITS</i> | <i>Q</i> | <i>DL</i> | <i>LOD</i> | <i>LOQ</i> |
|----------------|---------------|--------------|----------|-----------|------------|------------|
| Antimony | 1.00 | ug/L | U | 0.50 | 1.00 | 2.00 |
| Copper | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Lead | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Zinc | 10.0 | ug/L | U | 5.00 | 10.0 | 20.0 |

FORM III - IN

II
CONTINUING CALIBRATION VERIFICATION (CCV) STANDARD

Report No: 218081812 GCAL QC ID: 1800
Instrument ID: ICPMS2 Lab File ID: 2180820B_MS2.b\121287_CCV.d
Analyst: LWZ Analytical Batch: 642309
Analysis Date: 08/20/18 Time: 1838 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|-----------------------|--------------------|---------------------|-------------------------|-----------------|---------------------|
| Antimony | 20.0 | 19.7 | 99 | | ug/L |
| Copper | 10.0 | 10.4 | 104 | | ug/L |
| Lead | 10.0 | 9.61 | 96 | | ug/L |
| Zinc | 200 | 196 | 98 | | ug/L |

CONTROL LIMITS 90-110%

FORM II - IN

III
CONTINUING CALIBRATION BLANK

Report No: 218081812 Blank ID: 1900
Instrument ID: ICPMS2 Lab File ID: 2180820B_MS2.b\121288_CCB.d
Analyst: LWZ Analytical Batch: 642309
Analysis Date: 08/20/18 Time: 1841 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>RESULT</i> | <i>UNITS</i> | <i>Q</i> | <i>DL</i> | <i>LOD</i> | <i>LOQ</i> |
|----------------|---------------|--------------|----------|-----------|------------|------------|
| Antimony | 1.00 | ug/L | U | 0.50 | 1.00 | 2.00 |
| Copper | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Lead | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Zinc | 10.0 | ug/L | U | 5.00 | 10.0 | 20.0 |

FORM III - IN

XV (He)
ICP-MS INTERNAL STANDARDS RELATIVE INTENSITY SUMMARY

| | |
|------------------------------------|---------------------------------|
| Report No: <u>218081812</u> | Start Date: <u>08/20/18</u> |
| Instrument ID: <u>ICPMS2</u> | End Date: <u>08/20/18</u> |
| Analytical Method <u>EPA 6020B</u> | Analytical Batch: <u>642309</u> |

| |
|-----------------------------|
| Internal Standards %Rt For: |
|-----------------------------|

| CLIENT SAMPLE ID | GCAL SAMPLE ID | TIME | ISTD1 Q | ISTD2 Q | ISTD3 Q | ISTD4 Q | ISTD5 Q | ISTD6 Q | ISTD7 Q |
|------------------|----------------------|------|---------|---------|---------|---------|---------|---------|---------|
| WIL02DA02A | 21808181206 | 1802 | 107 | 105 | 104 | 110 | 104 | 105 | 108 |
| WIL02DA02A MS | 21808181207 | 1806 | 108 | 105 | 103 | 110 | 104 | 105 | 109 |
| WIL02DA02A MSD | 21808181208 | 1810 | 111 | 106 | 106 | 113 | 104 | 105 | 110 |
| WIL02DA02APDS | 1841231 | 1813 | 113 | 107 | 106 | 116 | 106 | 108 | 113 |
| WIL02DA02ASD | 1841232 | 1817 | 112 | 107 | 107 | 113 | 107 | 101 | 112 |
| WIL02DA01A | 21808181209 | 1820 | 112 | 106 | 105 | 113 | 104 | 106 | 110 |
| WIL02DA01B | 21808181210 | 1824 | 115 | 108 | 109 | 116 | 108 | 107 | 114 |

| | | |
|------------------------|-----------------------|----------------------|
| ISTD 1: Bismuth (He) | ISTD 4: Lutetium (He) | ISTD 7: Terbium (He) |
| ISTD 2: Germanium (He) | ISTD 5: Rhodium (He) | |
| ISTD 3: Indium (He) | ISTD 6: Scandium (He) | |

FORM XV - IN

XV (No Gas)
ICP-MS INTERNAL STANDARDS RELATIVE INTENSITY SUMMARY

| | |
|-------------------------------------|---------------------------------|
| Report No: <u>218081812</u> | Start Date: <u>08/20/18</u> |
| Instrument ID: <u>ICPMS2</u> | End Date: <u>08/20/18</u> |
| Analytical Method: <u>EPA 6020B</u> | Analytical Batch: <u>642309</u> |

| |
|-----------------------------|
| Internal Standards %RI For: |
|-----------------------------|

| CLIENT SAMPLE ID | GCAL SAMPLE ID | TIME | ISTD8 Q | ISTD9 Q | ISTD10 Q | ISTD11 Q | ISTD12 Q | ISTD13 Q | ISTD14 Q |
|------------------|----------------------|------|---------|---------|----------|----------|----------|----------|----------|
| WIL02DA02A | 21808181206 | 1802 | 92 | 99 | 98 | 99 | 97 | 106 | 99 |
| WIL02DA02A MS | 21808181207 | 1806 | 100 | 102 | 102 | 107 | 101 | 109 | 107 |
| WIL02DA02A MSD | 21808181208 | 1810 | 100 | 101 | 102 | 106 | 100 | 110 | 105 |
| WIL02DA02APDS | 1841231 | 1813 | 104 | 103 | 105 | 110 | 103 | 110 | 109 |
| WIL02DA02ASD | 1841232 | 1817 | 109 | 105 | 109 | 113 | 107 | 105 | 111 |
| WIL02DA01A | 21808181209 | 1820 | 104 | 103 | 106 | 110 | 102 | 110 | 108 |
| WIL02DA01B | 21808181210 | 1824 | 101 | 102 | 105 | 108 | 101 | 108 | 107 |

| | | |
|----------------------------|----------------------------|---------------------------|
| ISTD 8: Bismuth (No Gas) | ISTD 11: Lutetium (No Gas) | ISTD 14: Terbium (No Gas) |
| ISTD 9: Germanium (No Gas) | ISTD 12: Rhodium (No Gas) | |
| ISTD 10: Indium (No Gas) | ISTD 13: Scandium (No Gas) | |

FORM XV - IN

XV (He)
ICP-MS INTERNAL STANDARDS RELATIVE INTENSITY SUMMARY

| | |
|------------------------------------|---------------------------------|
| Report No: <u>218081813</u> | Start Date: <u>08/20/18</u> |
| Instrument ID: <u>ICPMS2</u> | End Date: <u>08/20/18</u> |
| Analytical Method <u>EPA 6020B</u> | Analytical Batch: <u>642309</u> |

Internal Standards %RI For:

| CLIENT SAMPLE ID | GCAL SAMPLE ID | TIME | ISTD1 Q | ISTD2 Q | ISTD3 Q | ISTD4 Q | ISTD5 Q | ISTD6 Q | ISTD7 Q |
|------------------|----------------------|------|---------|---------|---------|---------|---------|---------|---------|
| MB1841052 | 1841052 | 1827 | 114 | 102 | 107 | 111 | 108 | 96 | 110 |
| LCS1841053 | 1841053 | 1831 | 116 | 105 | 109 | 118 | 108 | 97 | 116 |

| | | |
|------------------------|-----------------------|----------------------|
| ISTD 1: Bismuth (He) | ISTD 4: Lutetium (He) | ISTD 7: Terbium (He) |
| ISTD 2: Germanium (He) | ISTD 5: Rhodium (He) | |
| ISTD 3: Indium (He) | ISTD 6: Scandium (He) | |

FORM XV - IN

XV (No Gas)

ICP-MS INTERNAL STANDARDS RELATIVE INTENSITY SUMMARY

Report No: 218081813 Start Date: 08/20/18
 Instrument ID: ICPMS2 End Date: 08/20/18
 Analytical Method: EPA 6020B Analytical Batch: 642309

Internal Standards %RI For:

| CLIENT SAMPLE ID | GCAL SAMPLE ID | TIME | ISTD8 Q | ISTD9 Q | ISTD10 Q | ISTD11 Q | ISTD12 Q | ISTD13 Q | ISTD14 Q |
|------------------|----------------------|------|---------|---------|----------|----------|----------|----------|----------|
| MB1841052 | 1841052 | 1827 | 105 | 101 | 107 | 109 | 107 | 102 | 108 |
| LCS1841053 | 1841053 | 1831 | 108 | 105 | 112 | 117 | 107 | 104 | 115 |

ISTD 8: Bismuth (No Gas) ISTD 11: Lutetium (No Gas) ISTD 14: Terbium (No Gas)
 ISTD 9: Germanium (No Gas) ISTD 12: Rhodium (No Gas)
 ISTD 10: Indium (No Gas) ISTD 13: Scandium (No Gas)

FORM XV - IN

XIV
ANALYSIS RUN LOG

Report No: 218081812 Analytical Batch: 642381 Start Date: 08/21/18
 Instrument ID: ICPMS1 Analytical Method: EPA 6020B End Date: 08/21/18

| CLIENT SAMPLE ID | GCAL SAMPLE ID | PF | D/F | TIME | Analyte Symbols | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|------------------|----------------|----|-----|------|-----------------|----|----|----|----|---|----|----|----|----|----|----|----|----|----|----|----|----|----|---|----|----|----|----|----|----|----|----|---|----|----|---|---|
| | | | | | Al | Sb | As | Ba | Be | B | Cd | Ca | Cr | Co | Cu | Fe | Pb | Li | Mg | Mn | Hg | Mo | Ni | K | Se | Si | Ag | Na | Sr | Tl | Sn | Ti | V | Zn | Zr | | |
| ITUNE | 1150 | * | 1 | 0930 | | X | | | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |
| IICALB | 1300 | * | 1 | 1023 | | X | | | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |
| IICAL2 | 1302 | * | 1 | 1027 | | X | | | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |
| IICAL4 | 1304 | * | 1 | 1033 | | X | | | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |
| IICAL5 | 1305 | * | 1 | 1037 | | X | | | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |
| IICAL6 | 1306 | * | 1 | 1042 | | X | | | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |
| ICV | 1600 | * | 1 | 1046 | | X | | | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |
| ICB | 1700 | * | 1 | 1055 | | X | | | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |
| LLCCV | 1803 | * | 1 | 1108 | | X | | | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |
| ICSA | 2000 | * | 1 | 1112 | | X | | | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |
| ICSAB | 2100 | * | 1 | 1117 | | X | | | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |
| LDR | 2500 | * | 1 | 1121 | | X | | | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |
| MB1840818 | 1840818 | * | 1 | 1138 | | X | | | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |
| LCS1840819 | 1840819 | * | 1 | 1143 | | X | | | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |
| CCV | 1800 | * | 1 | 1156 | | X | | | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |
| CCB | 1900 | * | 1 | 1200 | | X | | | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |

FORM XIV - IN

XIV
ANALYSIS RUN LOG

Report No: 218081813 Analytical Batch: 642381 Start Date: 08/21/18
 Instrument ID: ICPMS1 Analytical Method: EPA 6020B End Date: 08/21/18

| CLIENT SAMPLE ID | GCAL SAMPLE ID | PF | D/F | TIME | Analyte Symbols | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|------------------|----------------|----|-----|------|-----------------|----|----|----|----|---|----|----|----|----|----|----|----|----|----|----|----|----|----|----|---|----|----|----|----|----|----|----|----|---|----|----|---|--|
| | | | | | Al | Sb | As | Ba | Be | B | Cd | Ca | Cr | Co | Cu | Fe | Pb | Pb | Li | Mg | Mn | Hg | Mo | Ni | K | Se | Si | Ag | Na | Sr | Tl | Sn | Ti | V | Zn | Zr | | |
| ITUNE | 1150 | * | 1 | 0930 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | |
| ICALB | 1300 | * | 1 | 1023 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | |
| ICAL2 | 1302 | * | 1 | 1027 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | |
| ICAL4 | 1304 | * | 1 | 1033 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | |
| ICAL5 | 1305 | * | 1 | 1037 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | |
| ICAL6 | 1306 | * | 1 | 1042 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | |
| ICV | 1600 | * | 1 | 1046 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | |
| ICB | 1700 | * | 1 | 1055 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | |
| LLCCV | 1803 | * | 1 | 1108 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | |
| ICSA | 2000 | * | 1 | 1112 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | |
| ICSAB | 2100 | * | 1 | 1117 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | |
| LDR | 2500 | * | 1 | 1121 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | |
| WILO3IS00 | 21808181303 | * | 1 | 1151 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | |
| CCV | 1800 | * | 1 | 1156 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| CCB | 1900 | * | 1 | 1200 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |

FORM XIV - IN

VIII
ICP-MS TUNE

Report No: 218081812 GCAL QC ID: 1150
Instrument ID: ICPMS1 Lab File ID: 2180821A_MS1.b\QCTune\2180821A_MS1-QCTu
Analyst: AWG Analytical Batch: 642381
Analysis Date: 08/21/18 Time: 0931 Analytical Method: EPA 6020B

| <i>ELEMENT - MASS</i> | <i>AVG MEASURED MASS (amu)</i> | <i>PEAK WIDTH AT 5% PEAK HEIGHT (amu)</i> | <i>%RSD</i> |
|-----------------------|--|---|-------------|
| Be-9 | 9 | .7803 | .2982 |
| Mg-24 | 23.9 | .7774 | .8215 |
| Mg-25 | 24.9 | .7735 | .9524 |
| Mg-26 | 25.9 | .7497 | .5087 |
| Co-59 | 58.95 | .7688 | .6382 |
| In-115 | 115 | .7517 | .5548 |
| Pb-206 | 206 | .7916 | .8925 |
| Pb-207 | 207 | .8145 | 1.0512 |
| Pb-208 | 208 | .8111 | .8122 |

FORM VIII - IN

II
INITIAL CALIBRATION VERIFICATION (ICV) STANDARD

| | |
|--|--|
| Report No: <u>218081812</u> | GCAL QC ID: <u>1600</u> |
| Instrument ID: <u>ICPMS1</u> | Lab File ID: <u>2180821A_MS1.b\009_ICV.d</u> |
| Analyst: <u>LWZ</u> | Analytical Batch: <u>642381</u> |
| Analysis Date: <u>08/21/18</u> Time: <u>1046</u> | Analytical Method: <u>EPA 6020B</u> |

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|-----------------------|--------------------|---------------------|-------------------------|-----------------|---------------------|
| Antimony | 50.0 | 52.2 | 104 | | ug/L |
| Copper | 50.0 | 50.1 | 100 | | ug/L |
| Lead | 50.0 | 48.8 | 98 | | ug/L |
| Zinc | 1000 | 994 | 99 | | ug/L |

CONTROL LIMITS 90-110%

FORM II - IN

III
INITIAL CALIBRATION BLANK

Report No: 218081812 Blank ID: 1700
Instrument ID: ICPMS1 Lab File ID: 2180821A_MS1.b\011_ICB.d
Analyst: LWZ Analytical Batch: 642381
Analysis Date: 08/21/18 Time: 1055 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>RESULT</i> | <i>UNITS</i> | <i>Q</i> | <i>DL</i> | <i>LOD</i> | <i>LOQ</i> |
|----------------|---------------|--------------|----------|-----------|------------|------------|
| Antimony | 1.00 | ug/L | U | 0.50 | 1.00 | 2.00 |
| Copper | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Lead | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Zinc | 10.0 | ug/L | U | 5.00 | 10.0 | 20.0 |

FORM III - IN

II
LOW LEVEL CONTINUING CALIBRATION VERIFICATION (LLCCV) STANDARD

Report No: 218081812 GCAL QC ID: 1803
Instrument ID: ICPMS1 Lab File ID: 2180821A_MS1.b\014CCV1.d
Analyst: LWZ Analytical Batch: 642381
Analysis Date: 08/21/18 Time: 1108 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|----------------|-------------|--------------|------------------|----------|--------------|
| Antimony | 2.00 | 2.07 | 103 | | ug/L |
| Copper | 1.00 | 0.810 | 81 | | ug/L |
| Lead | 1.00 | 0.980 | 98 | | ug/L |
| Zinc | 20.0 | 20.8 | 104 | | ug/L |

CONTROL LIMITS 80-120%

FORM II - IN

IV
ICPMS INTERFERENCE CHECKS

| | |
|---|--|
| Report No: <u>218081812</u> | ICSA \ AB ID: <u>2000 \ 2100</u> |
| Instrument ID: <u>ICPMS1</u> | Analytical Batch: <u>642381</u> |
| Analyst: <u>LWZ</u> | Analytical Method: <u>EPA 6020B</u> |
| Lab File ID ICSA1: <u>2180821A_MS1.b\015\ICSA.d</u> | Lab File ID ICSAB1: <u>2180821A_MS1.b\016\ICSB.d</u> |
| Lab File ID ICSA2: _____ | Lab File ID ICSAB2: _____ |

Concentration Units: ug/L

| ANALYTE | Analyzed (A/AB): | | 08/21/18 1112 | | 08/21/18 1117 | | ICSA2 | ICSAB2 | %R |
|------------|------------------|---------|---------------|---------|---------------|--|-------|--------|----|
| | TRUE A | TRUE AB | ICSA1 | ICSAB1 | %R | | | | |
| Aluminum | 1000 | 1000 | 1010 | 1010 | 101 | | | | |
| Antimony | 0 | 0 | 0.063 | 0.065 | | | | | |
| Arsenic | 0 | 10.0 | -0.0010 | 10.3 | 103 | | | | |
| Barium | 0 | 0 | 0.039 | 0.034 | | | | | |
| Beryllium | 0 | 0 | 0.0 | 0.0 | | | | | |
| Boron | 0 | 20.0 | 1.21 | 20.2 | 101 | | | | |
| Cadmium | 0 | 10.0 | 0.027 | 10.1 | 101 | | | | |
| Calcium | 3000 | 3000 | 2960 | 3010 | 100 | | | | |
| Chromium | 0 | 20.0 | 0.0060 | 20.4 | 102 | | | | |
| Cobalt | 0 | 20.0 | 0.014 | 20.8 | 104 | | | | |
| Copper | 0 | 20.0 | -0.17 | 20.2 | 101 | | | | |
| Iron | 2500 | 2500 | 2530 | 2510 | 100 | | | | |
| Lead | 0 | 0 | -0.0040 | -0.0090 | | | | | |
| Lithium | 0 | 20.0 | -0.051 | 22.7 | 114 | | | | |
| Magnesium | 1000 | 1000 | 1010 | 1010 | 101 | | | | |
| Manganese | 0 | 20.0 | 0.031 | 20.8 | 104 | | | | |
| Molybdenum | 20.0 | 20.0 | 19.2 | 19.0 | 95 | | | | |
| Nickel | 0 | 20.0 | 0.022 | 21.2 | 106 | | | | |
| Potassium | 1000 | 1000 | 1020 | 992 | 99 | | | | |
| Selenium | 0 | 10.0 | 0.063 | 10.1 | 101 | | | | |
| Silicon | 0 | 1000 | -23 | 1100 | 110 | | | | |
| Silver | 0 | 5.00 | 0.0070 | 5.15 | 103 | | | | |
| Sodium | 2500 | 2500 | 2590 | 2570 | 103 | | | | |
| Strontium | 0 | 10.0 | 0.094 | 9.55 | 96 | | | | |
| Thallium | 0 | 0 | -0.067 | -0.069 | | | | | |
| Tin | 0 | 10.0 | 0.010 | 7.32 | 73 | | | | |
| Titanium | 20.0 | 20.0 | 20.5 | 20.0 | 100 | | | | |
| Vanadium | 0 | 20.0 | -0.0020 | 17.4 | 87 | | | | |
| Zinc | 0 | 20.0 | 0.27 | 21.6 | 108 | | | | |
| Zirconium | 0 | 20.0 | 0.027 | 18.8 | 94 | | | | |

FORM IV - IN

II
LINEAR DYNAMIC RANGE (LDR) STANDARD

Report No: 218081812 GCAL QC ID: 2500
Instrument ID: ICPMS1 Lab File ID: 2180821A_MS1.b\017_QC1.d
Analyst: LWZ Analytical Batch: 642381
Analysis Date: 08/21/18 Time: 1121 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|----------------|-------------|--------------|------------------|----------|--------------|
| Antimony | 1000 | 901 | 90 | | ug/L |
| Copper | 1000 | 1030 | 103 | | ug/L |
| Lead | 1000 | 928 | 93 | | ug/L |
| Zinc | 20000 | 18500 | 93 | | ug/L |

CONTROL LIMITS 90-110%

FORM II - IN

II
CONTINUING CALIBRATION VERIFICATION (CCV) STANDARD

| | |
|--|--|
| Report No: <u>218081812</u> | GCAL QC ID: <u>1800</u> |
| Instrument ID: <u>ICPMS1</u> | Lab File ID: <u>2180821A_MS1.b\024_CCV.d</u> |
| Analyst: <u>LWZ</u> | Analytical Batch: <u>642381</u> |
| Analysis Date: <u>08/21/18</u> Time: <u>1156</u> | Analytical Method: <u>EPA 6020B</u> |

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|-----------------------|--------------------|---------------------|-------------------------|-----------------|---------------------|
| Antimony | 20.0 | 20.0 | 100 | | ug/L |
| Copper | 10.0 | 10.1 | 101 | | ug/L |
| Lead | 10.0 | 9.57 | 96 | | ug/L |
| Zinc | 200 | 205 | 102 | | ug/L |

CONTROL LIMITS 90-110%

FORM II - IN

III
CONTINUING CALIBRATION BLANK

Report No: 218081812 Blank ID: 1900
Instrument ID: ICPMS1 Lab File ID: 2180821A_MS1.b\025_CCB.d
Analyst: LWZ Analytical Batch: 642381
Analysis Date: 08/21/18 Time: 1200 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>RESULT</i> | <i>UNITS</i> | <i>Q</i> | <i>DL</i> | <i>LOD</i> | <i>LOQ</i> |
|----------------|---------------|--------------|----------|-----------|------------|------------|
| Antimony | 1.00 | ug/L | U | 0.50 | 1.00 | 2.00 |
| Copper | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Lead | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Zinc | 10.0 | ug/L | U | 5.00 | 10.0 | 20.0 |

FORM III - IN

XV (He)
ICP-MS INTERNAL STANDARDS RELATIVE INTENSITY SUMMARY

| | |
|-------------------------------------|---------------------------------|
| Report No: <u>218081812</u> | Start Date: <u>08/21/18</u> |
| Instrument ID: <u>ICPMS1</u> | End Date: <u>08/21/18</u> |
| Analytical Method: <u>EPA 6020B</u> | Analytical Batch: <u>642381</u> |

Internal Standards %RI For:

| CLIENT SAMPLE ID | GCAL SAMPLE ID | TIME | ISTD1 Q | ISTD2 Q | ISTD3 Q | ISTD4 Q | ISTD5 Q | ISTD6 Q | ISTD7 Q |
|------------------|----------------------|------|---------|---------|---------|---------|---------|---------|---------|
| MB1840818 | 1840818 | 1138 | 103 | 114 | 113 | 119 | 108 | 117 | 109 |
| LCS1840819 | 1840819 | 1143 | 98 | 107 | 108 | 116 | 104 | 113 | 105 |

| | | |
|------------------------|-----------------------|----------------------|
| ISTD 1: Bismuth (He) | ISTD 4: Lutetium (He) | ISTD 7: Terbium (He) |
| ISTD 2: Germanium (He) | ISTD 5: Rhodium (He) | |
| ISTD 3: Indium (He) | ISTD 6: Scandium (He) | |

FORM XV - IN

XV (No Gas)
ICP-MS INTERNAL STANDARDS RELATIVE INTENSITY SUMMARY

| | |
|-------------------------------------|---------------------------------|
| Report No: <u>218081812</u> | Start Date: <u>08/21/18</u> |
| Instrument ID: <u>ICPMS1</u> | End Date: <u>08/21/18</u> |
| Analytical Method: <u>EPA 6020B</u> | Analytical Batch: <u>642381</u> |

Internal Standards %RI For:

| CLIENT SAMPLE ID | GCAL SAMPLE ID | TIME | ISTD8 Q | ISTD9 Q | ISTD10 Q | ISTD11 Q | ISTD12 Q | ISTD13 Q | ISTD14 Q |
|------------------|----------------------|------|---------|---------|----------|----------|----------|----------|----------|
| MB1840818 | 1840818 | 1138 | 95 | 105 | 101 | 97 | 104 | 108 | 100 |
| LCS1840819 | 1840819 | 1143 | 93 | 104 | 100 | 97 | 100 | 107 | 99 |

| | | |
|----------------------------|----------------------------|---------------------------|
| ISTD 8: Bismuth (No Gas) | ISTD 11: Lutetium (No Gas) | ISTD 14: Terbium (No Gas) |
| ISTD 9: Germanium (No Gas) | ISTD 12: Rhodium (No Gas) | |
| ISTD 10: Indium (No Gas) | ISTD 13: Scandium (No Gas) | |

FORM XV - IN

XV (He)
ICP-MS INTERNAL STANDARDS RELATIVE INTENSITY SUMMARY

| | |
|------------------------------------|---------------------------------|
| Report No: <u>218081813</u> | Start Date: <u>08/21/18</u> |
| Instrument ID: <u>ICPMS1</u> | End Date: <u>08/21/18</u> |
| Analytical Method <u>EPA 6020B</u> | Analytical Batch: <u>642381</u> |

Internal Standards %RI For:

| CLIENT SAMPLE ID | GCAL SAMPLE ID | TIME | ISTD1 Q | ISTD2 Q | ISTD3 Q | ISTD4 Q | ISTD5 Q | ISTD6 Q | ISTD7 Q |
|------------------|----------------------|------|---------|---------|---------|---------|---------|---------|---------|
| WIL03IS00 | 21808181303 | 1151 | 100 | 102 | 101 | 108 | 103 | 104 | 101 |

| | | |
|------------------------|-----------------------|----------------------|
| ISTD 1: Bismuth (He) | ISTD 4: Lutetium (He) | ISTD 7: Terbium (He) |
| ISTD 2: Germanium (He) | ISTD 5: Rhodium (He) | |
| ISTD 3: Indium (He) | ISTD 6: Scandium (He) | |

FORM XV - IN

XV (No Gas)
ICP-MS INTERNAL STANDARDS RELATIVE INTENSITY SUMMARY

| | |
|-------------------------------------|---------------------------------|
| Report No: <u>218081813</u> | Start Date: <u>08/21/18</u> |
| Instrument ID: <u>ICPMS1</u> | End Date: <u>08/21/18</u> |
| Analytical Method: <u>EPA 6020B</u> | Analytical Batch: <u>642381</u> |

Internal Standards %RI For:

| CLIENT SAMPLE ID | GCAL SAMPLE ID | TIME | ISTD8 Q | ISTD9 Q | ISTD10 Q | ISTD11 Q | ISTD12 Q | ISTD13 Q | ISTD14 Q |
|------------------|----------------------|------|---------|---------|----------|----------|----------|----------|----------|
| WIL03IS00 | 21808181303 | 1151 | 100 | 105 | 102 | 99 | 104 | 104 | 101 |

| | | |
|----------------------------|----------------------------|---------------------------|
| ISTD 8: Bismuth (No Gas) | ISTD 11: Lutetium (No Gas) | ISTD 14: Terbium (No Gas) |
| ISTD 9: Germanium (No Gas) | ISTD 12: Rhodium (No Gas) | |
| ISTD 10: Indium (No Gas) | ISTD 13: Scandium (No Gas) | |

FORM XV - IN

XIV
ANALYSIS RUN LOG

Report No: 218081813 Analytical Batch: 642536 Start Date: 08/23/18
 Instrument ID: ICPMS1 Analytical Method: EPA 6020B End Date: 08/23/18

| CLIENT SAMPLE ID | GCAL SAMPLE ID | PF | D/F | TIME | Analyte Symbols | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|------------------|----------------|----|-----|------|-----------------|----|----|----|----|---|----|----|----|----|----|----|----|----|----|----|----|----|----|---|----|----|----|----|----|----|----|----|---|----|----|---|---|---|---|
| | | | | | Al | Sb | As | Ba | Be | B | Cd | Ca | Cr | Co | Cu | Fe | Pb | Li | Mg | Mn | Hg | Mo | Ni | K | Se | Si | Ag | Na | Sr | Ti | Sn | Ti | V | Zn | Zr | | | | |
| ITUNE | 1150 | * | 1 | 0902 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X | | |
| ICALB | 1300 | * | 1 | 1014 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X | | |
| ICAL2 | 1302 | * | 1 | 1018 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X | | |
| ICAL4 | 1304 | * | 1 | 1023 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X | | |
| ICAL5 | 1305 | * | 1 | 1027 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X | | |
| ICAL6 | 1306 | * | 1 | 1032 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X | | |
| ICV | 1600 | * | 1 | 1036 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X | | |
| ICB | 1700 | * | 1 | 1041 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X | |
| LLCCV | 1803 | * | 1 | 1045 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X | |
| ICSA | 2000 | * | 1 | 1049 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X | |
| ICSAB | 2100 | * | 1 | 1054 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X | |
| LDR | 2500 | * | 1 | 1058 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X | |
| MB1841866 | 1841866 | * | 1 | 1129 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X | |
| LCS1841867 | 1841867 | * | 1 | 1134 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X | |
| CCV | 1800 | * | 1 | 1200 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X | |
| CCB | 1900 | * | 1 | 1205 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X |
| WIL01IS02 | 21808181302 | * | 10 | 1209 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X |
| WIL01IS02MS | 1842047 | * | 10 | 1214 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X |
| WIL01IS02MSD | 1842048 | * | 10 | 1218 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X |
| WIL01IS02PDS | 1842322 | * | 10 | 1222 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X |
| WIL01IS02SD | 1842323 | * | 50 | 1227 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X |
| WIL04IS03 | 21808181304 | * | 10 | 1231 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X |
| WIL04IS01 | 21808181305 | * | 10 | 1235 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X |
| WIL04IS02 | 21808181306 | * | 10 | 1240 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X |
| CCV | 1800 | * | 1 | 1328 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X |
| CCB | 1900 | * | 1 | 1332 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | X |

FORM XIV - IN

VIII
ICP-MS TUNE

Report No: 218081813 GCAL QC ID: 1150
Instrument ID: ICPMS1 Lab File ID: 2180823A_MS1.b\QCTune\2180823A_MS1-QCTu
Analyst: AWG Analytical Batch: 642536
Analysis Date: 08/23/18 Time: 0902 Analytical Method: EPA 6020B

| <i>ELEMENT - MASS</i> | <i>AVG MEASURED MASS (amu)</i> | <i>PEAK WIDTH AT 5% PEAK HEIGHT (amu)</i> | <i>%RSD</i> |
|-----------------------|--|---|-------------|
| Be-9 | 9 | .7775 | .4292 |
| Mg-24 | 23.9 | .7434 | .3034 |
| Mg-25 | 24.95 | .735 | 1.0936 |
| Mg-26 | 25.9 | .7422 | 1.4023 |
| Co-59 | 58.95 | .7335 | .5488 |
| In-115 | 115 | .7328 | .6735 |
| Pb-206 | 206 | .7673 | 1.1793 |
| Pb-207 | 207 | .7794 | .977 |
| Pb-208 | 208 | .8013 | .4759 |

FORM VIII - IN

II
INITIAL CALIBRATION VERIFICATION (ICV) STANDARD

| | |
|--|--|
| Report No: <u>218081813</u> | GCAL QC ID: <u>1600</u> |
| Instrument ID: <u>ICPMS1</u> | Lab File ID: <u>2180823A_MS1.b\017_ICV.d</u> |
| Analyst: <u>LWZ</u> | Analytical Batch: <u>642536</u> |
| Analysis Date: <u>08/23/18</u> Time: <u>1036</u> | Analytical Method: <u>EPA 6020B</u> |

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|-----------------------|--------------------|---------------------|-------------------------|-----------------|---------------------|
| Antimony | 50.0 | 49.7 | 99 | | ug/L |
| Copper | 50.0 | 50.9 | 102 | | ug/L |
| Lead | 50.0 | 48.8 | 98 | | ug/L |
| Zinc | 1000 | 1000 | 100 | | ug/L |

CONTROL LIMITS 90-110%

FORM II - IN

III
INITIAL CALIBRATION BLANK

Report No: 218081813 Blank ID: 1700
Instrument ID: ICPMS1 Lab File ID: 2180823A_MS1.b\018_ICB.d
Analyst: LWZ Analytical Batch: 642536
Analysis Date: 08/23/18 Time: 1041 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>RESULT</i> | <i>UNITS</i> | <i>Q</i> | <i>DL</i> | <i>LOD</i> | <i>LOQ</i> |
|----------------|---------------|--------------|----------|-----------|------------|------------|
| Antimony | 1.00 | ug/L | U | 0.50 | 1.00 | 2.00 |
| Copper | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Lead | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Zinc | 10.0 | ug/L | U | 5.00 | 10.0 | 20.0 |

II
LOW LEVEL CONTINUING CALIBRATION VERIFICATION (LLCCV) STANDARD

Report No: 218081813 GCAL QC ID: 1803
Instrument ID: ICPMS1 Lab File ID: 2180823A_MS1.b\019CCV1.d
Analyst: LWZ Analytical Batch: 642536
Analysis Date: 08/23/18 Time: 1045 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|----------------|-------------|--------------|------------------|----------|--------------|
| Antimony | 2.00 | 1.95 | 98 | | ug/L |
| Copper | 1.00 | 0.960 | 96 | | ug/L |
| Lead | 1.00 | 0.980 | 98 | | ug/L |
| Zinc | 20.0 | 20.4 | 102 | | ug/L |

CONTROL LIMITS 80-120%

FORM II - IN

IV
ICPMS INTERFERENCE CHECKS

| | |
|--|---|
| Report No: <u>218081813</u> | ICSA \ AB ID: <u>2000 \ 2100</u> |
| Instrument ID: <u>ICPMS1</u> | Analytical Batch: <u>642536</u> |
| Analyst: <u>LWZ</u> | Analytical Method: <u>EPA 6020B</u> |
| Lab File ID ICSA1: <u>2180823A_MS1.b\020ICSA.d</u> | Lab File ID ICSAB1: <u>2180823A_MS1.b\021ICSB.d</u> |
| Lab File ID ICSA2: _____ | Lab File ID ICSAB2: _____ |

Concentration Units: ug/L

| Analyzed (A/AB): | | | 08/23/18 1049 | 08/23/18 1054 | | | | |
|------------------|--------|---------|---------------|---------------|-----|-------|--------|----|
| ANALYTE | TRUE A | TRUE AB | ICSA1 | ICSAB1 | %R | ICSA2 | ICSAB2 | %R |
| Aluminum | 1000 | 1000 | 1010 | 999 | 100 | | | |
| Antimony | 0 | 0 | -0.0020 | -0.012 | | | | |
| Arsenic | 0 | 10.0 | 0.0040 | 10.0 | 100 | | | |
| Barium | 0 | 0 | 0.025 | 0.025 | | | | |
| Beryllium | 0 | 0 | -0.0010 | -0.0050 | | | | |
| Boron | 0 | 20.0 | 0.42 | 19.5 | 98 | | | |
| Cadmium | 0 | 10.0 | 0.020 | 10.1 | 101 | | | |
| Calcium | 3000 | 3000 | 2870 | 2960 | 99 | | | |
| Chromium | 0 | 20.0 | 0.0 | 19.9 | 100 | | | |
| Cobalt | 0 | 20.0 | 0.0040 | 20.2 | 101 | | | |
| Copper | 0 | 20.0 | -0.098 | 20.6 | 103 | | | |
| Iron | 2500 | 2500 | 2490 | 2480 | 99 | | | |
| Lead | 0 | 0 | 0.0010 | -0.010 | | | | |
| Lithium | 0 | 20.0 | -0.026 | 22.4 | 112 | | | |
| Magnesium | 1000 | 1000 | 997 | 991 | 99 | | | |
| Manganese | 0 | 20.0 | 0.028 | 20.2 | 101 | | | |
| Molybdenum | 20.0 | 20.0 | 19.5 | 19.3 | 96 | | | |
| Nickel | 0 | 20.0 | -0.0080 | 20.5 | 102 | | | |
| Potassium | 1000 | 1000 | 1010 | 998 | 100 | | | |
| Selenium | 0 | 10.0 | 0.021 | 10.4 | 104 | | | |
| Silicon | 0 | 1000 | -1.4 | 1090 | 109 | | | |
| Silver | 0 | 5.00 | 0.0070 | 5.16 | 103 | | | |
| Sodium | 2500 | 2500 | 2600 | 2520 | 101 | | | |
| Strontium | 0 | 10.0 | 0.10 | 9.92 | 99 | | | |
| Thallium | 0 | 0 | -0.026 | -0.034 | | | | |
| Tin | 0 | 10.0 | -0.038 | 7.41 | 74 | | | |
| Titanium | 20.0 | 20.0 | 20.0 | 20.8 | 104 | | | |
| Vanadium | 0 | 20.0 | -0.011 | 17.1 | 86 | | | |
| Zinc | 0 | 20.0 | -0.13 | 19.8 | 99 | | | |
| Zirconium | 0 | 20.0 | 0.021 | 19.1 | 96 | | | |

FORM IV - IN

II
LINEAR DYNAMIC RANGE (LDR) STANDARD

Report No: 218081813 GCAL QC ID: 2500
Instrument ID: ICPMS1 Lab File ID: 2180823A_MS1.b\022_QC1.d
Analyst: LWZ Analytical Batch: 642536
Analysis Date: 08/23/18 Time: 1058 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|----------------|-------------|--------------|------------------|----------|--------------|
| Antimony | 1000 | 909 | 91 | | ug/L |
| Copper | 1000 | 923 | 92 | | ug/L |
| Lead | 1000 | 952 | 95 | | ug/L |
| Zinc | 20000 | 19600 | 98 | | ug/L |

CONTROL LIMITS 90-110%

FORM II - IN

II
CONTINUING CALIBRATION VERIFICATION (CCV) STANDARD

| | | |
|--|--|--|
| Report No: <u>218081813</u> | GCAL QC ID: <u>1800</u> | |
| Instrument ID: <u>ICPMS1</u> | Lab File ID: <u>2180823A_MS1.b\036_CCV.d</u> | |
| Analyst: <u>LWZ</u> | Analytical Batch: <u>642536</u> | |
| Analysis Date: <u>08/23/18</u> Time: <u>1200</u> | Analytical Method: <u>EPA 6020B</u> | |

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|-----------------------|--------------------|---------------------|-------------------------|-----------------|---------------------|
| Antimony | 20.0 | 19.5 | 98 | | ug/L |
| Copper | 10.0 | 9.97 | 100 | | ug/L |
| Lead | 10.0 | 9.44 | 94 | | ug/L |
| Zinc | 200 | 198 | 99 | | ug/L |

CONTROL LIMITS 90-110%

III
CONTINUING CALIBRATION BLANK

Report No: 218081813 Blank ID: 1900
Instrument ID: ICPMS1 Lab File ID: 2180823A_MS1.b\037_CCB.d
Analyst: LWZ Analytical Batch: 642536
Analysis Date: 08/23/18 Time: 1205 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>RESULT</i> | <i>UNITS</i> | <i>Q</i> | <i>DL</i> | <i>LOD</i> | <i>LOQ</i> |
|----------------|---------------|--------------|----------|-----------|------------|------------|
| Antimony | 1.00 | ug/L | U | 0.50 | 1.00 | 2.00 |
| Copper | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Lead | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Zinc | 10.0 | ug/L | U | 5.00 | 10.0 | 20.0 |

FORM III - IN

II
CONTINUING CALIBRATION VERIFICATION (CCV) STANDARD

Report No: 218081813 GCAL QC ID: 1800
Instrument ID: ICPMS1 Lab File ID: 2180823A_MS1.b\056_CCV.d
Analyst: LWZ Analytical Batch: 642536
Analysis Date: 08/23/18 Time: 1328 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|----------------|-------------|--------------|------------------|----------|--------------|
| Antimony | 20.0 | 19.8 | 99 | | ug/L |
| Copper | 10.0 | 10.1 | 101 | | ug/L |
| Lead | 10.0 | 9.43 | 94 | | ug/L |
| Zinc | 200 | 199 | 100 | | ug/L |

CONTROL LIMITS 90-110%

FORM II - IN

III
CONTINUING CALIBRATION BLANK

Report No: 218081813 Blank ID: 1900
Instrument ID: ICPMS1 Lab File ID: 2180823A_MS1.b\057_CCB.d
Analyst: LWZ Analytical Batch: 642536
Analysis Date: 08/23/18 Time: 1332 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>RESULT</i> | <i>UNITS</i> | <i>Q</i> | <i>DL</i> | <i>LOD</i> | <i>LOQ</i> |
|----------------|---------------|--------------|----------|-----------|------------|------------|
| Antimony | 1.00 | ug/L | U | 0.50 | 1.00 | 2.00 |
| Copper | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Lead | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Zinc | 10.0 | ug/L | U | 5.00 | 10.0 | 20.0 |

FORM III - IN

XV (He)
ICP-MS INTERNAL STANDARDS RELATIVE INTENSITY SUMMARY

| | |
|------------------------------------|---------------------------------|
| Report No: <u>218081813</u> | Start Date: <u>08/23/18</u> |
| Instrument ID: <u>ICPMS1</u> | End Date: <u>08/23/18</u> |
| Analytical Method <u>EPA 6020B</u> | Analytical Batch: <u>642536</u> |

Internal Standards %RI For:

| CLIENT SAMPLE ID | GCAL SAMPLE ID | TIME | ISTD1 Q | ISTD2 Q | ISTD3 Q | ISTD4 Q | ISTD5 Q | ISTD6 Q | ISTD7 Q |
|------------------|----------------------|------|---------|---------|---------|---------|---------|---------|---------|
| MB1841866 | 1841866 | 1129 | 102 | 108 | 109 | 112 | 105 | 110 | 107 |
| LCS1841867 | 1841867 | 1134 | 99 | 105 | 106 | 107 | 101 | 109 | 105 |
| WIL01IS02 | 21808181302 | 1209 | 90 | 97 | 93 | 96 | 91 | 102 | 97 |
| WIL01IS02MS | 1842047 | 1214 | 84 | 95 | 92 | 95 | 88 | 100 | 95 |
| WIL01IS02MSD | 1842048 | 1218 | 87 | 96 | 92 | 95 | 88 | 101 | 95 |
| WIL01IS02PDS | 1842322 | 1222 | 86 | 97 | 92 | 95 | 89 | 101 | 96 |
| WIL01IS02SD | 1842323 | 1227 | 100 | 101 | 100 | 100 | 99 | 101 | 101 |
| WIL04IS03 | 21808181304 | 1231 | 93 | 99 | 95 | 97 | 91 | 101 | 98 |
| WIL04IS01 | 21808181305 | 1235 | 90 | 98 | 95 | 98 | 92 | 101 | 98 |
| WIL04IS02 | 21808181306 | 1240 | 93 | 98 | 95 | 98 | 94 | 102 | 98 |

| | | |
|------------------------|-----------------------|----------------------|
| ISTD 1: Bismuth (He) | ISTD 4: Lutetium (He) | ISTD 7: Terbium (He) |
| ISTD 2: Germanium (He) | ISTD 5: Rhodium (He) | |
| ISTD 3: Indium (He) | ISTD 6: Scandium (He) | |

FORM XV - IN

XV (No Gas)
ICP-MS INTERNAL STANDARDS RELATIVE INTENSITY SUMMARY

Report No: 218081813
Instrument ID: ICPMS1
Analytical Method EPA 6020B

Start Date: 08/23/18
End Date: 08/23/18
Analytical Batch: 642536

Internal Standards %RI For:

| CLIENT SAMPLE ID | GCAL SAMPLE ID | TIME | Internal Standards %RI For: | | | | | | | |
|------------------|----------------------|------|-----------------------------|---------|----------|----------|----------|----------|----------|--|
| | | | ISTD8 Q | ISTD9 Q | ISTD10 Q | ISTD11 Q | ISTD12 Q | ISTD13 Q | ISTD14 Q | |
| MB1841866 | 1841866 | 1129 | 98 | 103 | 103 | 101 | 102 | 106 | 102 | |
| LCS1841867 | 1841867 | 1134 | 96 | 103 | 101 | 102 | 101 | 105 | 103 | |
| WIL01IS02 | 21808181302 | 1209 | 96 | 105 | 101 | 103 | 100 | 109 | 104 | |
| WIL01IS02MS | 1842047 | 1214 | 95 | 103 | 100 | 102 | 98 | 108 | 102 | |
| WIL01IS02MSD | 1842048 | 1218 | 95 | 102 | 97 | 100 | 96 | 107 | 101 | |
| WIL01IS02PDS | 1842322 | 1222 | 95 | 104 | 99 | 101 | 98 | 108 | 102 | |
| WIL01IS02SD | 1842323 | 1227 | 101 | 103 | 103 | 103 | 103 | 104 | 104 | |
| WIL04IS03 | 21808181304 | 1231 | 97 | 104 | 101 | 102 | 100 | 108 | 104 | |
| WIL04IS01 | 21808181305 | 1235 | 97 | 105 | 101 | 103 | 100 | 108 | 104 | |
| WIL04IS02 | 21808181306 | 1240 | 98 | 105 | 102 | 104 | 102 | 108 | 105 | |

| | | |
|----------------------------|----------------------------|---------------------------|
| ISTD 8: Bismuth (No Gas) | ISTD 11: Lutetium (No Gas) | ISTD 14: Terbium (No Gas) |
| ISTD 9: Germanium (No Gas) | ISTD 12: Rhodium (No Gas) | |
| ISTD 10: Indium (No Gas) | ISTD 13: Scandium (No Gas) | |

FORM XV - IN

XIV
ANALYSIS RUN LOG

Report No: 218081813 Analytical Batch: 642829 Start Date: 08/28/18
 Instrument ID: ICPMS2 Analytical Method: EPA 6020B End Date: 08/28/18

| CLIENT SAMPLE ID | GCAL SAMPLE ID | PF | D/F | TIME | Analyte Symbols | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|------------------|----------------|----|-----|------|-----------------|----|----|----|----|---|----|----|----|----|----|----|----|----|----|----|----|----|----|---|----|----|----|----|----|----|----|----|---|----|
| | | | | | Al | Sb | As | Ba | Be | B | Cd | Ca | Cr | Co | Cu | Fe | Pb | Li | Mg | Mn | Hg | Mo | Ni | K | Se | Si | Ag | Na | Sr | Tl | Sn | Ti | V | Zn |
| ITUNE | 1150 | * | 1 | 0931 | | X | | | | | | | | X | X | | | | | | | | | | | | | | | | | | X | X |
| ICALB | 1300 | * | 1 | 1024 | X | | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X |
| ICAL2 | 1302 | * | 1 | 1028 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | | X | X |
| ICAL4 | 1304 | * | 1 | 1031 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | | X | X |
| ICAL5 | 1305 | * | 1 | 1035 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | | X | X |
| ICAL6 | 1306 | * | 1 | 1039 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | | X | X |
| ICV | 1600 | * | 1 | 1042 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| ICB | 1700 | * | 1 | 1050 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| LLCCV | 1803 | * | 1 | 1100 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| ICSA | 2000 | * | 1 | 1104 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| ICSAB | 2100 | * | 1 | 1107 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| LDR | 2500 | * | 1 | 1111 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| CCV | 1800 | * | 1 | 1240 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| CCB | 1900 | * | 1 | 1243 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| MB1842310 | 1842310 | * | 1 | 1247 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| LCS1842311 | 1842311 | * | 1 | 1250 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| WIL03IS01 | 21808181301 | * | 10 | 1305 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| WIL03IS02 | 21808181308 | * | 10 | 1308 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | | X | X |
| CCV | 1800 | * | 1 | 1337 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| CCB | 1900 | * | 1 | 1340 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| WIL03IS03 | 21808181307 | * | 10 | 1355 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| CCV | 1800 | * | 1 | 1358 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| CCB | 1900 | * | 1 | 1402 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| ICSA | 2000 | * | 1 | 1800 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |
| ICSAB | 2100 | * | 1 | 1804 | X | | | | | | | X | X | | | | | | | | | | | | | | | | | | | X | X | |

XIV
ANALYSIS RUN LOG

Report No: 218081814
Instrument ID: ICPMS2

Analytical Batch: 642829
Analytical Method: EPA 6020B

Start Date: 08/28/18
End Date: 08/28/18

| CLIENT SAMPLE ID | GCAL SAMPLE ID | PF | D/F | TIME | Analyte Symbols | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|------------------|----------------|----|-----|------|-----------------|----|----|----|----|---|----|----|----|----|----|----|----|----|----|----|----|----|----|---|----|----|----|----|----|----|----|----|---|----|----|---|---|---|---|--|
| | | | | | Al | Sb | As | Ba | Be | B | Cd | Ca | Cr | Co | Cu | Fe | Pb | Li | Mg | Mn | Hg | Mo | Ni | K | Se | Si | Ag | Na | Sr | Tl | Sn | Ti | V | Zn | Zr | | | | | |
| ITUNE | 1150 | * | 1 | 0931 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | | |
| IICALB | 1300 | * | 1 | 1024 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | |
| IICAL2 | 1302 | * | 1 | 1028 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | |
| IICAL4 | 1304 | * | 1 | 1031 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | |
| IICAL5 | 1305 | * | 1 | 1035 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | |
| IICAL6 | 1306 | * | 1 | 1039 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | |
| ICV | 1600 | * | 1 | 1042 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | |
| ICB | 1700 | * | 1 | 1050 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | |
| LLCCV | 1803 | * | 1 | 1100 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | |
| ICSA | 2000 | * | 1 | 1104 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | | |
| ICSAB | 2100 | * | 1 | 1107 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | |
| LDR | 2500 | * | 1 | 1111 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | |
| CCV | 1800 | * | 1 | 1240 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | |
| CCB | 1900 | * | 1 | 1243 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | |
| MB1842310 | 1842310 | * | 1 | 1247 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | |
| LCS1842311 | 1842311 | * | 1 | 1250 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| WIL02IS01 | 21808181401 | * | 10 | 1312 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | | |
| WIL02IS01 MS | 21808181402 | * | 10 | 1315 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| WIL02IS01 MSD | 21808181403 | * | 10 | 1319 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| WIL02IS01PDS | 1843727 | * | 10 | 1322 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| WIL02IS01SD | 1843728 | * | 50 | 1326 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| WIL02IS02 | 21808181404 | * | 10 | 1329 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| WIL01IS03 | 21808181405 | * | 10 | 1333 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| CCV | 1800 | * | 1 | 1337 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| CCB | 1900 | * | 1 | 1340 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| WIL01IS01 | 21808181406 | * | 10 | 1344 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| WIL02IS03 | 21808181407 | * | 10 | 1347 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| CCV | 1800 | * | 1 | 1358 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| CCB | 1900 | * | 1 | 1402 | X | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |

FORM XIV - IN

VIII
ICP-MS TUNE

Report No: 218081813 GCAL QC ID: 1150
Instrument ID: ICPMS2 Lab File ID: 2180828A_MS2.b\QCTune\2180828A_MS2-QCTu
Analyst: AWG Analytical Batch: 642829
Analysis Date: 08/28/18 Time: 0932 Analytical Method: EPA 6020B

| <i>ELEMENT - MASS</i> | <i>AVG MEASURED MASS (amu)</i> | <i>PEAK WIDTH AT 5% PEAK HEIGHT (amu)</i> | <i>%RSD</i> |
|-----------------------|--|---|-------------|
| Be-9 | 9 | .785 | .4858 |
| Mg-24 | 24 | .7864 | .6731 |
| Mg-25 | 25 | .7931 | 1.2316 |
| Mg-26 | 26 | .7857 | .6112 |
| Co-59 | 59 | .7733 | .4029 |
| In-115 | 115.05 | .7179 | .2559 |
| Pb-206 | 206.05 | .7639 | 1.0553 |
| Pb-207 | 207.05 | .744 | .4899 |
| Pb-208 | 208.05 | .7747 | .8738 |

FORM VIII - IN

II
INITIAL CALIBRATION VERIFICATION (ICV) STANDARD

| | |
|--|---|
| Report No: <u>218081813</u> | GCAL QC ID: <u>1600</u> |
| Instrument ID: <u>ICPMS2</u> | Lab File ID: <u>2180828B_MS2.b\015_ICV_2180828A_MS2.D</u> |
| Analyst: <u>LWZ</u> | Analytical Batch: <u>642829</u> |
| Analysis Date: <u>08/28/18</u> Time: <u>1042</u> | Analytical Method: <u>EPA 6020B</u> |

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|-----------------------|--------------------|---------------------|-------------------------|-----------------|---------------------|
| Antimony | 50.0 | 50.9 | 102 | | ug/L |
| Copper | 50.0 | 51.6 | 103 | | ug/L |
| Lead | 50.0 | 49.6 | 99 | | ug/L |
| Zinc | 1000 | 1010 | 101 | | ug/L |

CONTROL LIMITS 90-110%

FORM II - IN

III
INITIAL CALIBRATION BLANK

Report No: 218081813 Blank ID: 1700
Instrument ID: ICPMS2 Lab File ID: 2180828B_MS2.b\017_ICB_2180828A_MS2.D
Analyst: LWZ Analytical Batch: 642829
Analysis Date: 08/28/18 Time: 1050 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>RESULT</i> | <i>UNITS</i> | <i>Q</i> | <i>DL</i> | <i>LOD</i> | <i>LOQ</i> |
|----------------|---------------|--------------|----------|-----------|------------|------------|
| Antimony | 1.00 | ug/L | U | 0.50 | 1.00 | 2.00 |
| Copper | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Lead | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Zinc | 10.0 | ug/L | U | 5.00 | 10.0 | 20.0 |

FORM III - IN

II
 LOW LEVEL CONTINUING CALIBRATION VERIFICATION (LLCCV) STANDARD

| | |
|--|--|
| Report No: <u>218081813</u> | GCAL QC ID: <u>1803</u> |
| Instrument ID: <u>ICPMS2</u> | Lab File ID: <u>2180828B_MS2.b\1811CCV1_2180828A_MS2.D</u> |
| Analyst: <u>LWZ</u> | Analytical Batch: <u>642829</u> |
| Analysis Date: <u>08/28/18</u> Time: <u>1100</u> | Analytical Method: <u>EPA 6020B</u> |

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|----------------|-------------|--------------|------------------|----------|--------------|
| Antimony | 2.00 | 1.91 | 95 | | ug/L |
| Copper | 1.00 | 1.08 | 108 | | ug/L |
| Lead | 1.00 | 0.980 | 98 | | ug/L |
| Zinc | 20.0 | 19.1 | 95 | | ug/L |

CONTROL LIMITS 80-120%

FORM II - IN

IV
ICPMS INTERFERENCE CHECKS

| | |
|---|--|
| Report No: <u>218081813</u> | ICSA \ AB ID: <u>2000 \ 2100</u> |
| Instrument ID: <u>ICPMS2</u> | Analytical Batch: <u>642829</u> |
| Analyst: <u>LWZ</u> | Analytical Method: <u>EPA 6020B</u> |
| Lab File ID ICSA1: <u>2180828B_MS2.b\1812\ICSA_2180828A_MS2.D</u> | Lab File ID ICSAB1: <u>2180828B_MS2.b\1813\ICSB_2180828A_MS2.D</u> |
| Lab File ID ICSA2: <u>2180828C_MS2.b\070\ICSA.d</u> | Lab File ID ICSAB2: <u>2180828C_MS2.b\071\ICSB.d</u> |

Concentration Units: ug/L

| Analyzed (A/AB): | | | 08/28/18 1104 | 08/28/18 1107 | | 08/28/18 1800 | 08/28/18 1804 | |
|------------------|--------|---------|---------------|---------------|-----|---------------|---------------|------|
| ANALYTE | TRUE A | TRUE AB | ICSA1 | ICSAB1 | %R | ICSA2 | ICSAB2 | %R |
| Aluminum | 1000 | 1000 | 1020 | 1030 | 103 | 988 | 988 | 99 |
| Antimony | 0 | 0 | -0.059 | -0.057 | | -0.062 | -0.069 | |
| Arsenic | 0 | 10.0 | -0.0030 | 10.3 | 103 | -0.018 | 10.1 | 101 |
| Barium | 0 | 0 | -0.0040 | 0.014 | | 0.0060 | 0.021 | |
| Beryllium | 0 | 0 | -0.0070 | -0.0060 | | -0.0080 | -0.010 | |
| Boron | 0 | 20.0 | -0.11 | 20.1 | 100 | -0.71 | 18.5 | 92 |
| Cadmium | 0 | 10.0 | 0.014 | 9.83 | 98 | 0.0080 | 9.90 | 99 |
| Calcium | 3000 | 3000 | 2950 | 3030 | 101 | 2980 | 3010 | 100 |
| Chromium | 0 | 20.0 | 0.0010 | 20.4 | 102 | -0.11 | 19.4 | 97 |
| Cobalt | 0 | 20.0 | 0.0070 | 20.8 | 104 | 0.0050 | 20.2 | 101 |
| Copper | 0 | 20.0 | 0.027 | 21.7 | 108 | 0.098 | 21.1 | 106 |
| Iron | 2500 | 2500 | 2550 | 2550 | 102 | 2430 | 2450 | 98 |
| Lead | 0 | 0 | -0.0050 | -0.0030 | | -0.0050 | -0.0070 | |
| Lithium | 0 | 20.0 | 0.11 | 23.9 | 120 | -0.24 | 22.7 | 114 |
| Magnesium | 1000 | 1000 | 1020 | 1040 | 104 | 981 | 980 | 98 |
| Manganese | 0 | 20.0 | -0.010 | 20.4 | 102 | -0.028 | 19.7 | 98 |
| Molybdenum | 20.0 | 20.0 | 19.4 | 19.1 | 96 | 19.7 | 19.5 | 98 |
| Nickel | 0 | 20.0 | -0.0020 | 21.0 | 105 | -0.030 | 20.3 | 102 |
| Potassium | 1000 | 1000 | 991 | 1020 | 102 | 992 | 985 | 98 |
| Selenium | 0 | 10.0 | -0.020 | 9.83 | 98 | -0.0070 | 9.48 | 95 |
| Silicon | 0 | 1000 | -1200 | -80 | -8 | -5500 | -4500 | -454 |
| Silver | 0 | 5.00 | 0.0040 | 5.00 | 100 | 0.0030 | 5.06 | 101 |
| Sodium | 2500 | 2500 | 2520 | 2590 | 104 | 2450 | 2470 | 99 |
| Strontium | 0 | 10.0 | 0.10 | 10.3 | 103 | 0.11 | 10.4 | 104 |
| Thallium | 0 | 0 | -0.023 | -0.025 | | -0.043 | -0.041 | |
| Tin | 0 | 10.0 | -0.028 | 9.84 | 98 | 0.0060 | 10.1 | 101 |
| Titanium | 20.0 | 20.0 | 19.4 | 20.3 | 102 | 19.7 | 20.0 | 100 |
| Vanadium | 0 | 20.0 | -0.020 | 17.5 | 88 | -0.026 | 16.9 | 84 |
| Zinc | 0 | 20.0 | -0.92 | 19.5 | 98 | -1.0 | 18.8 | 94 |
| Zirconium | 0 | 20.0 | 0.023 | 19.4 | 97 | 0.0060 | 19.3 | 96 |

FORM IV - IN

II
 LINEAR DYNAMIC RANGE (LDR) STANDARD

| | | |
|--|--|--|
| Report No: <u>218081813</u> | GCAL QC ID: <u>2500</u> | |
| Instrument ID: <u>ICPMS2</u> | Lab File ID: <u>2180828B_MS2.b\1814_QC1_2180828A_MS2.D</u> | |
| Analyst: <u>LWZ</u> | Analytical Batch: <u>642829</u> | |
| Analysis Date: <u>08/28/18</u> Time: <u>1111</u> | Analytical Method: <u>EPA 6020B</u> | |

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|-----------------------|--------------------|---------------------|-------------------------|-----------------|---------------------|
| Antimony | 1000 | 955 | 96 | | ug/L |
| Copper | 1000 | 950 | 95 | | ug/L |
| Lead | 1000 | 972 | 97 | | ug/L |
| Zinc | 20000 | 18900 | 94 | | ug/L |

CONTROL LIMITS 90-110%

II
CONTINUING CALIBRATION VERIFICATION (CCV) STANDARD

| | |
|--|--|
| Report No: <u>218081813</u> | GCAL QC ID: <u>1800</u> |
| Instrument ID: <u>ICPMS2</u> | Lab File ID: <u>2180828B_MS2.b\1834_CCV_2180828A_MS2.D</u> |
| Analyst: <u>LWZ</u> | Analytical Batch: <u>642829</u> |
| Analysis Date: <u>08/28/18</u> Time: <u>1240</u> | Analytical Method: <u>EPA 6020B</u> |

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|-----------------------|--------------------|---------------------|-------------------------|-----------------|---------------------|
| Antimony | 20.0 | 20.1 | 100 | | ug/L |
| Copper | 10.0 | 10.7 | 107 | | ug/L |
| Lead | 10.0 | 9.83 | 98 | | ug/L |
| Zinc | 200 | 205 | 102 | | ug/L |

CONTROL LIMITS 90-110%

III
CONTINUING CALIBRATION BLANK

Report No: 218081813 Blank ID: 1900
Instrument ID: ICPMS2 Lab File ID: 2180828B_MS2.b\1835_CCB_2180828A_MS2.D
Analyst: LWZ Analytical Batch: 642829
Analysis Date: 08/28/18 Time: 1243 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>RESULT</i> | <i>UNITS</i> | <i>Q</i> | <i>DL</i> | <i>LOD</i> | <i>LOQ</i> |
|----------------|---------------|--------------|----------|-----------|------------|------------|
| Antimony | 1.00 | ug/L | U | 0.50 | 1.00 | 2.00 |
| Copper | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Lead | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Zinc | 10.0 | ug/L | U | 5.00 | 10.0 | 20.0 |

FORM III - IN

II
CONTINUING CALIBRATION VERIFICATION (CCV) STANDARD

| | | |
|--|--|--|
| Report No: <u>218081813</u> | GCAL QC ID: <u>1800</u> | |
| Instrument ID: <u>ICPMS2</u> | Lab File ID: <u>2180828B_MS2.b\1850_CCV_2180828A_MS2.D</u> | |
| Analyst: <u>LWZ</u> | Analytical Batch: <u>642829</u> | |
| Analysis Date: <u>08/28/18</u> Time: <u>1337</u> | Analytical Method: <u>EPA 6020B</u> | |

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|-----------------------|--------------------|---------------------|-------------------------|-----------------|---------------------|
| Antimony | 20.0 | 19.7 | 99 | | ug/L |
| Copper | 10.0 | 10.6 | 106 | | ug/L |
| Lead | 10.0 | 9.67 | 97 | | ug/L |
| Zinc | 200 | 203 | 102 | | ug/L |

CONTROL LIMITS 90-110%

III
CONTINUING CALIBRATION BLANK

Report No: 218081813 Blank ID: 1900
Instrument ID: ICPMS2 Lab File ID: 2180828B_MS2.b\1851_CCB_2180828A_MS2.D
Analyst: LWZ Analytical Batch: 642829
Analysis Date: 08/28/18 Time: 1340 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>RESULT</i> | <i>UNITS</i> | <i>Q</i> | <i>DL</i> | <i>LOD</i> | <i>LOQ</i> |
|----------------|---------------|--------------|----------|-----------|------------|------------|
| Antimony | 1.00 | ug/L | U | 0.50 | 1.00 | 2.00 |
| Copper | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Lead | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Zinc | 10.0 | ug/L | U | 5.00 | 10.0 | 20.0 |

FORM III - IN

II
CONTINUING CALIBRATION VERIFICATION (CCV) STANDARD

| | | |
|--|--|--|
| Report No: <u>218081813</u> | GCAL QC ID: <u>1800</u> | |
| Instrument ID: <u>ICPMS2</u> | Lab File ID: <u>2180828C_MS2.b\002_CCV.d</u> | |
| Analyst: <u>LWZ</u> | Analytical Batch: <u>642829</u> | |
| Analysis Date: <u>08/28/18</u> Time: <u>1358</u> | Analytical Method: <u>EPA 6020B</u> | |

| <i>ANALYTE</i> | <i>TRUE</i> | <i>FOUND</i> | <i>%RECOVERY</i> | <i>Q</i> | <i>UNITS</i> |
|-----------------------|--------------------|---------------------|-------------------------|-----------------|---------------------|
| Antimony | 20.0 | 20.2 | 101 | | ug/L |
| Copper | 10.0 | 10.5 | 105 | | ug/L |
| Lead | 10.0 | 9.64 | 96 | | ug/L |
| Zinc | 200 | 203 | 102 | | ug/L |

CONTROL LIMITS 90-110%

FORM II - IN

III
CONTINUING CALIBRATION BLANK

Report No: 218081813 Blank ID: 1900
Instrument ID: ICPMS2 Lab File ID: 2180828C_MS2.b\003_CCB.d
Analyst: LWZ Analytical Batch: 642829
Analysis Date: 08/28/18 Time: 1402 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>RESULT</i> | <i>UNITS</i> | <i>Q</i> | <i>DL</i> | <i>LOD</i> | <i>LOQ</i> |
|----------------|---------------|--------------|----------|-----------|------------|------------|
| Antimony | 1.00 | ug/L | U | 0.50 | 1.00 | 2.00 |
| Copper | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Lead | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Zinc | 10.0 | ug/L | U | 5.00 | 10.0 | 20.0 |

FORM III - IN

XV (He)
ICP-MS INTERNAL STANDARDS RELATIVE INTENSITY SUMMARY

| | |
|------------------------------------|---------------------------------|
| Report No: <u>218081813</u> | Start Date: <u>08/28/18</u> |
| Instrument ID: <u>ICPMS2</u> | End Date: <u>08/28/18</u> |
| Analytical Method <u>EPA 6020B</u> | Analytical Batch: <u>642829</u> |

| CLIENT SAMPLE ID | GCAL SAMPLE ID | TIME | Internal Standards %RI For: | | | | | | | | | |
|------------------|----------------------|------|-----------------------------|---------|---------|---------|---------|---------|---------|--|--|--|
| | | | ISTD1 Q | ISTD2 Q | ISTD3 Q | ISTD4 Q | ISTD5 Q | ISTD6 Q | ISTD7 Q | | | |
| MB1842310 | 1842310 | 1247 | 126 | 115 | 121 | 122 | 114 | 113 | 119 | | | |
| LCS1842311 | 1842311 | 1250 | 130 | 119 | 124 | 128 | 115 | 115 | 125 | | | |
| WIL03IS01 | 21808181301 | 1305 | 87 | 100 | 94 | 77 | 93 | 101 | 83 | | | |
| WIL03IS02 | 21808181308 | 1308 | 91 | 101 | 95 | 79 | 96 | 100 | 85 | | | |
| WIL03IS03 | 21808181307 | 1355 | 94 | 102 | 99 | 83 | 99 | 101 | 90 | | | |

| | | |
|------------------------|-----------------------|----------------------|
| ISTD 1: Bismuth (He) | ISTD 4: Lutetium (He) | ISTD 7: Terbium (He) |
| ISTD 2: Germanium (He) | ISTD 5: Rhodium (He) | |
| ISTD 3: Indium (He) | ISTD 6: Scandium (He) | |

FORM XV - IN

XV (No Gas)
ICP-MS INTERNAL STANDARDS RELATIVE INTENSITY SUMMARY

Report No: 218081813
Instrument ID: ICPMS2
Analytical Method EPA 6020B

Start Date: 08/28/18
End Date: 08/28/18
Analytical Batch: 642829

Internal Standards %RI For:

| CLIENT SAMPLE ID | GCAL SAMPLE ID | TIME | ISTD8 Q | ISTD9 Q | ISTD10 Q | ISTD11 Q | ISTD12 Q | ISTD13 Q | ISTD14 Q |
|------------------|----------------------|------|---------|---------|----------|----------|----------|----------|----------|
| MB1842310 | 1842310 | 1247 | 110 | 103 | 106 | 108 | 104 | 98 | 107 |
| LCS1842311 | 1842311 | 1250 | 113 | 104 | 108 | 112 | 105 | 99 | 110 |
| WIL03IS01 | 21808181301 | 1305 | 87 | 95 | 89 | 84 | 89 | 97 | 86 |
| WIL03IS02 | 21808181308 | 1308 | 90 | 96 | 92 | 86 | 91 | 97 | 88 |
| WIL03IS03 | 21808181307 | 1355 | 92 | 98 | 94 | 92 | 94 | 98 | 94 |

| | | |
|----------------------------|---------------------------|--------------------------|
| ISTD 8: Bismuth (No Gas) | ISTD 11 Lutetium (No Gas) | ISTD 14 Terbium (No Gas) |
| ISTD 9: Germanium (No Gas) | ISTD 12 Rhodium (No Gas) | |
| ISTD 10 Indium (No Gas) | ISTD 13 Scandium (No Gas) | |

FORM XV - IN

XV (He)
ICP-MS INTERNAL STANDARDS RELATIVE INTENSITY SUMMARY

Report No: 218081814
Instrument ID: ICPMS2
Analytical Method EPA 6020B

Start Date: 08/28/18
End Date: 08/28/18
Analytical Batch: 642829

Internal Standards %RI For:

| CLIENT SAMPLE ID | GCAL SAMPLE ID | TIME | Internal Standards %RI For: | | | | | | | | | |
|------------------|----------------------|------|-----------------------------|---------|---------|---------|---------|---------|---------|--|--|--|
| | | | ISTD1 Q | ISTD2 Q | ISTD3 Q | ISTD4 Q | ISTD5 Q | ISTD6 Q | ISTD7 Q | | | |
| MB1842310 | 1842310 | 1247 | 126 | 115 | 121 | 122 | 114 | 113 | 119 | | | |
| LCS1842311 | 1842311 | 1250 | 130 | 119 | 124 | 128 | 115 | 115 | 125 | | | |
| WIL02IS01 | 21808181401 | 1312 | 88 | 101 | 95 | 78 | 95 | 102 | 84 | | | |
| WIL02IS01 MS | 21808181402 | 1315 | 85 | 100 | 94 | 75 | 95 | 102 | 79 | | | |
| WIL02IS01 MSD | 21808181403 | 1319 | 87 | 101 | 94 | 76 | 94 | 102 | 81 | | | |
| WIL02IS01PDS | 1843727 | 1322 | 90 | 101 | 97 | 79 | 95 | 102 | 84 | | | |
| WIL02IS01SD | 1843728 | 1326 | 92 | 99 | 97 | 79 | 97 | 96 | 85 | | | |
| WIL02IS02 | 21808181404 | 1329 | 92 | 101 | 96 | 80 | 96 | 100 | 88 | | | |
| WIL01IS03 | 21808181405 | 1333 | 89 | 99 | 94 | 77 | 96 | 99 | 83 | | | |
| WIL01IS01 | 21808181406 | 1344 | 89 | 101 | 97 | 78 | 97 | 100 | 85 | | | |
| WIL02IS03 | 21808181407 | 1347 | 94 | 102 | 98 | 83 | 98 | 102 | 90 | | | |

ISTD 1: Bismuth (He) ISTD 4: Lutetium (He) ISTD 7: Terbium (He)
 ISTD 2: Germanium (He) ISTD 5: Rhodium (He)
 ISTD 3: Indium (He) ISTD 6: Scandium (He)

FORM XV - IN

XV (No Gas)
ICP-MS INTERNAL STANDARDS RELATIVE INTENSITY SUMMARY

Report No: 218081814
Instrument ID: ICPMS2
Analytical Method EPA 6020B

Start Date: 08/28/18
End Date: 08/28/18
Analytical Batch: 642829

Internal Standards %RI For:

| CLIENT SAMPLE ID | GCAL SAMPLE ID | TIME | Internal Standards %RI For: | | | | | | | | | |
|------------------|----------------------|------|-----------------------------|---------|----------|----------|----------|----------|----------|--|--|--|
| | | | ISTD8 Q | ISTD9 Q | ISTD10 Q | ISTD11 Q | ISTD12 Q | ISTD13 Q | ISTD14 Q | | | |
| MB1842310 | 1842310 | 1247 | 110 | 103 | 106 | 108 | 104 | 98 | 107 | | | |
| LCS1842311 | 1842311 | 1250 | 113 | 104 | 108 | 112 | 105 | 99 | 110 | | | |
| WIL02IS01 | 21808181401 | 1312 | 88 | 96 | 92 | 87 | 91 | 99 | 89 | | | |
| WIL02IS01 MS | 21808181402 | 1315 | 85 | 93 | 89 | 82 | 89 | 96 | 84 | | | |
| WIL02IS01 MSD | 21808181403 | 1319 | 89 | 96 | 91 | 85 | 91 | 98 | 87 | | | |
| WIL02IS01PDS | 1843727 | 1322 | 90 | 96 | 93 | 88 | 91 | 98 | 89 | | | |
| WIL02IS01SD | 1843728 | 1326 | 92 | 96 | 94 | 89 | 94 | 95 | 90 | | | |
| WIL02IS02 | 21808181404 | 1329 | 92 | 96 | 93 | 89 | 93 | 98 | 91 | | | |
| WIL01IS03 | 21808181405 | 1333 | 89 | 97 | 94 | 88 | 92 | 97 | 90 | | | |
| WIL01IS01 | 21808181406 | 1344 | 89 | 96 | 93 | 87 | 92 | 98 | 89 | | | |
| WIL02IS03 | 21808181407 | 1347 | 90 | 98 | 95 | 90 | 94 | 99 | 92 | | | |

ISTD 8: Bismuth (No Gas) ISTD 11: Lutetium (No Gas) ISTD 14: Terbium (No Gas)
 ISTD 9: Germanium (No Gas) ISTD 12: Rhodium (No Gas)
 ISTD 10: Indium (No Gas) ISTD 13: Scandium (No Gas)

FORM XV - IN

XIII
PREPARATION LOG

Report No: 218081812

Prep Method: EPA 3050B

Prep Batch: 642222

| <i>CLIENT SAMPLE ID</i> | <i>GCAL SAMPLE ID</i> | <i>PREP DATE</i> | <i>WEIGHT</i> | <i>UNITS</i> | <i>VOLUME</i> | <i>UNITS</i> |
|-------------------------|-----------------------|------------------|---------------|--------------|---------------|--------------|
| LCS1840819 | 1840819 | 08/20/18 | 1.25 | g | 50 | mL |
| MB1840818 | 1840818 | 08/20/18 | 1.25 | g | 50 | mL |
| WIL02DA01A | 21808181209 | 08/20/18 | 1.45 | g | 50 | mL |
| WIL02DA01B | 21808181210 | 08/20/18 | 1.33 | g | 50 | mL |
| WIL02DA02A | 21808181206 | 08/20/18 | 1.25 | g | 50 | mL |
| WIL02DA02A MS | 21808181207 | 08/20/18 | 1.25 | g | 50 | mL |
| WIL02DA02A MSD | 21808181208 | 08/20/18 | 1.25 | g | 50 | mL |

FORM XIII - IN

III
METHOD BLANK

Report No: 218081812 Blank ID: MB1840818
Instrument ID: ICPMS1 Lab File ID: 2180821A_MS1.b\020SMPL.d
Analyst: LWZ Analytical Batch: 642381
Analysis Date: 08/21/18 Time: 1138 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>RESULT</i> | <i>UNITS</i> | <i>Q</i> | <i>DL</i> | <i>LOD</i> | <i>LOQ</i> |
|----------------|---------------|--------------|----------|-----------|------------|------------|
| Antimony | 21.7 | ug/kg | J | 20.0 | 40.0 | 80.0 |
| Copper | 13.0 | ug/kg | J | 10.0 | 20.0 | 40.0 |
| Lead | 20.0 | ug/kg | U | 10.0 | 20.0 | 40.0 |
| Zinc | 518 | ug/kg | J | 200 | 400 | 800 |

FORM III - IN

VII
LABORATORY CONTROL SAMPLE

| | | | |
|--------------|-----------------------------------|--------------------|-----------------------------------|
| Report No: | <u>218081812</u> | GCAL ID: | <u>1840819 (LCS)</u> |
| Matrix: | <u>Solid</u> | Instrument ID: | <u>ICPMS1</u> |
| Analyst: | <u>LWZ</u> | Lab File ID: | <u>2180821A_MS1.b\021SMPL.d</u> |
| Prep Date: | <u>08/20/18</u> Time: <u>1100</u> | Analysis Date: | <u>08/21/18</u> Time: <u>1143</u> |
| Prep Batch: | <u>642222</u> | Analytical Batch: | <u>642381</u> |
| Prep Method: | <u>3050B</u> | Analytical Method: | <u>EPA 6020B</u> |

| ANALYTE | UNITS | TRUE | FOUND | % R | Q | LCL | UCL |
|----------|-------|-------|-------|-----|----|-----|-----|
| Antimony | ug/kg | 4000 | 3800 | 95 | 72 | 124 | |
| Copper | ug/kg | 2000 | 1990 | 100 | 84 | 119 | |
| Lead | ug/kg | 2000 | 1970 | 99 | 84 | 118 | |
| Zinc | ug/kg | 40000 | 34700 | 87 | 82 | 119 | |

FORM VII - IN

V1
MS/MSD RECOVERY

Report No: 218081812 Parent Sample ID: WIL02DA02A
 Prep Method: 3050B Parent GCAL ID: 21808181206
 Prep Date: 08/20/18 Time: 1100 Prep Batch: 642222
 Analytical Method: EPA 6020B Analytical Batch: 642309

| | |
|-----------------------------------|--|
| GCAL QC ID: 21808181207 MS | Instrument ID: ICPMS2 |
| Analyst: LWZ | Lab File ID: 2180820B_MS2.b\121278SMPL.d |
| Analysis Date: 08/20/18 1806 | Dilution: 10 |

| ANALYTE | UNITS | SPIKE ADDED | SAMPLE RESULT | MS RESULT | MS % REC | # | QC LIMITS |
|----------|-------|-------------|---------------|-----------|----------|---|-----------|
| Antimony | ug/kg | 5740 | 0 | 373 | 7 | * | 72 - 124 |
| Copper | ug/kg | 2870 | 35400 | 36500 | 38 | * | 84 - 119 |
| Lead | ug/kg | 2870 | 18900 | 20800 | 67 | * | 84 - 118 |
| Zinc | ug/kg | 57400 | 101000 | 157000 | 98 | | 82 - 119 |

| | |
|------------------------------------|--|
| GCAL QC ID: 21808181208 MSD | Instrument ID: ICPMS2 |
| Analyst: LWZ | Lab File ID: 2180820B_MS2.b\121279SMPL.d |
| Analysis Date: 08/20/18 1810 | Dilution: 10 |

| ANALYTE | UNITS | SPIKE ADDED | MSD RESULT | MSD % REC | # | % RPD | # | QC LIMITS | RPD |
|----------|-------|-------------|------------|-----------|---|-------|---|-----------|--------|
| Antimony | ug/kg | 5740 | 314 | 5 | * | 17 | | 72 - 124 | 0 - 20 |
| Copper | ug/kg | 2870 | 42100 | 232 | * | 14 | | 84 - 119 | 0 - 20 |
| Lead | ug/kg | 2870 | 23000 | 141 | * | 10 | | 84 - 118 | 0 - 20 |
| Zinc | ug/kg | 57400 | 170000 | 121 | * | 8 | | 82 - 119 | 0 - 20 |

RPD : 0 out of 4 outside limits
 Spike Recovery: 7 out of 8 outside limits

Column to be used to flag recovery and RPD values with an asterisk
 * Values outside of QC limits

FORM V (PART 1) - IN

V2
POST DIGEST SPIKE SAMPLE RECOVERY

| | | | |
|--------------------|-----------------------------------|-------------------|------------------------------------|
| Report No: | <u>218081812</u> | GCAL PDS ID: | <u>1841231</u> |
| Matrix: | <u>Solid</u> | Parent Sample ID: | <u>WIL02DA02A (21808181206)</u> |
| Analyst: | <u>LWZ</u> | Instrument ID: | <u>ICPMS2</u> |
| Analysis Date: | <u>08/20/18</u> Time: <u>1813</u> | Lab File ID: | <u>2180820B_MS2.b\121280SMPL.d</u> |
| Analytical Method: | <u>EPA 6020B</u> | Analytical Batch: | <u>642309</u> |

| ANALYTE | UNITS | SPIKED SAMPLE RESULT | C | SAMPLE RESULT | C | SPIKE ADDED | % R | Q | LCL | UCL |
|----------|-------|----------------------------|---|------------------|---|----------------|-----|---|-----|-----|
| Antimony | ug/kg | 54000 | | 0 | U | 57400 | 94 | | 80 | 120 |
| Copper | ug/kg | 64600 | | 35400 | | 28700 | 102 | | 80 | 120 |
| Lead | ug/kg | 48100 | | 18900 | | 28700 | 102 | | 80 | 120 |
| Zinc | ug/kg | 652000 | | 101000 | | 574000 | 96 | | 80 | 120 |

IX
SERIAL DILUTIONS

| | | | |
|--------------------|-----------------------------------|-------------------|------------------------------------|
| Report No: | <u>218081812</u> | GCAL SD ID: | <u>1841232</u> |
| Matrix: | <u>Solid</u> | Parent Sample ID: | <u>WIL02DA02A (21808181206)</u> |
| Analyst: | <u>LWZ</u> | Instrument ID: | <u>ICPMS2</u> |
| Analysis Date: | <u>08/20/18</u> Time: <u>1817</u> | Lab File ID: | <u>2180820B_MS2.b\121281SMPL.d</u> |
| Analytical Method: | <u>EPA 6020B</u> | Analytical Batch: | <u>642309</u> |

| ANALYTE | UNITS | PARENT | | SERIAL | | % DIFF | Q | LCL | UCL |
|----------|-------|--------|--------|----------|--------|--------|------|-----|-----|
| | | SAMPLE | RESULT | DILUTION | RESULT | | | | |
| Antimony | ug/kg | | 0 | U | 0 | U | | | |
| Copper | ug/kg | | 35400 | | 38200 | | 7.9 | 0 | 10 |
| Lead | ug/kg | | 18900 | | 19600 | | 3.7 | 0 | 10 |
| Zinc | ug/kg | | 101000 | | 112000 | | 10.9 | E | 0 |

FORM IX - IN



3010A Metals Water Preparation



| | | | | | | | |
|--------------|-----------|-----------------|--------------|---------------|-------|-------|--------|
| ANALYST/TECH | <i>JK</i> | START DATE/TIME | 11:30 9/7/08 | END DATE/TIME | 12:55 | BATCH | 642277 |
|--------------|-----------|-----------------|--------------|---------------|-------|-------|--------|

| # | CLIENT | TYPE | CLIENT ID | GCAL ID | INITIAL VOL (mL) | FINAL VOL (mL) | COMMENT | STANDARDS/ REAGENTS |
|----|--------|------|--------------------------|-------------|------------------|----------------|---------|------------------------|
| 1 | QC | MB | MB 1841052 | 1841052 | 80 | 90 | | GCAL - 8 - 250uL |
| 2 | QC | LCS | LCS 1841053 | 1841053 | ↓ | | | 7261918 |
| 3 | 4989 | SAMP | #14 | 21808166103 | 5mL | | | Sb,Ag,Se SPIKE - 250uL |
| 4 | 0042 | SAMP | Nickel Catalyst #2... | 21808173601 | 5mL | | | 316-652 |
| 5 | QC | MS | Nickel Catalyst #2...MS | 1841056 | | | | Li,B,Zr SPIKE - 250uL |
| 6 | QC | MSD | Nickel Catalyst #2...MSD | 1841057 | | | | 316 66-2 |
| 7 | 4629 | SAMP | PMW-3A | 21808180601 | 50mL | | | SI SPIKE - 250uL |
| 8 | 4629 | SAMP | MW-65 | 21808180602 | | | | 726775 |
| 9 | 4838 | SAMP | WIL03IS00 | 21808181303 | | | | 1:1 HNO3 |
| 10 | 0176 | SAMP | SARA Separator Eff... | 21808201101 | | | | 726 77-6 |
| 11 | 0176 | SAMP | SARA ACLA Sump | 21808201102 | | | | 1:1 HCL |
| 12 | 0176 | SAMP | U31PS201G | 21808201401 | ↓ | ↓ | | 726 78-1 |
| 13 | | | | | | | | |
| 14 | | | | | | | | |
| 15 | | | | | | | | |
| 16 | | | | | | | | |
| 17 | | | | | | | | |
| 18 | | | | | | | | |
| 19 | | | | | | | | |
| 20 | | | | | | | | |
| 21 | | | | | | | | |
| 22 | | | | | | | | |
| 23 | | | | | | | | |
| 24 | | | | | | | | |
| 25 | | | | | | | | |
| 26 | | | | | | | | |
| 27 | | | | | | | | Digestion Vessel Lot# |
| 28 | | | | | | | | 180516 |
| 29 | | | | | | | | |
| 30 | | | | | | | | |

EQUIPMENT/CONDITIONS

| | | | |
|-----------------|----|-------------|-----|
| DIGESTION BLOCK | A1 | TEMPERATURE | 94° |
|-----------------|----|-------------|-----|

NOTES

TCLPs are reduced in volume due to sample matrix. Matrix-Water. 6020_W_EX

XIII
PREPARATION LOG

Report No: 218081813

Prep Method: EPA 3010A

Prep Batch: 642277

| <i>CLIENT SAMPLE ID</i> | <i>GCAL SAMPLE ID</i> | <i>PREP DATE</i> | <i>WEIGHT</i> | <i>UNITS</i> | <i>VOLUME</i> | <i>UNITS</i> |
|-------------------------|-----------------------|------------------|---------------|--------------|---------------|--------------|
| LCS1841053 | 1841053 | 08/20/18 | | | 50 | mL |
| MB1841052 | 1841052 | 08/20/18 | | | 50 | mL |
| Nickel Catalyst ...MS | 1841056 | 08/20/18 | | | 50 | mL |
| Nickel Catalyst...MSD | 1841057 | 08/20/18 | | | 50 | mL |
| WIL03IS00 | 21808181303 | 08/20/18 | | | 50 | mL |

FORM XIII - IN

III
METHOD BLANK

Report No: 218081813 Blank ID: MB1841052
Instrument ID: ICPMS2 Lab File ID: 2180820B_MS2.b\121284SMPL.d
Analyst: LWZ Analytical Batch: 642309
Analysis Date: 08/20/18 Time: 1827 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>RESULT</i> | <i>UNITS</i> | <i>Q</i> | <i>DL</i> | <i>LOD</i> | <i>LOQ</i> |
|----------------|---------------|--------------|----------|-----------|------------|------------|
| Antimony | 1.00 | ug/L | U | 0.50 | 1.00 | 2.00 |
| Copper | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Lead | 0.50 | ug/L | U | 0.25 | 0.50 | 1.00 |
| Zinc | 10.0 | ug/L | U | 5.00 | 10.0 | 20.0 |

FORM III - IN

VII
LABORATORY CONTROL SAMPLE

| | |
|--|--|
| Report No: <u>218081813</u> | GCAL ID: <u>1841053 (LCS)</u> |
| Matrix: <u>Water</u> | Instrument ID: <u>ICPMS2</u> |
| Analyst: <u>LWZ</u> | Lab File ID: <u>2180820B_MS2.b\121285SMPL.d</u> |
| Prep Date: <u>08/20/18</u> Time: <u>1130</u> | Analysis Date: <u>08/20/18</u> Time: <u>1831</u> |
| Prep Batch: <u>642277</u> | Analytical Batch: <u>642309</u> |
| Prep Method: <u>3010A</u> | Analytical Method: <u>EPA 6020B</u> |

| ANALYTE | UNITS | TRUE | FOUND | % R | Q | LCL | UCL |
|----------|-------|------|-------|-----|---|-----|-----|
| Antimony | ug/L | 100 | 106 | 106 | | 85 | 117 |
| Copper | ug/L | 50.0 | 56.2 | 112 | | 85 | 118 |
| Lead | ug/L | 50.0 | 51.9 | 104 | | 88 | 115 |
| Zinc | ug/L | 1000 | 1020 | 102 | | 83 | 119 |

FORM VII - IN



3050B Metals Solid Preparation



| | | | |
|---------------------|--------------------------------|------------------------|-----------------|
| ANALYST/TECH JSL | START DATE/TIME 12:18 02-19 | END DATE/TIME 15:10 | BATCH 642442 |
|---------------------|--------------------------------|------------------------|-----------------|

| # | CLIENT | TYPE | CLIENT ID | GCAL ID | INITIAL WGT (g) | FINAL VOL (mL) | COMMENT | STANDARDS/ REAGENTS |
|----|--------|------|-----------------------|-------------|-----------------|----------------|---------|--------------------------|
| 1 | QC | MB | MB 1841866 | 1841866 | 1.25 | 50 | | GCAL - B - 250uL |
| 2 | QC | LCS | LCS 1841867 | 1841867 | 1.25 | ↓ | | 2176918 |
| 3 | 4838 | SAMP | WIL01IS02 | 21808181302 | 1.25 | | | Sb, Ag, Se SPIKE - 250uL |
| 4 | 4838 | SAMP | WIL04IS03 | 21808181304 | 1.33 | | | 311.69-2 |
| 5 | 4838 | SAMP | WIL04IS01 | 21808181305 | 1.35 | | | LI, B, Z SPIKE - 250uL |
| 6 | 4838 | SAMP | WIL04IS02 | 21808181306 | 1.32 | | | 311.66-2 |
| 7 | 4612 | SAMP | ADA-25X38-A03-SP07... | 21808210501 | 1.32 | | | SI SPIKE - 250uL |
| 8 | QC | MS | WIL01IS02 MS | 1842047 | 1.25 | | 50 | |
| 9 | QC | MSD | WIL01IS02 MSD | 1842048 | 1.25 | ↓ | | HNO3 |
| 10 | | | | | | | | 301 11-15 |
| 11 | | | | | | | | H2O2 |
| 12 | | | | | | | | 2176907 |
| 13 | | | | | | | | HCL |
| 14 | | | | | | | | N/A |
| 15 | | | | | | | | 1:1 HNO3 |
| 16 | | | | | | | | 217678-2 |
| 17 | | | | | | | | |
| 18 | | | | | | | | |
| 19 | | | | | | | | |
| 20 | | | | | | | | |
| 21 | | | | | | | | |
| 22 | | | | | | | | |
| 23 | | | | | | | | |
| 24 | | | | | | | | |
| 25 | | | | | | | | |
| 26 | | | | | | | | |
| 27 | | | | | | | | Digestion Vessel Lot # |
| 28 | | | | | | | | 188513 |
| 29 | | | | | | | | |
| 30 | | | | | | | | |

EQUIPMENT/CONDITIONS

| | |
|-----------------------|--------------------|
| DIGESTION BLOCK B2 | TEMPERATURE 93° |
|-----------------------|--------------------|

NOTES

Matrix-Soll. 6020_S_EX

XIII
PREPARATION LOG

Report No: 218081813

Prep Method: EPA 3050B

Prep Batch: 642442

| <i>CLIENT SAMPLE ID</i> | <i>GCAL SAMPLE ID</i> | <i>PREP DATE</i> | <i>WEIGHT</i> | <i>UNITS</i> | <i>VOLUME</i> | <i>UNITS</i> |
|-------------------------|-----------------------|------------------|---------------|--------------|---------------|--------------|
| LCS1841867 | 1841867 | 08/22/18 | 1.25 | g | 50 | mL |
| MB1841866 | 1841866 | 08/22/18 | 1.25 | g | 50 | mL |
| WIL01IS02 | 21808181302 | 08/22/18 | 1.25 | g | 50 | mL |
| WIL01IS02MS | 1842047 | 08/22/18 | 1.25 | g | 50 | mL |
| WIL01IS02MSD | 1842048 | 08/22/18 | 1.25 | g | 50 | mL |
| WIL04IS01 | 21808181305 | 08/22/18 | 1.35 | g | 50 | mL |
| WIL04IS02 | 21808181306 | 08/22/18 | 1.32 | g | 50 | mL |
| WIL04IS03 | 21808181304 | 08/22/18 | 1.33 | g | 50 | mL |

FORM XIII - IN

III
METHOD BLANK

Report No: 218081813 Blank ID: MB1841866
Instrument ID: ICPMS1 Lab File ID: 2180823A_MS1.b\029SMPL.d
Analyst: LWZ Analytical Batch: 642536
Analysis Date: 08/23/18 Time: 1129 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>RESULT</i> | <i>UNITS</i> | <i>Q</i> | <i>DL</i> | <i>LOD</i> | <i>LOQ</i> |
|----------------|---------------|--------------|----------|-----------|------------|------------|
| Antimony | 40.0 | ug/kg | U | 20.0 | 40.0 | 80.0 |
| Copper | 20.0 | ug/kg | U | 10.0 | 20.0 | 40.0 |
| Lead | 20.0 | ug/kg | U | 10.0 | 20.0 | 40.0 |
| Zinc | 400 | ug/kg | U | 200 | 400 | 800 |

FORM III - IN

VII
LABORATORY CONTROL SAMPLE

| | | | |
|--------------|-----------------------------------|--------------------|-----------------------------------|
| Report No: | <u>218081813</u> | GCAL ID: | <u>1841867 (LCS)</u> |
| Matrix: | <u>Solid</u> | Instrument ID: | <u>ICPMS1</u> |
| Analyst: | <u>LWZ</u> | Lab File ID: | <u>2180823A_MS1.b030SMPL.d</u> |
| Prep Date: | <u>08/22/18</u> Time: <u>1210</u> | Analysis Date: | <u>08/23/18</u> Time: <u>1134</u> |
| Prep Batch: | <u>642442</u> | Analytical Batch: | <u>642536</u> |
| Prep Method: | <u>3050B</u> | Analytical Method: | <u>EPA 6020B</u> |

| ANALYTE | UNITS | TRUE | FOUND | % R | Q | LCL | UCL |
|----------|-------|-------|-------|-----|---|-----|-----|
| Antimony | ug/kg | 4000 | 3820 | 96 | | 72 | 124 |
| Copper | ug/kg | 2000 | 1960 | 98 | | 84 | 119 |
| Lead | ug/kg | 2000 | 1990 | 99 | | 84 | 118 |
| Zinc | ug/kg | 40000 | 36700 | 92 | | 82 | 119 |

FORM VII - IN

V1
MS/MSD RECOVERY

Report No: 218081813 Parent Sample ID: WIL01IS02
 Prep Method: 3050B Parent GCAL ID: 21808181302
 Prep Date: 08/22/18 Time: 1210 Prep Batch: 642442
 Analytical Method: EPA 6020B Analytical Batch: 642536

GCAL QC ID: **1842047 MS** Instrument ID: ICPMS1
 Analyst: LWZ Lab File ID: 2180823A_MS1.b\039SMPL.d
 Analysis Date: 08/23/18 1214 Dilution: 10

| ANALYTE | UNITS | SPIKE ADDED | SAMPLE RESULT | MS RESULT | MS % REC | # | QC LIMITS |
|----------|-------|----------------|------------------|--------------|-------------|---|-----------|
| Antimony | ug/kg | 4270 | 0 | 471 | 11 | * | 72 - 124 |
| Copper | ug/kg | 2140 | 21100 | 23800 | 132 | * | 84 - 119 |
| Lead | ug/kg | 2140 | 63500 | 66800 | 152 | * | 84 - 118 |
| Zinc | ug/kg | 42700 | 61200 | 110000 | 114 | | 82 - 119 |

GCAL QC ID: **1842048 MSD** Instrument ID: ICPMS1
 Analyst: LWZ Lab File ID: 2180823A_MS1.b\040SMPL.d
 Analysis Date: 08/23/18 1218 Dilution: 10

| ANALYTE | UNITS | SPIKE ADDED | MSD RESULT | MSD % REC | # | % RPD | # | QC LIMITS %REC | RPD |
|----------|-------|----------------|---------------|--------------|---|----------|---|-------------------|--------|
| Antimony | ug/kg | 4270 | 496 | 12 | * | 5 | | 72 - 124 | 0 - 20 |
| Copper | ug/kg | 2140 | 23500 | 115 | | 2 | | 84 - 119 | 0 - 20 |
| Lead | ug/kg | 2140 | 61100 | 0 | * | 9 | | 84 - 118 | 0 - 20 |
| Zinc | ug/kg | 42700 | 112000 | 119 | | 2 | | 82 - 119 | 0 - 20 |

RPD: 0 out of 4 outside limits
 Spike Recovery: 5 out of 8 outside limits

Column to be used to flag recovery and RPD values with an asterisk
 * Values outside of QC limits

FORM V (PART 1) - IN

V2
POST DIGEST SPIKE SAMPLE RECOVERY

| | | | |
|--------------------|-----------------------------------|-------------------|---------------------------------|
| Report No: | <u>218081813</u> | GCAL PDS ID: | <u>1842322</u> |
| Matrix: | <u>Solid</u> | Parent Sample ID: | <u>WIL01IS02 (21808181302)</u> |
| Analyst: | <u>LWZ</u> | Instrument ID: | <u>ICPMS1</u> |
| Analysis Date: | <u>08/23/18</u> Time: <u>1222</u> | Lab File ID: | <u>2180823A_MS1.b\041SMPL.d</u> |
| Analytical Method: | <u>EPA 6020B</u> | Analytical Batch: | <u>642536</u> |

| ANALYTE | UNITS | SPIKED SAMPLE RESULT | C | SAMPLE RESULT | C | SPIKE ADDED | % R | Q | LCL | UCL |
|----------|-------|----------------------------|---|------------------|---|----------------|-----|---|-----|-----|
| Antimony | ug/kg | 38600 | | 0 | U | 42700 | 90 | | 80 | 120 |
| Copper | ug/kg | 40700 | | 21100 | | 21400 | 92 | | 80 | 120 |
| Lead | ug/kg | 84900 | | 63500 | | 21400 | 100 | | 80 | 120 |
| Zinc | ug/kg | 478000 | | 61200 | | 427000 | 97 | | 80 | 120 |

FORM V (PART 2) - IN

IX
SERIAL DILUTIONS

| | | | |
|--------------------|-----------------------------------|-------------------|---------------------------------|
| Report No: | <u>218081813</u> | GCAL SD ID: | <u>1842323</u> |
| Matrix: | <u>Solid</u> | Parent Sample ID: | <u>WIL01IS02 (21808181302)</u> |
| Analyst: | <u>LWZ</u> | Instrument ID: | <u>ICPMS1</u> |
| Analysis Date: | <u>08/23/18</u> Time: <u>1227</u> | Lab File ID: | <u>2180823A_MS1.b\042SMPL.d</u> |
| Analytical Method: | <u>EPA 6020B</u> | Analytical Batch: | <u>642536</u> |

| ANALYTE | UNITS | PARENT SAMPLE | | SERIAL DILUTION | | % DIFF | Q | LCL | UCL |
|----------|-------|---------------|---|-----------------|---|--------|---|-----|-----|
| | | RESULT | C | RESULT | C | | | | |
| Antimony | ug/kg | 0 | U | 2150 | J | | | | |
| Copper | ug/kg | 21000 | | 22200 | | 5.7 | | 0 | 10 |
| Lead | ug/kg | 63500 | | 62500 | | 1.6 | | 0 | 10 |
| Zinc | ug/kg | 61200 | | 63600 | | 3.9 | | 0 | 10 |

FORM IX - IN



| | | | |
|---------------------------|--|-------------------------------|------------------------|
| ANALYST/TECH <i>JL</i> | START DATE/TIME <i>8:15 8/21/18</i> | END DATE/TIME <i>11:15</i> | BATCH 642531 |
|---------------------------|--|-------------------------------|------------------------|

| # | CLIENT | TYPE | CLIENT ID | GCAL ID | INITIAL WGT (g) | FINAL VOL (mL) | COMMENT | STANDARDS/ REAGENTS |
|----|--------|------|---------------|-------------|-----------------|----------------|---------|--------------------------|
| 1 | QC | MB | MB 1842310 | 1842310 | <i>1.25</i> | <i>50</i> | | GCAL - 8 - 250uL |
| 2 | QC | LCS | LCS 1842311 | 1842311 | <i>1.25</i> | | | <i>7/26/18</i> |
| 3 | 4957 | SAMP | A013 D1 (0-1) | 21808223401 | <i>1.33</i> | | | Sb, Ag, Se SPIKE - 250uL |
| 4 | 4957 | SAMP | A013 D1 (1-2) | 21808223402 | <i>1.25</i> | | | <i>3/6/18</i> |
| 5 | 4957 | SAMP | A013 D1 (2-3) | 21808223403 | <i>1.46</i> | | | Li, B, Zr SPIKE - 250uL |
| 6 | 4838 | SAMP | WIL03IS01 | 21808181301 | <i>1.35</i> | | | <i>3/6/18</i> |
| 7 | 4838 | SAMP | WIL03IS02 | 21808181308 | <i>1.28</i> | | | Si SPIKE - 250uL |
| 8 | 4838 | SAMP | WIL02IS01 | 21808181401 | <i>1.25</i> | | | <i>7/26/18</i> |
| 9 | 4838 | MS | WIL02IS01 MS | 21808181402 | <i>1.25</i> | | | HNO3 |
| 10 | 4838 | MSD | WIL02IS01 MSD | 21808181403 | <i>1.25</i> | | | <i>301-11-16</i> |
| 11 | 4838 | SAMP | WIL02IS02 | 21808181404 | <i>1.31</i> | | | H2O2 |
| 12 | 4838 | SAMP | WIL01IS03 | 21808181405 | <i>1.40</i> | | | <i>7/26/18</i> |
| 13 | 4838 | SAMP | WIL01IS01 | 21808181406 | <i>1.26</i> | | | HCL |
| 14 | 4838 | SAMP | WIL02IS03 | 21808181407 | <i>1.34</i> | | | <i>N/A</i> |
| 15 | 4838 | SAMP | WIL03IS03 | 21808181307 | <i>1.26</i> | | | 1:1 HNO3 |
| 16 | | | | | | | | <i>7/26/18</i> |
| 17 | | | | | | | | |
| 18 | | | | | | | | |
| 19 | | | | | | | | |
| 20 | | | | | | | | |
| 21 | | | | | | | | |
| 22 | | | | | | | | |
| 23 | | | | | | | | |
| 24 | | | | | | | | |
| 25 | | | | | | | | |
| 26 | | | | | | | | |
| 27 | | | | | | | | Digestion Vessel Lot # |
| 28 | | | | | | | | <i>180521</i> |
| 29 | | | | | | | | |
| 30 | | | | | | | | |

EQUIPMENT/CONDITIONS

| | |
|------------------------------|---------------------------|
| DIGESTION BLOCK <i>A1</i> | TEMPERATURE <i>94°</i> |
|------------------------------|---------------------------|

NOTES

Matrix-Soil. 6020_S_EX

XIII
PREPARATION LOG

Report No: 218081813

Prep Method: EPA 3050B

Prep Batch: 642531

| <i>CLIENT SAMPLE ID</i> | <i>GCAL SAMPLE ID</i> | <i>PREP DATE</i> | <i>WEIGHT</i> | <i>UNITS</i> | <i>VOLUME</i> | <i>UNITS</i> |
|-------------------------|-----------------------|------------------|---------------|--------------|---------------|--------------|
| LCS1842311 | 1842311 | 08/25/18 | 1.25 | g | 50 | mL |
| MB1842310 | 1842310 | 08/25/18 | 1.25 | g | 50 | mL |
| WIL02IS01 MS | 21808181402 | 08/25/18 | 1.25 | g | 50 | mL |
| WIL02IS01 MSD | 21808181403 | 08/25/18 | 1.25 | g | 50 | mL |
| WIL03IS01 | 21808181301 | 08/25/18 | 1.35 | g | 50 | mL |
| WIL03IS02 | 21808181308 | 08/25/18 | 1.28 | g | 50 | mL |
| WIL03IS03 | 21808181307 | 08/25/18 | 1.26 | g | 50 | mL |

FORM XIII - IN

XIII
PREPARATION LOG

Report No: 218081814

Prep Method: EPA 3050B

Prep Batch: 642531

| <i>CLIENT SAMPLE ID</i> | <i>GCAL SAMPLE ID</i> | <i>PREP DATE</i> | <i>WEIGHT</i> | <i>UNITS</i> | <i>VOLUME</i> | <i>UNITS</i> |
|-------------------------|-----------------------|------------------|---------------|--------------|---------------|--------------|
| LCS1842311 | 1842311 | 08/25/18 | 1.25 | g | 50 | mL |
| MB1842310 | 1842310 | 08/25/18 | 1.25 | g | 50 | mL |
| WIL01IS01 | 21808181406 | 08/25/18 | 1.26 | g | 50 | mL |
| WIL01IS03 | 21808181405 | 08/25/18 | 1.4 | g | 50 | mL |
| WIL02IS01 | 21808181401 | 08/25/18 | 1.25 | g | 50 | mL |
| WIL02IS01 MS | 21808181402 | 08/25/18 | 1.25 | g | 50 | mL |
| WIL02IS01 MSD | 21808181403 | 08/25/18 | 1.25 | g | 50 | mL |
| WIL02IS02 | 21808181404 | 08/25/18 | 1.31 | g | 50 | mL |
| WIL02IS03 | 21808181407 | 08/25/18 | 1.34 | g | 50 | mL |

FORM XIII - IN

III
METHOD BLANK

Report No: 218081813 Blank ID: MB1842310
Instrument ID: ICPMS2 Lab File ID: 2180828B_MS2.b\1836SMPL_2180828A_MS2.D
Analyst: LWZ Analytical Batch: 642829
Analysis Date: 08/28/18 Time: 1247 Analytical Method: EPA 6020B

| <i>ANALYTE</i> | <i>RESULT</i> | <i>UNITS</i> | <i>Q</i> | <i>DL</i> | <i>LOD</i> | <i>LOQ</i> |
|----------------|---------------|--------------|----------|-----------|------------|------------|
| Antimony | 40.0 | ug/kg | U | 20.0 | 40.0 | 80.0 |
| Copper | 20.0 | ug/kg | U | 10.0 | 20.0 | 40.0 |
| Lead | 20.0 | ug/kg | U | 10.0 | 20.0 | 40.0 |
| Zinc | 400 | ug/kg | U | 200 | 400 | 800 |

FORM III - IN

VII
LABORATORY CONTROL SAMPLE

| | | | |
|--------------|-----------------------------------|--------------------|---|
| Report No: | <u>218081813</u> | GCAL ID: | <u>1842311 (LCS)</u> |
| Matrix: | <u>Solid</u> | Instrument ID: | <u>ICPMS2</u> |
| Analyst: | <u>LWZ</u> | Lab File ID: | <u>2180828B_MS2.b\1837SMPL_2180828A_MS2.D</u> |
| Prep Date: | <u>08/25/18</u> Time: <u>0815</u> | Analysis Date: | <u>08/28/18</u> Time: <u>1250</u> |
| Prep Batch: | <u>642531</u> | Analytical Batch: | <u>642829</u> |
| Prep Method: | <u>3050B</u> | Analytical Method: | <u>EPA 6020B</u> |

| ANALYTE | UNITS | TRUE | FOUND | % R | Q | LCL | UCL |
|----------|-------|-------|-------|-----|---|-----|-----|
| Antimony | ug/kg | 4000 | 4190 | 105 | | 72 | 124 |
| Copper | ug/kg | 2000 | 2010 | 101 | | 84 | 119 |
| Lead | ug/kg | 2000 | 2030 | 102 | | 84 | 118 |
| Zinc | ug/kg | 40000 | 40300 | 101 | | 82 | 119 |

FORM VII - IN

V1
MS/MSD RECOVERY

Report No: 218081814 Parent Sample ID: WIL02IS01
 Prep Method: EPA 3050B \ ISM Parent GCAL ID: 21808181401
 Prep Date: 08/25/18 Time: 0815 Prep Batch: 642531
 Analytical Method: EPA 6020B Analytical Batch: 642829

| | |
|-----------------------------------|---|
| GCAL QC ID: 21808181402 MS | Instrument ID: ICPMS2 |
| Analyst: LWZ | Lab File ID: 2180828B_MS2.b\1844SMPL_2180828A_MS2.D |
| Analysis Date: 08/28/18 1315 | Dilution: 10 |

| ANALYTE | UNITS | SPIKE ADDED | SAMPLE RESULT | MS RESULT | MS % REC | # | QC LIMITS |
|----------|-------|-------------|---------------|-----------|----------|---|-----------|
| Antimony | ug/kg | 4690 | 0 | 315 | 7 | * | 72 - 124 |
| Copper | ug/kg | 2340 | 38400 | 40000 | 67 | * | 84 - 119 |
| Lead | ug/kg | 2340 | 15800 | 18400 | 109 | | 84 - 118 |
| Zinc | ug/kg | 46900 | 88500 | 136000 | 101 | | 82 - 119 |

| | |
|------------------------------------|---|
| GCAL QC ID: 21808181403 MSD | Instrument ID: ICPMS2 |
| Analyst: LWZ | Lab File ID: 2180828B_MS2.b\1845SMPL_2180828A_MS2.D |
| Analysis Date: 08/28/18 1319 | Dilution: 10 |

| ANALYTE | UNITS | SPIKE ADDED | MSD RESULT | MSD % REC | # | % RPD | # | QC LIMITS | RPD |
|----------|-------|-------------|------------|-----------|---|-------|---|-----------|--------|
| Antimony | ug/kg | 4690 | 361 | 8 | * | 14 | | 72 - 124 | 0 - 20 |
| Copper | ug/kg | 2340 | 39900 | 66 | * | 0 | | 84 - 119 | 0 - 20 |
| Lead | ug/kg | 2340 | 17600 | 75 | * | 4 | | 84 - 118 | 0 - 20 |
| Zinc | ug/kg | 46900 | 133000 | 95 | | 2 | | 82 - 119 | 0 - 20 |

RPD : 0 out of 4 outside limits
 Spike Recovery: 5 out of 8 outside limits

Column to be used to flag recovery and RPD values with an asterisk
 * Values outside of QC limits

V2
POST DIGEST SPIKE SAMPLE RECOVERY

| | | | |
|--------------------|-----------------------------------|-------------------|---|
| Report No: | <u>218081814</u> | GCAL PDS ID: | <u>1843727</u> |
| Matrix: | <u>Solid</u> | Parent Sample ID: | <u>WIL02IS01 (21808181401)</u> |
| Analyst: | <u>LWZ</u> | Instrument ID: | <u>ICPMS2</u> |
| Analysis Date: | <u>08/28/18</u> Time: <u>1322</u> | Lab File ID: | <u>2180828C_MS2.b\1846SMPL_2180828A_MS2.D</u> |
| Analytical Method: | <u>EPA 6020B</u> | Analytical Batch: | <u>642829</u> |

| ANALYTE | UNITS | SPIKED SAMPLE RESULT | C | SAMPLE RESULT | C | SPIKE ADDED | % R | Q | LCL | UCL |
|----------|-------|----------------------------|---|------------------|---|----------------|-----|---|-----|-----|
| Antimony | ug/kg | 47000 | | 0 | U | 46900 | 100 | | 80 | 120 |
| Copper | ug/kg | 61200 | | 38400 | | 23400 | 97 | | 80 | 120 |
| Lead | ug/kg | 38400 | | 15800 | | 23400 | 96 | | 80 | 120 |
| Zinc | ug/kg | 553000 | | 88500 | | 469000 | 99 | | 80 | 120 |

IX
SERIAL DILUTIONS

| | | | |
|--------------------|-----------------------------------|-------------------|---|
| Report No: | <u>218081814</u> | GCAL SD ID: | <u>1843728</u> |
| Matrix: | <u>Solid</u> | Parent Sample ID: | <u>WIL02IS01 (21808181401)</u> |
| Analyst: | <u>LWZ</u> | Instrument ID: | <u>ICPMS2</u> |
| Analysis Date: | <u>08/28/18</u> Time: <u>1326</u> | Lab File ID: | <u>2180828C_MS2.b\1847SMPL_2180828A_MS2.D</u> |
| Analytical Method: | <u>EPA 6020B</u> | Analytical Batch: | <u>642829</u> |

| ANALYTE | UNITS | PARENT SAMPLE | | SERIAL DILUTION | | % DIFF | Q | LCL | UCL |
|----------|-------|---------------|---|-----------------|---|--------|---|-----|-----|
| | | RESULT | C | RESULT | C | | | | |
| Antimony | ug/kg | 0 | U | 0 | U | | | | |
| Copper | ug/kg | 38400 | | 40100 | | 4.4 | | 0 | 10 |
| Lead | ug/kg | 15900 | | 15900 | | 0 | | 0 | 10 |
| Zinc | ug/kg | 88500 | | 88300 | | .2 | | 0 | 10 |

FORM IX - IN

Sample Summary

| GCAL ID | Client ID | Matrix | Collect Date/Time | Receive Date/Time |
|----------------|---------------------|---------------|--------------------------|--------------------------|
| 21808181206 | WIL02DA02A | Solid | 08/16/2018 10:20 | 08/18/2018 10:45 |
| 21808181207 | WIL02DA02A MS | Solid | 08/16/2018 10:20 | 08/18/2018 10:45 |
| 21808181208 | WIL02DA02A MSD | Solid | 08/16/2018 10:20 | 08/18/2018 10:45 |
| 21808181209 | WIL02DA01A | Solid | 08/16/2018 09:30 | 08/18/2018 10:45 |
| 21808181210 | WIL02DA01B | Solid | 08/16/2018 09:35 | 08/18/2018 10:45 |
| 21808181213 | WIL02DA02A (RE) | Solid | 08/16/2018 10:20 | 08/18/2018 10:45 |
| 21808181214 | WIL02DA02A MS (RE) | Solid | 08/16/2018 10:20 | 08/18/2018 10:45 |
| 21808181215 | WIL02DA02A MSD (RE) | Solid | 08/16/2018 10:20 | 08/18/2018 10:45 |
| 21808181216 | WIL02DA01A (RE) | Solid | 08/16/2018 09:30 | 08/18/2018 10:45 |
| 21808181217 | WIL02DA01B (RE) | Solid | 08/16/2018 09:35 | 08/18/2018 10:45 |

Sample Summary

| GCAL ID | Client ID | Matrix | Collect Date/Time | Receive Date/Time |
|----------------|------------------|---------------|--------------------------|--------------------------|
| 21808181301 | WIL03IS01 | Solid | 08/14/2018 16:00 | 08/18/2018 10:45 |
| 21808181302 | WIL01IS02 | Solid | 08/14/2018 13:50 | 08/18/2018 10:45 |
| 21808181303 | WIL03IS00 | Water | 08/16/2018 08:20 | 08/18/2018 10:45 |
| 21808181304 | WIL04IS03 | Solid | 08/15/2018 16:10 | 08/18/2018 10:45 |
| 21808181305 | WIL04IS01 | Solid | 08/15/2018 16:00 | 08/18/2018 10:45 |
| 21808181306 | WIL04IS02 | Solid | 08/15/2018 16:05 | 08/18/2018 10:45 |
| 21808181307 | WIL03IS03 | Solid | 08/14/2018 16:10 | 08/18/2018 10:45 |
| 21808181308 | WIL03IS02 | Solid | 08/14/2018 16:05 | 08/18/2018 10:45 |

Sample Summary

| GCAL ID | Client ID | Matrix | Collect Date/Time | Receive Date/Time |
|----------------|------------------|---------------|--------------------------|--------------------------|
| 21808181401 | WIL02IS01 | Solid | 08/15/2018 12:00 | 08/18/2018 10:45 |
| 21808181402 | WIL02IS01 MS | Solid | 08/15/2018 12:00 | 08/18/2018 10:45 |
| 21808181403 | WIL02IS01 MSD | Solid | 08/15/2018 12:00 | 08/18/2018 10:45 |
| 21808181404 | WIL02IS02 | Solid | 08/15/2018 12:10 | 08/18/2018 10:45 |
| 21808181405 | WIL01IS03 | Solid | 08/14/2018 14:00 | 08/18/2018 10:45 |
| 21808181406 | WIL01IS01 | Solid | 08/14/2018 13:40 | 08/18/2018 10:45 |
| 21808181407 | WIL02IS03 | Solid | 08/15/2018 12:20 | 08/18/2018 10:45 |

Case Narrative

Client: AECOM **Report:** 218081812

Gulf Coast Analytical Laboratories received and analyzed the sample(s) listed on the Report Sample Summary page of this report. Receipt of the sample(s) is documented by the attached chain of custody. This applies only to the sample(s) listed in this report. No sample integrity or quality control exceptions were identified unless noted below.

This report was completed in accordance with DOD QSM 5.1 as specified in the contract.

PROJECT MANAGER COMMENTS

Per Jennifer Li's email on 10/5/18, hold samples do not need to be analyzed. (Amanda Cobb 10/05/2018 10:30)

HIGH PERFORMANCE LIQUID CHROMATOGRAPHY

In the EPA 8330B analysis for prep batch 642698, the MS/MSD exhibited RPD failures. The LCS and/or LCSD (standard reference material) recoveries for all analytes were below the lower DOD control limit. The LCS/LCSD exhibited RPD failures. All samples were re-extracted outside holding time in prep batch 643342. The recovery for the surrogate is below the lower control limit for the LCSD (1843191).

In the EPA 8330B analysis for prep batch 643342, all samples were outside the holding time. The samples are reported as samples 21808181213-21808181217 with (RE) added to the client ID. The MS/MSD exhibited recovery and RPD failures. The LCS and/or LCSD (standard reference material) recoveries are above the upper control limit for 1,3,5-Trinitrobenzene, 1,3-Dinitrobenzene, 3-Nitrotoluene, 4-Nitrotoluene, and Pentaerythritol Tetranitrate. These analytes were not detected in the associated samples. The LCS/LCSD recoveries are 0% for Tetryl. This is a poor performing analyte.

METALS

In the EPA 6020B analysis for prep batch 642222, the MS and/or MSD recovery is outside the control limits for Antimony and Zinc. The LCS recovery is within control limits. This indicates the analysis is in control and the sample is affected by matrix interference or the element is non-homogeneous in the sample. A post-digestion spike was performed. The MS/MSD recoveries and RPD are not applicable for Copper and Lead because the sample concentration is greater than four times the spike concentration. Zinc is flagged E on the serial dilution form due to the fact that the % difference between the original result and the serial dilution result for the batch QC sample is greater than 10.

Case Narrative

Client: AECOM **Report:** 218081813

Gulf Coast Analytical Laboratories received and analyzed the sample(s) listed on the Report Sample Summary page of this report. Receipt of the sample(s) is documented by the attached chain of custody. This applies only to the sample(s) listed in this report. No sample integrity or quality control exceptions were identified unless noted below.

This report was completed in accordance with DOD QSM 5.1 as specified in the contract.

METALS

In the EPA 6020B analysis for prep batch 642442, the MS and/or MSD recovery is outside the control limits for Antimony. The LCS recovery is within control limits. This indicates the analysis is in control and the sample is affected by matrix interference or the element is non-homogeneous in the sample. A post-digestion spike was performed. The MS/MSD recoveries and RPD are not applicable for Copper and Lead because the sample concentration is greater than four times the spike concentration.

Case Narrative

Client: AECOM **Report:** 218081814

Gulf Coast Analytical Laboratories received and analyzed the sample(s) listed on the Report Sample Summary page of this report. Receipt of the sample(s) is documented by the attached chain of custody. This applies only to the sample(s) listed in this report. No sample integrity or quality control exceptions were identified unless noted below.

This report was completed in accordance with DOD QSM 5.1 as specified in the contract.

HIGH PERFORMANCE LIQUID CHROMATOGRAPHY

In the EPA 8330B analysis for prep batch 642680, the LCS/LCSD RPD is above the control limit for 4-Nitrotoluene. All recoveries are acceptable.

METALS

In the EPA 6020B analysis for prep batch 642531, the MS and/or MSD recovery is outside the control limits for Antimony. The LCS recovery is within control limits. This indicates the analysis is in control and the sample is affected by matrix interference or the element is non-homogeneous in the sample. A post-digestion spike was performed. The MS/MSD recoveries and RPD are not applicable for Copper and Lead because the sample concentration is greater than four times the spike concentration.

Q Flag Summary

Client Sample ID: **WIL02DA02A** Lab Sample ID: **21808181206**

| Method: EPA 8330B | | Analysis Date: 9/13/2018 5:40:10 AM | | | | |
|------------------------------|------------|-------------------------------------|--------------|---------------|--------|-----------|
| Analyte | CAS | CCV OUL | LCS/LCSD OUL | SURROGATE OUL | IS OUL | CLCCV OUL |
| 1,3,5-Trinitrobenzene | 99-35-4 | | X | | | |
| 1,3-Dinitrobenzene | 99-65-0 | | X | | | |
| 2,4,6-Trinitrotoluene | 118-96-7 | | X | | | |
| 2,4-Dinitrotoluene | 121-14-2 | | X | | | |
| 2,6-Dinitrotoluene | 606-20-2 | | X | | | |
| 2-Amino-4,6-dinitrotoluene | 35572-78-2 | | X | | | |
| 2-Nitrotoluene | 88-72-2 | | X | | | |
| 3,5-Dinitroaniline | 618-87-1 | | X | | | |
| 3-Nitrotoluene | 99-08-1 | | X | | | |
| 4-Amino-2,6-dinitrotoluene | 19406-51-0 | | X | | | |
| 4-Nitrotoluene | 99-99-0 | | X | | | |
| HMX | 2691-41-0 | | X | | | |
| Nitrobenzene | 98-95-3 | | X | | | |
| Nitroglycerin | 55-63-0 | | X | | | |
| Pentaerythritol Tetranitrate | 78-11-5 | | X | | | |
| RDX | 121-82-4 | | X | | | |
| Tetryl | 479-45-8 | | X | | | |

Client Sample ID: **WIL02DA01A** Lab Sample ID: **21808181209**

| Method: EPA 8330B | | Analysis Date: 9/13/2018 6:39:30 AM | | | | |
|------------------------------|------------|-------------------------------------|--------------|---------------|--------|-----------|
| Analyte | CAS | CCV OUL | LCS/LCSD OUL | SURROGATE OUL | IS OUL | CLCCV OUL |
| 1,3,5-Trinitrobenzene | 99-35-4 | | X | | | |
| 1,3-Dinitrobenzene | 99-65-0 | | X | | | |
| 2,4,6-Trinitrotoluene | 118-96-7 | | X | | | |
| 2,4-Dinitrotoluene | 121-14-2 | | X | | | |
| 2,6-Dinitrotoluene | 606-20-2 | | X | | | |
| 2-Amino-4,6-dinitrotoluene | 35572-78-2 | | X | | | |
| 2-Nitrotoluene | 88-72-2 | | X | | | |
| 3,5-Dinitroaniline | 618-87-1 | | X | | | |
| 3-Nitrotoluene | 99-08-1 | | X | | | |
| 4-Amino-2,6-dinitrotoluene | 19406-51-0 | | X | | | |
| 4-Nitrotoluene | 99-99-0 | | X | | | |
| HMX | 2691-41-0 | | X | | | |
| Nitrobenzene | 98-95-3 | | X | | | |
| Nitroglycerin | 55-63-0 | | X | | | |
| Pentaerythritol Tetranitrate | 78-11-5 | | X | | | |
| RDX | 121-82-4 | | X | | | |
| Tetryl | 479-45-8 | | X | | | |

Client Sample ID: **WIL02DA01B** Lab Sample ID: **21808181210**

| Method: EPA 8330B | | Analysis Date: 9/13/2018 6:59:16 AM | | | | |
|----------------------------|------------|-------------------------------------|--------------|---------------|--------|-----------|
| Analyte | CAS | CCV OUL | LCS/LCSD OUL | SURROGATE OUL | IS OUL | CLCCV OUL |
| 1,3,5-Trinitrobenzene | 99-35-4 | | X | | | |
| 1,3-Dinitrobenzene | 99-65-0 | | X | | | |
| 2,4,6-Trinitrotoluene | 118-96-7 | | X | | | |
| 2,4-Dinitrotoluene | 121-14-2 | | X | | | |
| 2,6-Dinitrotoluene | 606-20-2 | | X | | | |
| 2-Amino-4,6-dinitrotoluene | 35572-78-2 | | X | | | |
| 2-Nitrotoluene | 88-72-2 | | X | | | |
| 3,5-Dinitroaniline | 618-87-1 | | X | | | |
| 3-Nitrotoluene | 99-08-1 | | X | | | |
| 4-Amino-2,6-dinitrotoluene | 19406-51-0 | | X | | | |

| | | | | | | |
|------------------------------|-----------|--|---|--|--|--|
| 4-Nitrotoluene | 99-99-0 | | X | | | |
| HMX | 2691-41-0 | | X | | | |
| Nitrobenzene | 98-95-3 | | X | | | |
| Nitroglycerin | 55-63-0 | | X | | | |
| Pentaerythritol Tetranitrate | 78-11-5 | | X | | | |
| RDX | 121-82-4 | | X | | | |
| Tetryl | 479-45-8 | | X | | | |

Client Sample ID: **WIL02DA02A (RE)** Lab Sample ID: **21808181213**

| Method: EPA 8330B Analysis Date: 9/13/2018 8:18:22 AM | | | | | | |
|---|----------|---------|--------------|---------------|--------|-----------|
| Analyte | CAS | CCV OUL | LCS/LCSD OUL | SURROGATE OUL | IS OUL | CLCCV OUL |
| 1,3,5-Trinitrobenzene | 99-35-4 | | X | | | |
| 3-Nitrotoluene | 99-08-1 | | X | | | |
| 4-Nitrotoluene | 99-99-0 | | X | | | |
| Pentaerythritol Tetranitrate | 78-11-5 | | X | | | |
| Tetryl | 479-45-8 | | X | | | |

Client Sample ID: **WIL02DA01A (RE)** Lab Sample ID: **21808181216**

| Method: EPA 8330B Analysis Date: 9/13/2018 9:17:42 AM | | | | | | |
|---|----------|---------|--------------|---------------|--------|-----------|
| Analyte | CAS | CCV OUL | LCS/LCSD OUL | SURROGATE OUL | IS OUL | CLCCV OUL |
| 1,3,5-Trinitrobenzene | 99-35-4 | | X | | | |
| 3-Nitrotoluene | 99-08-1 | | X | | | |
| 4-Nitrotoluene | 99-99-0 | | X | | | |
| Pentaerythritol Tetranitrate | 78-11-5 | | X | | | |
| Tetryl | 479-45-8 | | X | | | |

Client Sample ID: **WIL02DA01B (RE)** Lab Sample ID: **21808181217**

| Method: EPA 8330B Analysis Date: 9/13/2018 9:37:28 AM | | | | | | |
|---|----------|---------|--------------|---------------|--------|-----------|
| Analyte | CAS | CCV OUL | LCS/LCSD OUL | SURROGATE OUL | IS OUL | CLCCV OUL |
| 1,3,5-Trinitrobenzene | 99-35-4 | | X | | | |
| 3-Nitrotoluene | 99-08-1 | | X | | | |
| 4-Nitrotoluene | 99-99-0 | | X | | | |
| Pentaerythritol Tetranitrate | 78-11-5 | | X | | | |
| Tetryl | 479-45-8 | | X | | | |

CCV OUL=CCV out of limits
LCS/LCSD OUL=LCS/LCSD out of limits
SURROGATE OUL=Surrogate out of limits
IS OUL=Internal Standard out of limits
CLCCV OUL=Closing CCV out of limits

Q Flag Summary

Client Sample ID: **WIL02IS01** Lab Sample ID: **21808181401**

| Method: EPA 8330B Analysis Date: 8/30/2018 5:41:00 PM | | | | | | |
|--|---------|---------|--------------|---------------|--------|-----------|
| Analyte | CAS | CCV OUL | LCS/LCSD OUL | SURROGATE OUL | IS OUL | CLCCV OUL |
| 4-Nitrotoluene | 99-99-0 | | X | | | |

Client Sample ID: **WIL02IS02** Lab Sample ID: **21808181404**

| Method: EPA 8330B Analysis Date: 8/30/2018 6:40:20 PM | | | | | | |
|--|---------|---------|--------------|---------------|--------|-----------|
| Analyte | CAS | CCV OUL | LCS/LCSD OUL | SURROGATE OUL | IS OUL | CLCCV OUL |
| 4-Nitrotoluene | 99-99-0 | | X | | | |

Client Sample ID: **WIL02IS03** Lab Sample ID: **21808181407**

| Method: EPA 8330B Analysis Date: 8/30/2018 7:00:07 PM | | | | | | |
|--|---------|---------|--------------|---------------|--------|-----------|
| Analyte | CAS | CCV OUL | LCS/LCSD OUL | SURROGATE OUL | IS OUL | CLCCV OUL |
| 4-Nitrotoluene | 99-99-0 | | X | | | |

CCV OUL=CCV out of limits
 LCS/LCSD OUL=LCS/LCSD out of limits
 SURROGATE OUL=Surrogate out of limits
 IS OUL=Internal Standard out of limits
 CLCCV OUL=Closing CCV out of limits



SAMPLE RECEIVING CHECKLIST



| | | | | |
|--|---|----------------------------------|-------------------------------------|-------------------------------------|
| SAMPLE DELIVERY GROUP 218081813 | | Transport Method FEDEX | YES | NO |
| Client 4838 - AECOM | PM AEC | | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Profile Number 277537 | Received By Savage, Tiffany R | | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Line Item(s) 1 - ISM- Explosives/Metals 3 - Water | Receive Date(s) 08/18/18 | | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| CHECKLIST | | | YES | NO |
| Samples received with proper thermal preservation? | | | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Radioactivity is <1600 cpm? if no, record cpm value in notes section. | | | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| COC relinquished and complete (including sampleIDs, collect times, and sampler)? | | | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| All containers received in good condition and within hold time? | | | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| All sample labels and containers received match the chain of custody? | | | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Preservative added to any containers? | | | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| If received, was headspace for VOC water containers < 6mm? | | | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Samples collected in containers provided by GCAL? | | | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| COOLERS | | | LAB PRESERVATIONS | |
| Airbill 7823-5315-3431 | Thermometer ID: E29 | Temp °C | None | |
| | | 3.2 2.4 1.9 | | |
| DISCREPANCIES | | | | |
| None | | | | |
| NOTES | | | | |

Revision 1.6

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SAMPLE RECEIVING CHECKLIST



| SAMPLE DELIVERY GROUP 218081814 | | CHECKLIST | YES | NO |
|---|-------------------------------------|--|-------------------------------------|-------------------------------------|
| Client 4838 - AECOM | PM AEC Transport Method FEDEX | Samples received with proper thermal preservation? | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Profile Number 277537 | Received By Savage, Tiffany R | Radioactivity is <1600 cpm? If no, record cpm value in notes section. | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Line Item(s) 1 - ISM - Explosives/Metals | Receive Date(s) 08/18/18 | COC relinquished and complete (including sampleIDs, collect times, and sampler)? | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| | | All containers received in good condition and within hold time? | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| | | All sample labels and containers received match the chain of custody? | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| | | Preservative added to any containers? | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| | | If received, was headspace for VOC water containers < 6mm? | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| | | Samples collected in containers provided by GCAL? | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

| COOLERS | DISCREPANCIES | LAB PRESERVATIONS |
|---|---------------|-------------------|
| Airbill: 7823-5315-3431 Thermometer ID: E29 Temp °C: 3.2 2.4 1.9 | None | None |

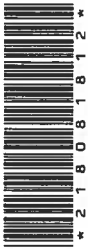
| | |
|--------------|--|
| NOTES | |
|--------------|--|

Revision 1.6

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SAMPLE RECEIVING CHECKLIST



| SAMPLE DELIVERY GROUP 218081812 | | CHECKLIST | YES | NO |
|---|-------------------------------------|--|-------------------------------------|-------------------------------------|
| Client 4838 - AECOM | PM AEC Transport Method FEDEX | Samples received with proper thermal preservation? | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Profile Number 277537 | Received By Savage, Tiffany R | Radioactivity is <1600 cpm? If no, record cpm value in notes section. | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| Line Item(s) 2 - S-Metals/Explosives | Receive Date(s) 08/18/18 | COC relinquished and complete (including sampleIDs, collect times, and sampler)? | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| | | All containers received in good condition and within hold time? | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| | | All sample labels and containers received match the chain of custody? | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| | | Preservative added to any containers? | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| | | If received, was headspace for VOC water containers < 6mm? | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| | | Samples collected in containers provided by GCAL? | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| COOLERS | | DISCREPANCIES | LAB PRESERVATIONS | |
| Airbill 7823-5315-3431 | Thermometer ID: E29 | None | None | |
| | Temp °C 3.2 2.4 1.9 | | | |
| NOTES | | | | |



ANALYTICAL LABORATORIES, LLC
 7979 Innovation Park Dr., Baton Rouge, LA 70820-7402
 Phone: 225.769.4900 • Fax: 225.767.5717 • www.gcal.com

CHAIN OF CUSTODY RECORD

Client ID: 4888 - AECOM

SDG: 218081813

PM: AEC



Report to:

Client: AECOM
 Address: 12420 Milestone Center Dr, 15D
 Germantown, MD 20876
 Contact: Jennifer L.
 Phone: 301-820-3476
 E-mail: jennifer.j.l@aecom.com

Bill to:

Client: Same as
 Address: Report to:
 Contact:
 Phone:
 E-mail:

RO Number

Project Name/Number
 Williston LTA, # 60520956

Sampled By:

Analytical Requests & Method

GCAL use only:
 Custody Seal used yes no
 intact yes no
 Temperature °C 32, 24, 19
 33, 38, 45, 6PM

Dissolved Analysis Requested
 Field filtered
 Lab filtered

Preservative

WHITE: CLIENT FINAL REPORT - CANARY; CLIENT

| Matrix ¹ | Date | Time (2400) | Comp | Grab | Sample Description | No. Containers | Total Metres (6020B) 30' 30" 10' | Explosives (8330B) |
|---------------------|---------|-------------|------|------|--------------------|----------------|----------------------------------|--------------------|
| SOIL | 8-14-18 | 1600 | X | | WIL03IS01 | 1 | X | -1 |
| SOIL | 8-14-18 | 1350 | X | | WIL01IS02 | 1 | X | -2 |
| WATER | 8-16-18 | 0820 | | X | WIL03IS00 | 1 | X | -3 |
| SOIL | 8-15-18 | 1610 | X | | WIL04IS03 | 1 | X | -4 |
| SOIL | 8-15-18 | 1600 | X | | WIL04IS01 | 1 | X | -5 |
| SOIL | 8-15-18 | 1605 | X | | WIL04IS02 | 1 | X | -6 |
| SOIL | 8-14-18 | 1610 | X | | WIL03IS03 | 1 | X | -7 |
| SOIL | 8-14-18 | 1605 | X | | WIL03IS02 | 1 | X | -8 |

Air Bill No: 7822-5315-3431

Turn Around Time (Business Days): 24h* 48h* 3 days* 1 week* Standard (Per Contract/Quote)

Note:

| Requested by (Signature) | Time | Received by (Signature) | Time |
|--------------------------|--------------|-------------------------|--------------|
| | 8/17/18 1100 | | 8/18/18 1045 |
| Requested by (Signature) | Time | Received by (Signature) | Time |
| | 8/18/18 1045 | | 8/18/18 1045 |

Matrix: W = water, S = solid, L = liquid, T = tissue

*Requires prior approval, rush charges may apply.

By submitting these samples, you agree to GCAL's terms and conditions contained in our most recent schedule of services.

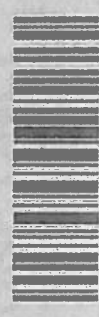
We cannot accept verbal changes. Please email written changes to your PM.



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 Phone: 225-769-4900 • Fax: 225-767-5717 • www.gcal.com

CHAIN OF CUSTODY RECORD

Client ID: 4838 - AECOM
 SDG: 218081814
 PM: AEC



Report to:
 Client: AECOM
 Address: 12420 Milestone Center Dr.
 St. 150, Germantown, MD 20876
 Contact: Jennifer Li
 Phone: 301-820-3476
 E-mail: jennifer.li@aecom.com

Bill to:
 Client:
 Address:
 Contact: "Same as"
 Phone:
 E-mail:

Project Name/Number
 Williston LTA # 60520456

Analytical Requests & Method

GCAL use only:
 Custody Seal used yes no intact yes no
 Temperature °C 32, 24, 19
 33, 38, 45 CPM
 Dissolved Analysis Requested
 Field filtered
 Lab filtered

| Matrix | Date | Time (2400) | Comp | Grab | Sample Description | No. Containers |
|--------|---------|-------------|------|------|--------------------|----------------|
| SOIL | 8-15-18 | 1200 | X | | W1LO2 IS01 | 1 |
| SOIL | 8-15-18 | 1210 | X | | W1LO2 IS02 | 1 |
| SOIL | 8-14-18 | 1400 | X | | W1LO1 IS03 | 1 |
| SOIL | 8-14-18 | 1340 | X | | W1LO1 IS01 | 1 |
| SOIL | 8-6-18 | 1220 | X | | W1LO2 IS03 | 1 |

Total Metals (60208) Pb, Cu, Zn, Cr, Ni
 Explosives (83308) B

| Preservative |
|---------------------|
| MS/MSD extra volume |
| -4 |
| -5 |
| -6 |
| -7 |

Sampled By:

Standard (Per Contract/Quote)
 24hr 48hr 3 days* 1 week* Standard

Air Bill No: 1823-5315-3431

Note: 1 of 4 Sample coolers
 MS/MSD from W1LO2IS01

By submitting these samples, you agree to GCAL's terms and conditions contained in our most recent schedule of services.

We cannot accept verbal changes. Please email written changes to your PM.

Matrix: W = water, S = solid, L = liquid, T = tissue



7979 Innovation Park Dr., Baton Rouge, LA 70820-7402
 Phone: 225.769.4900 • Fax: 225.787.5717 • www.gcal.com

CHAIN OF CUSTODY RECORD

Client ID: 4838 - AECOM

SDG: 218081812

PM: AEC



Report to:

Client: **AECOM**
 Address: **12420 Milestone Center Dr. S1160
 Gaithersburg, MD, 20876**
 Contact: **Jennifer Li**
 Phone: **301-820-3476**
 E-mail: **Jennifer.Li@aecocom.com**

Bill to:

Client: **Same as**
 Address: **" Report to: "**
 Contact:
 Phone:
 E-mail:

P.O. Number: **Williston LTA, # 60520956**

Sampled By:

| Matrix | Date | Time (2400) | Comp | Grab | Sample Description | No. Containers | Total Metals (6020B) Sr, Cu, Pb, Cr | Exposives (8330B) | TCP Metals (lead) 1311/6020B | Agalytical Requests & Method | GCAL use only: |
|--------|---------|-------------|------|------|--------------------|----------------|-------------------------------------|-------------------|------------------------------|------------------------------|---|
| SOIL | 8-16-18 | 1100 | | X | WIL03X18E | 1 | | | | | Custody Seal used <input checked="" type="checkbox"/> yes <input type="checkbox"/> no intact <input type="checkbox"/> yes <input type="checkbox"/> no Temperature °C 3.2, 2.4, 1.9 33, 38, 45 CPM <input type="checkbox"/> Dissolved Analysis Requested <input type="checkbox"/> Field filtered <input type="checkbox"/> Lab filtered |
| SOIL | 8-16-18 | 0920 | | X | WIL02X22E | 1 | | | | | Preservative Hold pending ISM Result -1 |
| SOIL | 8-16-18 | 1115 | | X | WIL01X21E | 1 | | | | | Hold " " |
| SOIL | 8-16-18 | 1030 | | X | WIL02DB02A | 2 | X | X | | | Hold pending "DA" result -1 |
| SOIL | 8-16-18 | 0930 | | X | WIL02DB01A | 2 | X | X | | | Hold " " |
| SOIL | 8-16-18 | 1020 | | X | WIL02DA02A | 6 | X | X | | | Hold -0.778 |
| SOIL | 8-16-18 | 0930 | | X | WIL02DA01A | 2 | X | X | | | Hold -9 |
| SOIL | 8-16-18 | 0935 | | X | WIL02DA01B | 2 | X | X | | | Hold -10 |

Air Bill No: **7823-5315-3431**

Turn Around Time (Business Days): 24h 48h 3 days* 1 week* Standard (Per Contract/Quote)

| | | |
|--------------------------|---------------------------|-------|
| Received by (Signature): | Date: 8/17/18 1100 | Time: |
| Received by (Signature): | Date: 8/18/18 1045 | Time: |
| Received by (Signature): | Date: 8/18/18 1045 | Time: |

Note: **WIL02DA02A is an MS/MSD**

By submitting these samples, you agree to GCAL's terms and conditions contained in our most recent schedule of services.

*Requires prior approval, rush charges may apply.

Matrix: W = water, S = solid, L = liquid, T = tissue

We cannot accept verbal changes. Please email written changes to your PM.

WHITNEY FLETCHER - PRIMARY CLIENT

Data Qualifying Codes

Two types of data qualifying codes or flags are applied in the course of the data review. The data validation flags indicate data that are not usable for decision-making, more than normally biased and/or variable, or not representative of field conditions. These codes and their definitions are presented below in the hierarchy stipulated in the USEPA Contract Laboratory Program National Functional Guidelines for Organic and Inorganic (January 2017) Data Review.

Data Validation Flags

| Flag | Interpretation |
|------|---|
| R | The sample results are unusable due to the quality of the data generated because certain criteria were not met. The analyte may or may not be present in the sample. |
| U | The analyte was analyzed for, but not detected at a level greater than or equal to the level of the adjusted Detection Limit (DL) for sample and method. |
| J+ | Inorganic analyte present. Reported value may not be accurate or precise, but the result may be biased high. |
| J- | Inorganic analyte present. Reported value may not be accurate or precise, but the result may be biased low. |
| J | The analyte was positively identified and the associated numerical value is the approximate concentration of the analyte in the sample (due either to the quality of the data generated because certain quality control criteria were not met, or the concentration of the analyte was below the Limit of Detection (LOD)). |
| NJ | The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration. |
| UJ | The analyte was not detected at a level greater than or equal to the adjusted DL. However, the reported adjusted DL is approximate and may be inaccurate or imprecise. |
| C | This qualifier applies to pesticide and Aroclor results when the identification has been confirmed by gas Chromatograph/Mass Spectrometer (GC/MS) |
| X | This qualifier applies to pesticide and Aroclor results when GC/MS analysis was attempted but was unsuccessful. |

The other type of code used by URS is a "Reason Code". The reason code indicates the type of quality control failure that led to the application of the data validation flag.

Reason Codes

| GC/MS Organics | | GC and HPLC Organics | | Inorganics and Conventionals | |
|----------------|--|----------------------|--|------------------------------|---|
| Code | Interpretation | Code | Interpretation | Code | Interpretation |
| a | Incorrect or incomplete analytical sequence | a | Incorrect or incomplete analytical sequence | a | Incorrect or incomplete analytical sequence |
| b | Bubble found in vial >6mm | b | Instrument performance failure | b | Laboratory duplicate imprecision |
| c | Calibration failure; poor or unstable response | c | Calibration failure; poor or unstable response | c | Calibration failure |
| d | MS/MSD imprecision | d | MS/MSD imprecision | d | MS/MSD imprecision |
| e | LCSD imprecision | e | LCSD imprecision | e | LCSD imprecision |
| f | Field duplicate imprecision | f | Field duplicate imprecision | f | Field duplicate imprecision |
| g | Tuning failure or poor mass spec performance | g | Dual column confirmation imprecision | g | Dual isotope imprecision |
| h | Holding time violation | h | Holding time violation | h | Holding time violation |
| i | Internal standard failure | i | Internal standard failure | k | Cooler receipt temperature exceeds limits |
| k | Cooler receipt temperature exceeds limits | k | Cooler receipt temperature exceeds limits | l | LCSD recovery failure |
| l | LCSD recovery failure | l | LCSD recovery failure | m | MS/MSD recovery failure |
| m | MS/MSD recovery failure | m | MS/MSD recovery failure | n | ICS failure |
| p | Poor chromatography | p | Poor chromatography | o | Calibration blank contamination |
| q | Concentration exceeded the linear range | q | Concentration exceeded the linear range | q | Concentration exceeded the linear range |
| r | Linearity failure in initial calibration | r | Linearity failure in initial calibration | r | Linearity failure in calibration or MSA |
| s | Surrogate failure | s | Surrogate failure | s | Serial dilution failure |
| t | TIC | t | Blender blank contamination | t | Carboy Lot detection |
| w | Identification criteria failure | u | No confirmation column | u | BOD minimum depletion did not exceed 2mg/L |
| x | Field blank contamination | w | Retention time failure | v | Post-digestion spike failure |
| y | Trip blank contamination | x | Field blank contamination | w | CRDL Standard Failure |
| z | Method blank contamination | z | Method blank contamination | x | Field blank contamination |
| | | | | z | Preparation/Method blank contamination |