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Automated Report

Technical Report for

Hydro-Environmental Technology, Inc.

8060.00 Indigo-Desoto Parish, LA

SGS Job Number: LA49126

Sampling Date: 10/23/18

Report to:

**Hydro-Environmental Technology
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Lafayette, LA 70596
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ATTN: Stewart L Stover, Jr.

Total number of pages in report: 77



Test results contained within this data package meet the requirements of the National Environmental Laboratory Accreditation Program and/or state specific certification programs as applicable.

Ron Benjamin
Ron Benjamin
Lab Director

Client Service contact: Ralph Frye 337-237-4775

Certifications: LDEQ(2048), LDHH(LA150012), AR(14-045-04), AZ(AZ0805), FL(E87657), IL(200082), KY(#31), NC(487), SC(73004001), NJ(LA007), TX(T104704186-15-7), WV(257)

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Test results relate only to samples analyzed.

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Sample Summary

Hydro-Environmental Technology, Inc.

Job No: LA49126

8060.00 Indigo-Desoto Parish, LA

| Sample Number | Collected Date | Time By | Received | Matrix Code | Type | Client Sample ID |
|---------------|----------------|---------|---------------|-------------|-------------------|---|
| LA49126-1 | 10/23/18 | 11:25 | LV/EM10/25/18 | AQ | Water | 031-8321Z (NELSON-MASON 1 ALT RIG SUPPLY) |
| LA49126-1F | 10/23/18 | 11:25 | LV/EM10/25/18 | AQ | Water Filtered | 031-8321Z (NELSON-MASON 1 ALT RIG SUPPLY) |
| LA49126-2 | 10/23/18 | 14:20 | LV/EM10/25/18 | AQ | Water | 031-7776Z (NELSON-JOHNSON #3 RIG SUPPLY) |
| LA49126-2F | 10/23/18 | 14:20 | LV/EM10/25/18 | AQ | Water Filtered | 031-7776Z (NELSON-JOHNSON #3 RIG SUPPLY) |
| LA49126-3 | 10/23/18 | 07:15 | LV/EM10/25/18 | AQ | Field Blank Water | FIELD BLANK |
| LA49126-4 | 10/23/18 | 06:30 | LV/EM10/25/18 | AQ | Trip Blank Water | TRIP BLANK |

Sample Results

Report of Analysis

Report of Analysis

| | | | |
|-------------------|---|-----------------|----------|
| Client Sample ID: | 031-8321Z (NELSON-MASON 1 ALT RIG SUPPLY) | Date Sampled: | 10/23/18 |
| Lab Sample ID: | LA49126-1 | Date Received: | 10/25/18 |
| Matrix: | AQ - Water | Percent Solids: | n/a |
| Method: | SW846 8270D SW846 3510C | | |
| Project: | 8060.00 Indigo-Desoto Parish, LA | | |

| Run # | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------|------------|----|----------------|----|----------------|------------|------------------|
| Run #1 | C0039925.D | 1 | 10/31/18 05:35 | IK | 10/29/18 07:19 | OP12657 | EC1681 |
| Run #2 | | | | | | | |

| Run # | Initial Volume | Final Volume |
|--------|----------------|--------------|
| Run #1 | 1040 ml | 1.0 ml |
| Run #2 | | |

ABN RECAP LIST

| CAS No. | Compound | Result | RL | Units | Q |
|----------|------------------------------|--------|---------|-------|---|
| 95-57-8 | 2-Chlorophenol | ND | 0.0048 | mg/l | |
| 120-83-2 | 2,4-Dichlorophenol | ND | 0.0048 | mg/l | |
| 105-67-9 | 2,4-Dimethylphenol | ND | 0.0048 | mg/l | |
| 51-28-5 | 2,4-Dinitrophenol | ND | 0.019 | mg/l | |
| 100-02-7 | 4-Nitrophenol | ND | 0.024 | mg/l | |
| 87-86-5 | Pentachlorophenol | ND | 0.00096 | mg/l | |
| 108-95-2 | Phenol | ND | 0.0048 | mg/l | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | ND | 0.0048 | mg/l | |
| 95-95-4 | 2,4,5-Trichlorophenol | ND | 0.0048 | mg/l | |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | 0.0048 | mg/l | |
| 83-32-9 | Acenaphthene | ND | 0.00019 | mg/l | |
| 208-96-8 | Acenaphthylene | ND | 0.00019 | mg/l | |
| 62-53-3 | Aniline | ND | 0.0048 | mg/l | |
| 120-12-7 | Anthracene | ND | 0.00019 | mg/l | |
| 56-55-3 | Benzo(a)anthracene | ND | 0.00019 | mg/l | |
| 50-32-8 | Benzo(a)pyrene | ND | 0.00019 | mg/l | |
| 205-99-2 | Benzo(b)fluoranthene | ND | 0.00019 | mg/l | |
| 207-08-9 | Benzo(k)fluoranthene | ND | 0.00019 | mg/l | |
| 92-52-4 | 1,1'-Biphenyl | ND | 0.0096 | mg/l | |
| 85-68-7 | Butyl Benzyl Phthalate | ND | 0.0048 | mg/l | |
| 106-47-8 | 4-Chloroaniline ^a | ND | 0.0048 | mg/l | |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | 0.0048 | mg/l | |
| 108-60-1 | 2,2'-Oxybis(1-chloropropane) | ND | 0.0048 | mg/l | |
| 91-58-7 | 2-Chloronaphthalene | ND | 0.0048 | mg/l | |
| 218-01-9 | Chrysene | ND | 0.00019 | mg/l | |
| 53-70-3 | Dibenzo(a,h)anthracene | ND | 0.00019 | mg/l | |
| 132-64-9 | Dibenzofuran | ND | 0.0048 | mg/l | |
| 91-94-1 | 3,3'-Dichlorobenzidine | ND | 0.0096 | mg/l | |
| 84-66-2 | Diethyl Phthalate | ND | 0.0048 | mg/l | |
| 131-11-3 | Dimethyl Phthalate | ND | 0.0048 | mg/l | |
| 117-84-0 | Di-n-octyl Phthalate | ND | 0.0048 | mg/l | |
| 99-65-0 | 1,3-Dinitrobenzene | ND | 0.0048 | mg/l | |

ND = Not detected

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

Report of Analysis

| | | | |
|-------------------|---|-----------------|----------|
| Client Sample ID: | 031-8321Z (NELSON-MASON 1 ALT RIG SUPPLY) | Date Sampled: | 10/23/18 |
| Lab Sample ID: | LA49126-1 | Date Received: | 10/25/18 |
| Matrix: | AQ - Water | Percent Solids: | n/a |
| Method: | SW846 8270D SW846 3510C | | |
| Project: | 8060.00 Indigo-Desoto Parish, LA | | |

ABN RECAP LIST

| CAS No. | Compound | Result | RL | Units | Q |
|----------|----------------------------|--------|---------|-------|---|
| 121-14-2 | 2,4-Dinitrotoluene | ND | 0.0048 | mg/l | |
| 606-20-2 | 2,6-Dinitrotoluene | ND | 0.0048 | mg/l | |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | ND | 0.0048 | mg/l | |
| 206-44-0 | Fluoranthene | ND | 0.00019 | mg/l | |
| 86-73-7 | Fluorene | ND | 0.00019 | mg/l | |
| 118-74-1 | Hexachlorobenzene | ND | 0.00096 | mg/l | |
| 87-68-3 | Hexachlorobutadiene | ND | 0.00048 | mg/l | |
| 77-47-4 | Hexachlorocyclopentadiene | ND | 0.0096 | mg/l | |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | ND | 0.00019 | mg/l | |
| 78-59-1 | Isophorone | ND | 0.0048 | mg/l | |
| 91-57-6 | 2-Methylnaphthalene | ND | 0.00019 | mg/l | |
| 91-20-3 | Naphthalene | ND | 0.00019 | mg/l | |
| 88-74-4 | 2-Nitroaniline | ND | 0.0048 | mg/l | |
| 99-09-2 | 3-Nitroaniline | ND | 0.0048 | mg/l | |
| 100-01-6 | 4-Nitroaniline | ND | 0.0048 | mg/l | |
| 98-95-3 | Nitrobenzene | ND | 0.00096 | mg/l | |
| 621-64-7 | N-Nitroso-di-n-propylamine | ND | 0.0048 | mg/l | |
| 86-30-6 | N-Nitrosodiphenylamine | ND | 0.0048 | mg/l | |
| 85-01-8 | Phenanthrene | ND | 0.00019 | mg/l | |
| 129-00-0 | Pyrene | ND | 0.00019 | mg/l | |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | ND | 0.00096 | mg/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 0.0048 | mg/l | |

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|-----------|----------------------|--------|--------|---------|
| 367-12-4 | 2-Fluorophenol | 44% | | 23-85% |
| 4165-62-2 | Phenol-d5 | 29% | | 10-69% |
| 118-79-6 | 2,4,6-Tribromophenol | 83% | | 48-138% |
| 4165-60-0 | Nitrobenzene-d5 | 82% | | 51-128% |
| 321-60-8 | 2-Fluorobiphenyl | 90% | | 55-122% |
| 1718-51-0 | Terphenyl-d14 | 84% | | 43-138% |

(a) Associated CCV outside of control limits high, sample was ND.

ND = Not detected
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

Report of Analysis

| | |
|--|--------------------------------|
| Client Sample ID: 031-8321Z (NELSON-MASON 1 ALT RIG SUPPLY) | |
| Lab Sample ID: LA49126-1 | Date Sampled: 10/23/18 |
| Matrix: AQ - Water | Date Received: 10/25/18 |
| Method: MADEP VPH REV 1.1 | Percent Solids: n/a |
| Project: 8060.00 Indigo-Desoto Parish, LA | |

| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------|------------|----|----------------|----|-----------|------------|------------------|
| Run #1 | LC381173.D | 1 | 10/27/18 23:09 | SV | n/a | n/a | GLC1882 |
| Run #2 | | | | | | | |

| | Purge Volume |
|--------|--------------|
| Run #1 | 5.0 ml |
| Run #2 | |

Volatile Petroleum Hydrocarbons (VPH)

| CAS No. | Compound | Result | RL | Units | Q |
|---------|------------------------------|--------|------|-------|---|
| | Aliphatics C6-C8 (Unadj.) | ND | 0.15 | mg/l | |
| | Aliphatics > C8-C10 (Unadj.) | ND | 0.15 | mg/l | |
| | Aromatics > C8-C10 (Unadj.) | ND | 0.15 | mg/l | |

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|----------|----------------------|-------------------|--------|---------|
| 615-59-8 | 2,5-Dibromotoluene | 105% ^a | | 70-130% |
| 615-59-8 | 2,5-Dibromotoluene | 103% ^b | | 70-130% |

- (a) Recovery from Aromatics fraction.
- (b) Recovery from Aliphatics fraction.

ND = Not detected
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

Report of Analysis

| | |
|--|--------------------------------|
| Client Sample ID: 031-8321Z (NELSON-MASON 1 ALT RIG SUPPLY) | |
| Lab Sample ID: LA49126-1 | Date Sampled: 10/23/18 |
| Matrix: AQ - Water | Date Received: 10/25/18 |
| Method: MADEP EPH REV 1.1 SW846 3511 | Percent Solids: n/a |
| Project: 8060.00 Indigo-Desoto Parish, LA | |

| Run # | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------|------------|----|----------------|----|----------------|------------|------------------|
| Run #1 | X0005567.D | 1 | 10/29/18 23:08 | JT | 10/26/18 14:00 | OP12635 | GLB1651 |
| Run #2 | Y0005567.D | 1 | 10/29/18 23:09 | JT | 10/26/18 14:00 | OP12635 | GLB1652 |

| Run # | Initial Volume | Final Volume |
|--------|----------------|--------------|
| Run #1 | 54.7 ml | 4.0 ml |
| Run #2 | 54.7 ml | 4.0 ml |

Louisiana EPH Ranges

| CAS No. | Compound | Result | RL | Units | Q |
|---------|-------------------------------|-----------------|------|-------|---|
| | Aliphatics > C10-C12 (Unadj.) | ND ^a | 0.14 | mg/l | |
| | Aliphatics > C12-C16 (Unadj.) | ND ^a | 0.14 | mg/l | |
| | Aliphatics > C16-C35 (Unadj.) | ND ^a | 0.14 | mg/l | |
| | Aromatics > C10-C12 (Unadj.) | ND | 0.14 | mg/l | |
| | Aromatics > C12-C16 (Unadj.) | ND | 0.14 | mg/l | |
| | Aromatics > C16-C21 (Unadj.) | ND | 0.14 | mg/l | |
| | Aromatics > C21-C35 (Unadj.) | ND | 0.14 | mg/l | |

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|-----------|----------------------|--------|--------|---------|
| 3386-33-2 | 1-Chlorooctadecane | | 86% | 40-140% |
| 84-15-1 | o-Terphenyl | 76% | | 40-140% |
| 321-60-8 | 2-Fluorobiphenyl | 78% | | 40-140% |

(a) Result is from Run# 2

ND = Not detected
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

Report of Analysis

| | |
|--|--------------------------------|
| Client Sample ID: 031-8321Z (NELSON-MASON 1 ALT RIG SUPPLY) | Date Sampled: 10/23/18 |
| Lab Sample ID: LA49126-1 | Date Received: 10/25/18 |
| Matrix: AQ - Water | Percent Solids: n/a |
| Project: 8060.00 Indigo-Desoto Parish, LA | |

Total Metals Analysis

| Analyte | Result | RL | Units | DF | Prep | Analyzed By | Method | Prep Method |
|-----------|-----------|---------|-------|----|----------|-------------|--------------------------|--------------------------|
| Aluminum | 1.06 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Arsenic | < 0.010 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Barium | 0.168 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Cadmium | < 0.0050 | 0.0050 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Calcium | 17.4 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Chromium | < 0.010 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Iron | < 1.0 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Lead | < 0.010 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Magnesium | 4.26 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Manganese | 0.0302 | 0.020 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Mercury | < 0.00020 | 0.00020 | mg/l | 1 | 10/29/18 | 10/29/18 SA | SW846 7470A ² | SW846 7470A ⁴ |
| Potassium | 2.74 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Selenium | < 0.050 | 0.050 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Silver | < 0.010 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Sodium | 247 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Strontium | 0.831 | 0.020 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Zinc | < 0.050 | 0.050 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |

- (1) Instrument QC Batch: MA13785
- (2) Instrument QC Batch: MA13797
- (3) Prep QC Batch: MP13152
- (4) Prep QC Batch: MP13160

RL = Reporting Limit

Report of Analysis

| | |
|--|--------------------------------|
| Client Sample ID: 031-8321Z (NELSON-MASON 1 ALT RIG SUPPLY) | Date Sampled: 10/23/18 |
| Lab Sample ID: LA49126-1 | Date Received: 10/25/18 |
| Matrix: AQ - Water | Percent Solids: n/a |
| Project: 8060.00 Indigo-Desoto Parish, LA | |

General Chemistry

| Analyte | Result | RL | Units | DF | Analyzed | By | Method |
|---|--------|------|----------|----|----------------|-----|-------------------|
| Alkalinity, Bicarbonate ^a | 300 | 5.0 | mg/l | 1 | 10/29/18 16:00 | ATX | SM18 2320B |
| Alkalinity, Carbonate ^a | < 5.0 | 5.0 | mg/l | 1 | 10/29/18 16:00 | ATX | SM18 2320B |
| Alkalinity, Total as CaCO ₃ ^a | 305 | 5.0 | mg/l | 1 | 10/29/18 16:00 | ATX | SM 2320B-2011 |
| Bromide ^a | 0.60 | 0.50 | mg/l | 1 | 11/07/18 01:33 | ATX | SW846 9056A |
| Chloride ^a | 87.7 | 5.0 | mg/l | 10 | 11/06/18 15:24 | ATX | SW846 9056A |
| Silica, Dissolved ^a | 7.9 | 0.70 | mg/l | 10 | 11/01/18 | ATX | SM4500SIO2 C-2011 |
| Solids, Total Dissolved ^a | 559 | 10 | mg/l | 1 | 10/29/18 | ATX | SM 2540C-2011 |
| Specific Conductivity ^b | 962 | 1.0 | umhos/cm | 1 | 10/29/18 14:45 | ATX | EPA 120.1 |
| Sulfate ^a | 48.5 | 5.0 | mg/l | 10 | 11/06/18 15:24 | ATX | SW846 9056A |

(a) Analysis performed at SGS Houston, TX.

(b) Conductivity results corrected to 25 degrees Celsius. Analysis performed at SGS Houston, TX.

RL = Reporting Limit

Report of Analysis

| | |
|--|--------------------------------|
| Client Sample ID: 031-8321Z (NELSON-MASON 1 ALT RIG SUPPLY) | Date Sampled: 10/23/18 |
| Lab Sample ID: LA49126-1F | Date Received: 10/25/18 |
| Matrix: AQ - Water Filtered | Percent Solids: n/a |
| Project: 8060.00 Indigo-Desoto Parish, LA | |

Dissolved Metals Analysis

| Analyte | Result | RL | Units | DF | Prep | Analyzed By | Method | Prep Method |
|-----------|-----------|---------|-------|----|----------|-------------|--------------------------|--------------------------|
| Aluminum | < 1.0 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Arsenic | < 0.010 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Barium | 0.149 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Cadmium | < 0.0050 | 0.0050 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Calcium | 14.3 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Chromium | < 0.010 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Iron | < 1.0 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Lead | < 0.010 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Magnesium | 4.37 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Manganese | 0.0234 | 0.020 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Mercury | < 0.00020 | 0.00020 | mg/l | 1 | 10/29/18 | 10/29/18 SA | SW846 7470A ² | SW846 7470A ⁴ |
| Potassium | 2.56 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Selenium | < 0.050 | 0.050 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Silver | < 0.010 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Sodium | 258 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Strontium | 0.886 | 0.020 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Zinc | < 0.050 | 0.050 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |

- (1) Instrument QC Batch: MA13785
- (2) Instrument QC Batch: MA13797
- (3) Prep QC Batch: MP13152
- (4) Prep QC Batch: MP13160

RL = Reporting Limit

Report of Analysis

| | | | | |
|-------------------|--|--|-----------------|----------|
| Client Sample ID: | 031-7776Z (NELSON-JOHNSON #3 RIG SUPPLY) | | Date Sampled: | 10/23/18 |
| Lab Sample ID: | LA49126-2 | | Date Received: | 10/25/18 |
| Matrix: | AQ - Water | | Percent Solids: | n/a |
| Method: | SW846 8270D SW846 3510C | | | |
| Project: | 8060.00 Indigo-Desoto Parish, LA | | | |

| Run # | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------|------------|----|----------------|----|----------------|------------|------------------|
| Run #1 | C0039926.D | 1 | 10/31/18 05:57 | IK | 10/29/18 07:19 | OP12657 | EC1681 |
| Run #2 | | | | | | | |

| Run # | Initial Volume | Final Volume |
|--------|----------------|--------------|
| Run #1 | 1040 ml | 1.0 ml |
| Run #2 | | |

ABN RECAP LIST

| CAS No. | Compound | Result | RL | Units | Q |
|----------|------------------------------|--------|---------|-------|---|
| 95-57-8 | 2-Chlorophenol | ND | 0.0048 | mg/l | |
| 120-83-2 | 2,4-Dichlorophenol | ND | 0.0048 | mg/l | |
| 105-67-9 | 2,4-Dimethylphenol | ND | 0.0048 | mg/l | |
| 51-28-5 | 2,4-Dinitrophenol | ND | 0.019 | mg/l | |
| 100-02-7 | 4-Nitrophenol | ND | 0.024 | mg/l | |
| 87-86-5 | Pentachlorophenol | ND | 0.00096 | mg/l | |
| 108-95-2 | Phenol | ND | 0.0048 | mg/l | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | ND | 0.0048 | mg/l | |
| 95-95-4 | 2,4,5-Trichlorophenol | ND | 0.0048 | mg/l | |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | 0.0048 | mg/l | |
| 83-32-9 | Acenaphthene | ND | 0.00019 | mg/l | |
| 208-96-8 | Acenaphthylene | ND | 0.00019 | mg/l | |
| 62-53-3 | Aniline | ND | 0.0048 | mg/l | |
| 120-12-7 | Anthracene | ND | 0.00019 | mg/l | |
| 56-55-3 | Benzo(a)anthracene | ND | 0.00019 | mg/l | |
| 50-32-8 | Benzo(a)pyrene | ND | 0.00019 | mg/l | |
| 205-99-2 | Benzo(b)fluoranthene | ND | 0.00019 | mg/l | |
| 207-08-9 | Benzo(k)fluoranthene | ND | 0.00019 | mg/l | |
| 92-52-4 | 1,1'-Biphenyl | ND | 0.0096 | mg/l | |
| 85-68-7 | Butyl Benzyl Phthalate | ND | 0.0048 | mg/l | |
| 106-47-8 | 4-Chloroaniline ^a | ND | 0.0048 | mg/l | |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | 0.0048 | mg/l | |
| 108-60-1 | 2,2'-Oxybis(1-chloropropane) | ND | 0.0048 | mg/l | |
| 91-58-7 | 2-Chloronaphthalene | ND | 0.0048 | mg/l | |
| 218-01-9 | Chrysene | ND | 0.00019 | mg/l | |
| 53-70-3 | Dibenzo(a,h)anthracene | ND | 0.00019 | mg/l | |
| 132-64-9 | Dibenzofuran | ND | 0.0048 | mg/l | |
| 91-94-1 | 3,3'-Dichlorobenzidine | ND | 0.0096 | mg/l | |
| 84-66-2 | Diethyl Phthalate | ND | 0.0048 | mg/l | |
| 131-11-3 | Dimethyl Phthalate | ND | 0.0048 | mg/l | |
| 117-84-0 | Di-n-octyl Phthalate | ND | 0.0048 | mg/l | |
| 99-65-0 | 1,3-Dinitrobenzene | ND | 0.0048 | mg/l | |

ND = Not detected

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

Report of Analysis

| | | | |
|-------------------|--|-----------------|----------|
| Client Sample ID: | 031-7776Z (NELSON-JOHNSON #3 RIG SUPPLY) | Date Sampled: | 10/23/18 |
| Lab Sample ID: | LA49126-2 | Date Received: | 10/25/18 |
| Matrix: | AQ - Water | Percent Solids: | n/a |
| Method: | SW846 8270D SW846 3510C | | |
| Project: | 8060.00 Indigo-Desoto Parish, LA | | |

ABN RECAP LIST

| CAS No. | Compound | Result | RL | Units | Q |
|----------|----------------------------|--------|---------|-------|---|
| 121-14-2 | 2,4-Dinitrotoluene | ND | 0.0048 | mg/l | |
| 606-20-2 | 2,6-Dinitrotoluene | ND | 0.0048 | mg/l | |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | ND | 0.0048 | mg/l | |
| 206-44-0 | Fluoranthene | ND | 0.00019 | mg/l | |
| 86-73-7 | Fluorene | ND | 0.00019 | mg/l | |
| 118-74-1 | Hexachlorobenzene | ND | 0.00096 | mg/l | |
| 87-68-3 | Hexachlorobutadiene | ND | 0.00048 | mg/l | |
| 77-47-4 | Hexachlorocyclopentadiene | ND | 0.0096 | mg/l | |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | ND | 0.00019 | mg/l | |
| 78-59-1 | Isophorone | ND | 0.0048 | mg/l | |
| 91-57-6 | 2-Methylnaphthalene | ND | 0.00019 | mg/l | |
| 91-20-3 | Naphthalene | ND | 0.00019 | mg/l | |
| 88-74-4 | 2-Nitroaniline | ND | 0.0048 | mg/l | |
| 99-09-2 | 3-Nitroaniline | ND | 0.0048 | mg/l | |
| 100-01-6 | 4-Nitroaniline | ND | 0.0048 | mg/l | |
| 98-95-3 | Nitrobenzene | ND | 0.00096 | mg/l | |
| 621-64-7 | N-Nitroso-di-n-propylamine | ND | 0.0048 | mg/l | |
| 86-30-6 | N-Nitrosodiphenylamine | ND | 0.0048 | mg/l | |
| 85-01-8 | Phenanthrene | ND | 0.00019 | mg/l | |
| 129-00-0 | Pyrene | ND | 0.00019 | mg/l | |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | ND | 0.00096 | mg/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 0.0048 | mg/l | |

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|-----------|----------------------|--------|--------|---------|
| 367-12-4 | 2-Fluorophenol | 49% | | 23-85% |
| 4165-62-2 | Phenol-d5 | 34% | | 10-69% |
| 118-79-6 | 2,4,6-Tribromophenol | 85% | | 48-138% |
| 4165-60-0 | Nitrobenzene-d5 | 85% | | 51-128% |
| 321-60-8 | 2-Fluorobiphenyl | 92% | | 55-122% |
| 1718-51-0 | Terphenyl-d14 | 86% | | 43-138% |

(a) Associated CCV outside of control limits high, sample was ND.

ND = Not detected

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

Report of Analysis

| | |
|---|--------------------------------|
| Client Sample ID: 031-7776Z (NELSON-JOHNSON #3 RIG SUPPLY) | Date Sampled: 10/23/18 |
| Lab Sample ID: LA49126-2 | Date Received: 10/25/18 |
| Matrix: AQ - Water | Percent Solids: n/a |
| Method: MADEP VPH REV 1.1 | |
| Project: 8060.00 Indigo-Desoto Parish, LA | |

| Run # | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------|------------|----|----------------|----|-----------|------------|------------------|
| Run #1 | LC381174.D | 1 | 10/27/18 23:53 | SV | n/a | n/a | GLC1882 |
| Run #2 | | | | | | | |

| Run # | Purge Volume |
|--------|--------------|
| Run #1 | 5.0 ml |
| Run #2 | |

Volatile Petroleum Hydrocarbons (VPH)

| CAS No. | Compound | Result | RL | Units | Q |
|---------|------------------------------|--------|------|-------|---|
| | Aliphatics C6-C8 (Unadj.) | ND | 0.15 | mg/l | |
| | Aliphatics > C8-C10 (Unadj.) | ND | 0.15 | mg/l | |
| | Aromatics > C8-C10 (Unadj.) | ND | 0.15 | mg/l | |

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|----------|----------------------|-------------------|--------|---------|
| 615-59-8 | 2,5-Dibromotoluene | 102% ^a | | 70-130% |
| 615-59-8 | 2,5-Dibromotoluene | 104% ^b | | 70-130% |

- (a) Recovery from Aromatics fraction.
- (b) Recovery from Aliphatics fraction.

ND = Not detected
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

Report of Analysis

| | |
|---|--------------------------------|
| Client Sample ID: 031-7776Z (NELSON-JOHNSON #3 RIG SUPPLY) | |
| Lab Sample ID: LA49126-2 | Date Sampled: 10/23/18 |
| Matrix: AQ - Water | Date Received: 10/25/18 |
| Method: MADEP EPH REV 1.1 SW846 3511 | Percent Solids: n/a |
| Project: 8060.00 Indigo-Desoto Parish, LA | |

| Run # | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------|------------|----|----------------|----|----------------|------------|------------------|
| Run #1 | X0005568.D | 1 | 10/29/18 23:33 | JT | 10/26/18 14:00 | OP12635 | GLB1651 |
| Run #2 | Y0005568.D | 1 | 10/29/18 23:34 | JT | 10/26/18 14:00 | OP12635 | GLB1652 |

| Run # | Initial Volume | Final Volume |
|--------|----------------|--------------|
| Run #1 | 55.3 ml | 4.0 ml |
| Run #2 | 55.3 ml | 4.0 ml |

Louisiana EPH Ranges

| CAS No. | Compound | Result | RL | Units | Q |
|---------|-------------------------------|-----------------|------|-------|---|
| | Aliphatics > C10-C12 (Unadj.) | ND ^a | 0.14 | mg/l | |
| | Aliphatics > C12-C16 (Unadj.) | ND ^a | 0.14 | mg/l | |
| | Aliphatics > C16-C35 (Unadj.) | ND ^a | 0.14 | mg/l | |
| | Aromatics > C10-C12 (Unadj.) | ND | 0.14 | mg/l | |
| | Aromatics > C12-C16 (Unadj.) | ND | 0.14 | mg/l | |
| | Aromatics > C16-C21 (Unadj.) | ND | 0.14 | mg/l | |
| | Aromatics > C21-C35 (Unadj.) | ND | 0.14 | mg/l | |

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|-----------|----------------------|--------|--------|---------|
| 3386-33-2 | 1-Chlorooctadecane | | 85% | 40-140% |
| 84-15-1 | o-Terphenyl | 81% | | 40-140% |
| 321-60-8 | 2-Fluorobiphenyl | 81% | | 40-140% |

(a) Result is from Run# 2

ND = Not detected
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

Report of Analysis

| | |
|---|--------------------------------|
| Client Sample ID: 031-7776Z (NELSON-JOHNSON #3 RIG SUPPLY) | Date Sampled: 10/23/18 |
| Lab Sample ID: LA49126-2 | Date Received: 10/25/18 |
| Matrix: AQ - Water | Percent Solids: n/a |
| Project: 8060.00 Indigo-Desoto Parish, LA | |

Total Metals Analysis

| Analyte | Result | RL | Units | DF | Prep | Analyzed By | Method | Prep Method |
|-----------|-----------|---------|-------|----|----------|-------------|--------------------------|--------------------------|
| Aluminum | < 1.0 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Arsenic | < 0.010 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Barium | 0.0322 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Cadmium | < 0.0050 | 0.0050 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Calcium | 1.44 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Chromium | < 0.010 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Iron | < 1.0 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Lead | < 0.010 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Magnesium | < 1.0 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Manganese | < 0.020 | 0.020 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Mercury | < 0.00020 | 0.00020 | mg/l | 1 | 10/29/18 | 10/29/18 SA | SW846 7470A ² | SW846 7470A ⁴ |
| Potassium | 1.27 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Selenium | < 0.050 | 0.050 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Silver | < 0.010 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Sodium | 324 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Strontium | 0.103 | 0.020 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Zinc | < 0.050 | 0.050 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |

- (1) Instrument QC Batch: MA13785
- (2) Instrument QC Batch: MA13797
- (3) Prep QC Batch: MP13152
- (4) Prep QC Batch: MP13160

RL = Reporting Limit

Report of Analysis

| | |
|---|--------------------------------|
| Client Sample ID: 031-7776Z (NELSON-JOHNSON #3 RIG SUPPLY) | Date Sampled: 10/23/18 |
| Lab Sample ID: LA49126-2 | Date Received: 10/25/18 |
| Matrix: AQ - Water | Percent Solids: n/a |
| Project: 8060.00 Indigo-Desoto Parish, LA | |

General Chemistry

| Analyte | Result | RL | Units | DF | Analyzed | By | Method |
|---|--------|------|----------|----|----------------|-----|-------------------|
| Alkalinity, Bicarbonate ^a | 372 | 5.0 | mg/l | 1 | 10/29/18 16:00 | ATX | SM18 2320B |
| Alkalinity, Carbonate ^a | 17.9 | 5.0 | mg/l | 1 | 10/29/18 16:00 | ATX | SM18 2320B |
| Alkalinity, Total as CaCO ₃ ^a | 390 | 5.0 | mg/l | 1 | 10/29/18 16:00 | ATX | SM 2320B-2011 |
| Bromide ^a | 0.50 | 0.50 | mg/l | 1 | 11/07/18 12:58 | ATX | SW846 9056A |
| Chloride ^a | 69.5 | 2.5 | mg/l | 5 | 11/06/18 16:15 | ATX | SW846 9056A |
| Silica, Dissolved ^a | 6.8 | 0.70 | mg/l | 10 | 11/01/18 | ATX | SM4500SIO2 C-2011 |
| Solids, Total Dissolved ^a | 638 | 10 | mg/l | 1 | 10/29/18 | ATX | SM 2540C-2011 |
| Specific Conductivity ^b | 1080 | 1.0 | umhos/cm | 1 | 10/29/18 14:45 | ATX | EPA 120.1 |
| Sulfate ^a | 6.4 | 0.50 | mg/l | 1 | 11/07/18 12:58 | ATX | SW846 9056A |

(a) Analysis performed at SGS Houston, TX.

(b) Conductivity results corrected to 25 degrees Celsius. Analysis performed at SGS Houston, TX.

RL = Reporting Limit

Report of Analysis

| | |
|---|--------------------------------|
| Client Sample ID: 031-7776Z (NELSON-JOHNSON #3 RIG SUPPLY) | Date Sampled: 10/23/18 |
| Lab Sample ID: LA49126-2F | Date Received: 10/25/18 |
| Matrix: AQ - Water Filtered | Percent Solids: n/a |
| Project: 8060.00 Indigo-Desoto Parish, LA | |

Dissolved Metals Analysis

| Analyte | Result | RL | Units | DF | Prep | Analyzed By | Method | Prep Method |
|-----------|-----------|---------|-------|----|----------|-------------|--------------------------|--------------------------|
| Aluminum | < 1.0 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Arsenic | < 0.010 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Barium | 0.0243 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Cadmium | < 0.0050 | 0.0050 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Calcium | 1.11 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Chromium | < 0.010 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Iron | < 1.0 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Lead | < 0.010 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Magnesium | < 1.0 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Manganese | < 0.020 | 0.020 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Mercury | < 0.00020 | 0.00020 | mg/l | 1 | 10/29/18 | 10/29/18 SA | SW846 7470A ² | SW846 7470A ⁴ |
| Potassium | 1.07 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Selenium | < 0.050 | 0.050 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Silver | < 0.010 | 0.010 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Sodium | 294 | 1.0 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Strontium | 0.0895 | 0.020 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |
| Zinc | < 0.050 | 0.050 | mg/l | 10 | 10/26/18 | 10/29/18 RT | SW846 6020A ¹ | SW846 3010A ³ |

- (1) Instrument QC Batch: MA13785
- (2) Instrument QC Batch: MA13797
- (3) Prep QC Batch: MP13152
- (4) Prep QC Batch: MP13160

RL = Reporting Limit

Report of Analysis

| | | | |
|-------------------|----------------------------------|-----------------|----------|
| Client Sample ID: | FIELD BLANK | Date Sampled: | 10/23/18 |
| Lab Sample ID: | LA49126-3 | Date Received: | 10/25/18 |
| Matrix: | AQ - Field Blank Water | Percent Solids: | n/a |
| Method: | SW846 8260B | | |
| Project: | 8060.00 Indigo-Desoto Parish, LA | | |

| Run # | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------|------------|----|----------------|----|-----------|------------|------------------|
| Run #1 | 11040657.D | 1 | 10/29/18 18:11 | LS | n/a | n/a | V111896 |
| Run #2 | | | | | | | |

| Run # | Purge Volume |
|--------|--------------|
| Run #1 | 5.0 ml |
| Run #2 | |

VOA RECAP List

| CAS No. | Compound | Result | RL | Units | Q |
|------------|-----------------------------|--------|--------|-------|---|
| 67-64-1 | Acetone | ND | 0.050 | mg/l | |
| 71-43-2 | Benzene | ND | 0.0050 | mg/l | |
| 75-27-4 | Bromodichloromethane | ND | 0.0010 | mg/l | |
| 75-25-2 | Bromoform | ND | 0.0010 | mg/l | |
| 75-15-0 | Carbon Disulfide | ND | 0.0010 | mg/l | |
| 56-23-5 | Carbon Tetrachloride | ND | 0.0010 | mg/l | |
| 108-90-7 | Chlorobenzene | ND | 0.0010 | mg/l | |
| 75-00-3 | Chloroethane | ND | 0.0010 | mg/l | |
| 67-66-3 | Chloroform | ND | 0.0010 | mg/l | |
| 124-48-1 | Dibromochloromethane | ND | 0.0010 | mg/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 0.0010 | mg/l | |
| 541-73-1 | m-Dichlorobenzene | ND | 0.0010 | mg/l | |
| 95-50-1 | o-Dichlorobenzene | ND | 0.0010 | mg/l | |
| 106-46-7 | p-Dichlorobenzene | ND | 0.0010 | mg/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 0.0010 | mg/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 0.0010 | mg/l | |
| 75-35-4 | 1,1-Dichloroethylene | ND | 0.0010 | mg/l | |
| 156-59-2 | cis-1,2-Dichloroethylene | ND | 0.0010 | mg/l | |
| 156-60-5 | trans-1,2-Dichloroethylene | ND | 0.0010 | mg/l | |
| 540-59-0 | 1,2-Dichloroethene (total) | ND | 0.0010 | mg/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 0.0010 | mg/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 0.0010 | mg/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 0.0010 | mg/l | |
| 542-75-6 | 1,3-Dichloropropene (total) | ND | 0.0010 | mg/l | |
| 100-41-4 | Ethylbenzene | ND | 0.0050 | mg/l | |
| 67-72-1 | Hexachloroethane | ND | 0.0010 | mg/l | |
| 78-83-1 | Isobutyl Alcohol | ND | 0.10 | mg/l | |
| 74-83-9 | Methyl Bromide | ND | 0.0010 | mg/l | |
| 74-87-3 | Methyl Chloride | ND | 0.0010 | mg/l | |
| 75-09-2 | Methylene Chloride | ND | 0.0010 | mg/l | |
| 78-93-3 | Methyl Ethyl Ketone | ND | 0.013 | mg/l | |
| 108-10-1 | 4-Methyl-2-pentanone | ND | 0.013 | mg/l | |

ND = Not detected

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

Report of Analysis

| | | |
|--|--|--------------------------------|
| Client Sample ID: FIELD BLANK | | Date Sampled: 10/23/18 |
| Lab Sample ID: LA49126-3 | | Date Received: 10/25/18 |
| Matrix: AQ - Field Blank Water | | Percent Solids: n/a |
| Method: SW846 8260B | | |
| Project: 8060.00 Indigo-Desoto Parish, LA | | |

VOA RECAP List

| CAS No. | Compound | Result | RL | Units | Q |
|-----------|---------------------------|--------|---------|-------|---|
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 0.0050 | mg/l | |
| 100-42-5 | Styrene | ND | 0.0010 | mg/l | |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | ND | 0.0010 | mg/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 0.00050 | mg/l | |
| 127-18-4 | Tetrachloroethylene | ND | 0.0010 | mg/l | |
| 108-88-3 | Toluene | ND | 0.0050 | mg/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 0.0010 | mg/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 0.0010 | mg/l | |
| 79-01-6 | Trichloroethylene | ND | 0.0010 | mg/l | |
| 75-69-4 | Trichlorofluoromethane | ND | 0.0010 | mg/l | |
| 75-01-4 | Vinyl Chloride | ND | 0.0010 | mg/l | |
| | m,p-Xylene | ND | 0.0050 | mg/l | |
| 95-47-6 | o-Xylene | ND | 0.0050 | mg/l | |
| 1330-20-7 | Xylene (total) | ND | 0.0050 | mg/l | |

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|------------|-----------------------|--------|--------|---------|
| 17060-07-0 | 1,2-Dichloroethane-D4 | 107% | | 84-124% |
| 2037-26-5 | Toluene-D8 | 98% | | 83-115% |
| 460-00-4 | 4-Bromofluorobenzene | 94% | | 89-111% |

ND = Not detected
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

Report of Analysis

| | | |
|--|--|--------------------------------|
| Client Sample ID: FIELD BLANK | | |
| Lab Sample ID: LA49126-3 | | Date Sampled: 10/23/18 |
| Matrix: AQ - Field Blank Water | | Date Received: 10/25/18 |
| Method: MADEP VPH REV 1.1 | | Percent Solids: n/a |
| Project: 8060.00 Indigo-Desoto Parish, LA | | |

| | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------|------------|----|----------------|----|-----------|------------|------------------|
| Run #1 | LC381179.D | 1 | 10/28/18 03:34 | SV | n/a | n/a | GLC1882 |
| Run #2 | | | | | | | |

| | Purge Volume |
|--------|--------------|
| Run #1 | 5.0 ml |
| Run #2 | |

Volatile Petroleum Hydrocarbons (VPH)

| CAS No. | Compound | Result | RL | Units | Q |
|---------|------------------------------|--------|------|-------|---|
| | Aliphatics C6-C8 (Unadj.) | ND | 0.15 | mg/l | |
| | Aliphatics > C8-C10 (Unadj.) | ND | 0.15 | mg/l | |
| | Aromatics > C8-C10 (Unadj.) | ND | 0.15 | mg/l | |

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|----------|----------------------|-------------------|--------|---------|
| 615-59-8 | 2,5-Dibromotoluene | 100% ^a | | 70-130% |
| 615-59-8 | 2,5-Dibromotoluene | 96% ^b | | 70-130% |

- (a) Recovery from Aromatics fraction.
- (b) Recovery from Aliphatics fraction.

ND = Not detected
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

Report of Analysis

| | | | |
|-------------------|----------------------------------|-----------------|----------|
| Client Sample ID: | TRIP BLANK | Date Sampled: | 10/23/18 |
| Lab Sample ID: | LA49126-4 | Date Received: | 10/25/18 |
| Matrix: | AQ - Trip Blank Water | Percent Solids: | n/a |
| Method: | SW846 8260B | | |
| Project: | 8060.00 Indigo-Desoto Parish, LA | | |

| Run # | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|---------------------|------------|----|----------------|----|-----------|------------|------------------|
| Run #1 | 2G090289.D | 1 | 10/31/18 21:46 | LS | n/a | n/a | V2G4705 |
| Run #2 ^a | 1I040655.D | 1 | 10/29/18 17:43 | LS | n/a | n/a | V1I1896 |

| Run # | Purge Volume |
|--------|--------------|
| Run #1 | 5.0 ml |
| Run #2 | 5.0 ml |

VOA RECAP List

| CAS No. | Compound | Result | RL | Units | Q |
|------------|-----------------------------|--------|--------|-------|---|
| 67-64-1 | Acetone | ND | 0.050 | mg/l | |
| 71-43-2 | Benzene | ND | 0.0050 | mg/l | |
| 75-27-4 | Bromodichloromethane | ND | 0.0010 | mg/l | |
| 75-25-2 | Bromoform | ND | 0.0010 | mg/l | |
| 75-15-0 | Carbon Disulfide | ND | 0.0010 | mg/l | |
| 56-23-5 | Carbon Tetrachloride | ND | 0.0010 | mg/l | |
| 108-90-7 | Chlorobenzene ^b | 0.0010 | 0.0010 | mg/l | |
| 75-00-3 | Chloroethane | ND | 0.0010 | mg/l | |
| 67-66-3 | Chloroform | ND | 0.0010 | mg/l | |
| 124-48-1 | Dibromochloromethane | ND | 0.0010 | mg/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 0.0010 | mg/l | |
| 541-73-1 | m-Dichlorobenzene | ND | 0.0010 | mg/l | |
| 95-50-1 | o-Dichlorobenzene | ND | 0.0010 | mg/l | |
| 106-46-7 | p-Dichlorobenzene | ND | 0.0010 | mg/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 0.0010 | mg/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 0.0010 | mg/l | |
| 75-35-4 | 1,1-Dichloroethylene | ND | 0.0010 | mg/l | |
| 156-59-2 | cis-1,2-Dichloroethylene | ND | 0.0010 | mg/l | |
| 156-60-5 | trans-1,2-Dichloroethylene | ND | 0.0010 | mg/l | |
| 540-59-0 | 1,2-Dichloroethene (total) | ND | 0.0010 | mg/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 0.0010 | mg/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 0.0010 | mg/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 0.0010 | mg/l | |
| 542-75-6 | 1,3-Dichloropropene (total) | ND | 0.0010 | mg/l | |
| 100-41-4 | Ethylbenzene | ND | 0.0050 | mg/l | |
| 67-72-1 | Hexachloroethane | ND | 0.0010 | mg/l | |
| 78-83-1 | Isobutyl Alcohol | ND | 0.10 | mg/l | |
| 74-83-9 | Methyl Bromide | ND | 0.0010 | mg/l | |
| 74-87-3 | Methyl Chloride | ND | 0.0010 | mg/l | |
| 75-09-2 | Methylene Chloride | ND | 0.0010 | mg/l | |
| 78-93-3 | Methyl Ethyl Ketone | ND | 0.013 | mg/l | |
| 108-10-1 | 4-Methyl-2-pentanone | ND | 0.013 | mg/l | |

ND = Not detected

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

Report of Analysis

| | | | |
|-------------------|----------------------------------|-----------------|----------|
| Client Sample ID: | TRIP BLANK | Date Sampled: | 10/23/18 |
| Lab Sample ID: | LA49126-4 | Date Received: | 10/25/18 |
| Matrix: | AQ - Trip Blank Water | Percent Solids: | n/a |
| Method: | SW846 8260B | | |
| Project: | 8060.00 Indigo-Desoto Parish, LA | | |

VOA RECAP List

| CAS No. | Compound | Result | RL | Units | Q |
|-----------|---------------------------|--------|---------|-------|---|
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 0.0050 | mg/l | |
| 100-42-5 | Styrene | ND | 0.0010 | mg/l | |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | ND | 0.0010 | mg/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 0.00050 | mg/l | |
| 127-18-4 | Tetrachloroethylene | ND | 0.0010 | mg/l | |
| 108-88-3 | Toluene | ND | 0.0050 | mg/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 0.0010 | mg/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 0.0010 | mg/l | |
| 79-01-6 | Trichloroethylene | ND | 0.0010 | mg/l | |
| 75-69-4 | Trichlorofluoromethane | ND | 0.0010 | mg/l | |
| 75-01-4 | Vinyl Chloride | ND | 0.0010 | mg/l | |
| | m,p-Xylene | ND | 0.0050 | mg/l | |
| 95-47-6 | o-Xylene | ND | 0.0050 | mg/l | |
| 1330-20-7 | Xylene (total) | ND | 0.0050 | mg/l | |

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|------------|-----------------------|--------|--------|---------|
| 17060-07-0 | 1,2-Dichloroethane-D4 | 101% | 104% | 84-124% |
| 2037-26-5 | Toluene-D8 | 98% | 99% | 83-115% |
| 460-00-4 | 4-Bromofluorobenzene | 103% | 94% | 89-111% |

- (a) Confirmation run.
(b) Confirmed by reanalysis.

ND = Not detected
RL = Reporting Limit
E = Indicates value exceeds calibration range

J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

Report of Analysis

| | |
|--|--------------------------------|
| Client Sample ID: TRIP BLANK | Date Sampled: 10/23/18 |
| Lab Sample ID: LA49126-4 | Date Received: 10/25/18 |
| Matrix: AQ - Trip Blank Water | Percent Solids: n/a |
| Method: MADEP VPH REV 1.1 | |
| Project: 8060.00 Indigo-Desoto Parish, LA | |

| Run # | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------|------------|----|----------------|----|-----------|------------|------------------|
| Run #1 | LC381180.D | 1 | 10/28/18 04:18 | SV | n/a | n/a | GLC1882 |
| Run #2 | | | | | | | |

| Run # | Purge Volume |
|--------|--------------|
| Run #1 | 5.0 ml |
| Run #2 | |

Volatile Petroleum Hydrocarbons (VPH)

| CAS No. | Compound | Result | RL | Units | Q |
|---------|------------------------------|--------|------|-------|---|
| | Aliphatics C6-C8 (Unadj.) | ND | 0.15 | mg/l | |
| | Aliphatics > C8-C10 (Unadj.) | ND | 0.15 | mg/l | |
| | Aromatics > C8-C10 (Unadj.) | ND | 0.15 | mg/l | |

| CAS No. | Surrogate Recoveries | Run# 1 | Run# 2 | Limits |
|----------|----------------------|-------------------|--------|---------|
| 615-59-8 | 2,5-Dibromotoluene | 105% ^a | | 70-130% |
| 615-59-8 | 2,5-Dibromotoluene | 100% ^b | | 70-130% |

- (a) Recovery from Aromatics fraction.
- (b) Recovery from Aliphatics fraction.

ND = Not detected
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

Misc. Forms

Custody Documents and Other Forms

Includes the following where applicable:

- Chain of Custody



HYDRO-ENVIRONMENTAL TECHNOLOGY, INC.
 Environmental Consultants
 P.O. Box 60295
 Lafayette, LA 70596-0295
 Phone (337) 261-1963 FAX (337) 261-1953

Handwritten signature/initials

SAMPLE CHAIN-OF-CUSTODY RECORD

Project Name: Indigo
 Project Number: 8060.00
 Project Location: DeSoto Parish, Louisiana
 Laboratory: SGS Lafayette
 Collected By: LY/EM
 Company: Hydro-Environmental Technology, Inc.
 Date: 10/23/2018

| Sample I.D. | Type | Date/Time Sampled | Containers | Analysis Requested/Method | Comments |
|---|------|-------------------|--|--|--|
| 031-8321Z (Nelson-Mason Tilt Rig Supply) | AQ | 10/23/2018 11:25 | (4) 40mL Glass HCl (3) 60mL Amber Glass HCl (1) L Amber Glass (1) 500mL Plastic (2) 250mL Plastic HNO3 | SVOC 8270, VPH, EPH, Chlorides, TDS, Specific Conductance, Silica, Cations*, Anions*, Total Metals*, Dissolved Metals* | 4°C Field filtered: Dissolved metals |
| 031-776Z (Nelson-Johnson #3 Rig Supply) | AQ | 10/23/2018 14:20 | (4) 40mL Glass HCl (3) 60mL Amber Glass HCl (1) L Amber Glass (1) 500mL Plastic (2) 250mL Plastic HNO4 | SVOC 8270, VPH, EPH, Chlorides, TDS, Specific Conductance, Silica, Cations*, Anions*, Total Metals*, Dissolved Metals* | 4°C Field filtered: Dissolved metals |
| Field Blank | AQ | 10/23/2018 7:15 | (8) 40mL Glass HCL | VOC 8260, VPH | 4°C |
| Trip Blank | AQ | 10/23/2018 6:30 | (8) 40mL Glass HCL | VOC 8260, VPH | 4°C |

*Metals: arsenic, barium, cadmium, chromium, lead, mercury, selenium, silver, strontium, zinc
 *Cations: aluminum, calcium, iron, magnesium, manganese, potassium, sodium
 *Anions: bromide, sulfate, carbonate alkalinity, bicarbonate alkalinity

Relinquished By: *[Signature]*
 Date/Time: 10/24/18 1300
 Received By: *[Signature]*
 Date/Time: 10/25/18 1445
 Written: *[Signature]*
 Date/Time: 10/25/18 1445
 Date/Time: 10/25/18 1445
 Date/Time: 10/25/18 1445

SGS Sample Receipt Summary

Job Number: LA49126

Client: HYDRO ENV

Project: INDIGO

Date / Time Received: 10/25/2018 2:45:00 PM

Delivery Method: Accutest Courier

Airbill #s: _____

Cooler Temps: #1: (2.5/2.5);

Cooler Security

| | <u>Y or N</u> | | <u>Y or N</u> | |
|---------------------------|--------------------------|-------------------------------------|----------------------|--|
| 1. Custody Seals Present: | <input type="checkbox"/> | <input checked="" type="checkbox"/> | 3. COC Present: | <input checked="" type="checkbox"/> <input type="checkbox"/> |
| 2. Custody Seals Intact: | <input type="checkbox"/> | <input checked="" type="checkbox"/> | 4. Smp Dates/Time OK | <input checked="" type="checkbox"/> <input type="checkbox"/> |

Cooler Temperature

| | <u>Y or N</u> | |
|------------------------------|-------------------------------------|--------------------------|
| 1. Temp criteria achieved: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. Cooler temp verification: | <u>IR Gun</u> | |
| 3. Cooler media: | <u>Ice (direct contact)</u> | |
| 4. No. Coolers: | <u>1</u> | |

Quality Control Preservation

| | <u>Y</u> | <u>or</u> | <u>N</u> | <u>N/A</u> |
|---------------------------------|-------------------------------------|-----------|--------------------------|--------------------------|
| 1. Trip Blank present / cooler: | <input checked="" type="checkbox"/> | | <input type="checkbox"/> | <input type="checkbox"/> |
| 2. Trip Blank listed on COC: | <input checked="" type="checkbox"/> | | <input type="checkbox"/> | <input type="checkbox"/> |
| 3. Samples preserved properly: | <input checked="" type="checkbox"/> | | <input type="checkbox"/> | |
| 4. VOCs headspace free: | <input checked="" type="checkbox"/> | | <input type="checkbox"/> | <input type="checkbox"/> |

Sample Integrity - Documentation

| | <u>Y or N</u> | |
|--|-------------------------------------|--------------------------|
| 1. Sample labels present on bottles: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. Container labeling complete: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 3. Sample container label / COC agree: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |

Sample Integrity - Condition

| | <u>Y or N</u> | |
|----------------------------------|-------------------------------------|--------------------------|
| 1. Sample recvd within HT: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. All containers accounted for: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 3. Condition of sample: | <u>Intact</u> | |

Sample Integrity - Instructions

| | <u>Y</u> | <u>or</u> | <u>N</u> | <u>N/A</u> |
|---|-------------------------------------|-----------|-------------------------------------|-------------------------------------|
| 1. Analysis requested is clear: | <input checked="" type="checkbox"/> | | <input type="checkbox"/> | |
| 2. Bottles received for unspecified tests | <input type="checkbox"/> | | <input checked="" type="checkbox"/> | |
| 3. Sufficient volume recvd for analysis: | <input checked="" type="checkbox"/> | | <input type="checkbox"/> | |
| 4. Compositing instructions clear: | <input type="checkbox"/> | | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| 5. Filtering instructions clear: | <input type="checkbox"/> | | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

Comments: Field Blank and Trip Blank, rec'd 6 vials per ID, coc states 8.

Responded to by: RF

Response Date: 10/26/2018

No issue

LA49126: Chain of Custody

Page 2 of 2

MS Volatiles

QC Data Summaries

Includes the following where applicable:

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries

Method Blank Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|-------------|------------|----|----------|----|-----------|------------|------------------|
| V111896-MB2 | 11040653.D | 1 | 10/29/18 | LS | n/a | n/a | V111896 |

The QC reported here applies to the following samples:

Method: SW846 8260B

LA49126-3

| CAS No. | Compound | Result | RL | Units | Q |
|------------|-----------------------------|--------|------|-------|---|
| 67-64-1 | Acetone | ND | 50 | ug/l | |
| 71-43-2 | Benzene | ND | 1.0 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | ug/l | |
| 75-15-0 | Carbon Disulfide | ND | 1.0 | ug/l | |
| 56-23-5 | Carbon Tetrachloride | ND | 1.0 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | ug/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 1.0 | ug/l | |
| 541-73-1 | m-Dichlorobenzene | ND | 1.0 | ug/l | |
| 95-50-1 | o-Dichlorobenzene | ND | 1.0 | ug/l | |
| 106-46-7 | p-Dichlorobenzene | ND | 1.0 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | ug/l | |
| 75-35-4 | 1,1-Dichloroethylene | ND | 1.0 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethylene | ND | 1.0 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethylene | ND | 1.0 | ug/l | |
| 540-59-0 | 1,2-Dichloroethene (total) | ND | 1.0 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | ug/l | |
| 542-75-6 | 1,3-Dichloropropene (total) | ND | 1.0 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | ug/l | |
| 67-72-1 | Hexachloroethane | ND | 1.0 | ug/l | |
| 78-83-1 | Isobutyl Alcohol | ND | 100 | ug/l | |
| 74-83-9 | Methyl Bromide | ND | 1.0 | ug/l | |
| 74-87-3 | Methyl Chloride | ND | 1.0 | ug/l | |
| 75-09-2 | Methylene Chloride | ND | 1.0 | ug/l | |
| 78-93-3 | Methyl Ethyl Ketone | ND | 13 | ug/l | |
| 108-10-1 | 4-Methyl-2-pentanone | ND | 13 | ug/l | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.0 | ug/l | |
| 100-42-5 | Styrene | ND | 1.0 | ug/l | |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | ND | 1.0 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 0.50 | ug/l | |

4.1.1
4

Method Blank Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|-------------|------------|----|----------|----|-----------|------------|------------------|
| V111896-MB2 | 11040653.D | 1 | 10/29/18 | LS | n/a | n/a | V111896 |

The QC reported here applies to the following samples:

Method: SW846 8260B

LA49126-3

| CAS No. | Compound | Result | RL | Units | Q |
|-----------|------------------------|--------|-----|-------|---|
| 127-18-4 | Tetrachloroethylene | ND | 1.0 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | ug/l | |
| 79-01-6 | Trichloroethylene | ND | 1.0 | ug/l | |
| 75-69-4 | Trichlorofluoromethane | ND | 1.0 | ug/l | |
| 75-01-4 | Vinyl Chloride | ND | 1.0 | ug/l | |
| | m,p-Xylene | ND | 2.0 | ug/l | |
| 95-47-6 | o-Xylene | ND | 1.0 | ug/l | |
| 1330-20-7 | Xylene (total) | ND | 2.0 | ug/l | |

| CAS No. | Surrogate Recoveries | Limits | |
|------------|-----------------------|--------|---------|
| 17060-07-0 | 1,2-Dichloroethane-D4 | 101% | 84-124% |
| 2037-26-5 | Toluene-D8 | 99% | 83-115% |
| 460-00-4 | 4-Bromofluorobenzene | 94% | 89-111% |

4.1.1
4

Method Blank Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|-------------|------------|----|----------|----|-----------|------------|------------------|
| V2G4705-MB2 | 2G090283.D | 1 | 10/31/18 | LS | n/a | n/a | V2G4705 |

The QC reported here applies to the following samples:

Method: SW846 8260B

LA49126-4

| CAS No. | Compound | Result | RL | Units | Q |
|------------|-----------------------------|--------|------|-------|---|
| 67-64-1 | Acetone | ND | 50 | ug/l | |
| 71-43-2 | Benzene | ND | 1.0 | ug/l | |
| 75-27-4 | Bromodichloromethane | ND | 1.0 | ug/l | |
| 75-25-2 | Bromoform | ND | 1.0 | ug/l | |
| 75-15-0 | Carbon Disulfide | ND | 1.0 | ug/l | |
| 56-23-5 | Carbon Tetrachloride | ND | 1.0 | ug/l | |
| 108-90-7 | Chlorobenzene | ND | 1.0 | ug/l | |
| 75-00-3 | Chloroethane | ND | 1.0 | ug/l | |
| 67-66-3 | Chloroform | ND | 1.0 | ug/l | |
| 124-48-1 | Dibromochloromethane | ND | 1.0 | ug/l | |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 1.0 | ug/l | |
| 541-73-1 | m-Dichlorobenzene | ND | 1.0 | ug/l | |
| 95-50-1 | o-Dichlorobenzene | ND | 1.0 | ug/l | |
| 106-46-7 | p-Dichlorobenzene | ND | 1.0 | ug/l | |
| 75-34-3 | 1,1-Dichloroethane | ND | 1.0 | ug/l | |
| 107-06-2 | 1,2-Dichloroethane | ND | 1.0 | ug/l | |
| 75-35-4 | 1,1-Dichloroethylene | ND | 1.0 | ug/l | |
| 156-59-2 | cis-1,2-Dichloroethylene | ND | 1.0 | ug/l | |
| 156-60-5 | trans-1,2-Dichloroethylene | ND | 1.0 | ug/l | |
| 540-59-0 | 1,2-Dichloroethene (total) | ND | 1.0 | ug/l | |
| 78-87-5 | 1,2-Dichloropropane | ND | 1.0 | ug/l | |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 1.0 | ug/l | |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 1.0 | ug/l | |
| 542-75-6 | 1,3-Dichloropropene (total) | ND | 1.0 | ug/l | |
| 100-41-4 | Ethylbenzene | ND | 1.0 | ug/l | |
| 67-72-1 | Hexachloroethane | ND | 1.0 | ug/l | |
| 78-83-1 | Isobutyl Alcohol | ND | 100 | ug/l | |
| 74-83-9 | Methyl Bromide | ND | 1.0 | ug/l | |
| 74-87-3 | Methyl Chloride | ND | 1.0 | ug/l | |
| 75-09-2 | Methylene Chloride | 0.48 | 1.0 | ug/l | J |
| 78-93-3 | Methyl Ethyl Ketone | ND | 13 | ug/l | |
| 108-10-1 | 4-Methyl-2-pentanone | ND | 13 | ug/l | |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 1.0 | ug/l | |
| 100-42-5 | Styrene | ND | 1.0 | ug/l | |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | ND | 1.0 | ug/l | |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 0.50 | ug/l | |

4.1.2
4

Method Blank Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|-------------|------------|----|----------|----|-----------|------------|------------------|
| V2G4705-MB2 | 2G090283.D | 1 | 10/31/18 | LS | n/a | n/a | V2G4705 |

The QC reported here applies to the following samples:

Method: SW846 8260B

LA49126-4

| CAS No. | Compound | Result | RL | Units | Q |
|-----------|------------------------|--------|-----|-------|---|
| 127-18-4 | Tetrachloroethylene | ND | 1.0 | ug/l | |
| 108-88-3 | Toluene | ND | 1.0 | ug/l | |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 1.0 | ug/l | |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 1.0 | ug/l | |
| 79-01-6 | Trichloroethylene | ND | 1.0 | ug/l | |
| 75-69-4 | Trichlorofluoromethane | ND | 1.0 | ug/l | |
| 75-01-4 | Vinyl Chloride | ND | 1.0 | ug/l | |
| | m,p-Xylene | ND | 2.0 | ug/l | |
| 95-47-6 | o-Xylene | ND | 1.0 | ug/l | |
| 1330-20-7 | Xylene (total) | ND | 2.0 | ug/l | |

| CAS No. | Surrogate Recoveries | Limits | |
|------------|-----------------------|--------|---------|
| 17060-07-0 | 1,2-Dichloroethane-D4 | 101% | 84-124% |
| 2037-26-5 | Toluene-D8 | 97% | 83-115% |
| 460-00-4 | 4-Bromofluorobenzene | 102% | 89-111% |

4.1.2
4

Blank Spike/Blank Spike Duplicate Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------------|------------|----|----------|----|-----------|------------|------------------|
| V111896-BS1 | 11040647.D | 1 | 10/29/18 | LS | n/a | n/a | V111896 |
| V111896-BSD1 | 11040649.D | 1 | 10/29/18 | LS | n/a | n/a | V111896 |

The QC reported here applies to the following samples:

Method: SW846 8260B

LA49126-3

| CAS No. | Compound | Spike ug/l | BSP ug/l | BSP % | BSD ug/l | BSD % | RPD | Limits Rec/RPD |
|------------|-----------------------------|---------------|-------------|----------|-------------|----------|-----|------------------------|
| 67-64-1 | Acetone | 50 | 43.8 | 88 | 44.2 | 88 | 1 | 38-178/30 |
| 71-43-2 | Benzene | 20 | 19.1 | 96 | 19.6 | 98 | 3 | 82-119/30 |
| 75-27-4 | Bromodichloromethane | 20 | 18.4 | 92 | 19.1 | 96 | 4 | 79-120/30 |
| 75-25-2 | Bromoform | 20 | 15.5 | 78 | 15.9 | 80 | 3 | 68-128/30 |
| 75-15-0 | Carbon Disulfide | 20 | 18.3 | 92 | 19.2 | 96 | 5 | 64-133/30 |
| 56-23-5 | Carbon Tetrachloride | 20 | 17.7 | 89 | 18.3 | 92 | 3 | 69-132/30 |
| 108-90-7 | Chlorobenzene | 20 | 18.5 | 93 | 19.0 | 95 | 3 | 85-120/30 |
| 75-00-3 | Chloroethane | 20 | 18.6 | 93 | 19.2 | 96 | 3 | 33-170/30 |
| 67-66-3 | Chloroform | 20 | 17.8 | 89 | 18.6 | 93 | 4 | 80-122/30 |
| 124-48-1 | Dibromochloromethane | 20 | 17.3 | 87 | 17.7 | 89 | 2 | 73-125/30 |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | 20 | 13.6 | 68 | 13.8 | 69 | 1 | 67-131/30 |
| 541-73-1 | m-Dichlorobenzene | 20 | 19.4 | 97 | 19.8 | 99 | 2 | 84-121/30 |
| 95-50-1 | o-Dichlorobenzene | 20 | 19.6 | 98 | 20.2 | 101 | 3 | 83-120/30 |
| 106-46-7 | p-Dichlorobenzene | 20 | 18.9 | 95 | 18.7 | 94 | 1 | 83-122/30 |
| 75-34-3 | 1,1-Dichloroethane | 20 | 17.7 | 89 | 18.6 | 93 | 5 | 78-124/30 |
| 107-06-2 | 1,2-Dichloroethane | 20 | 18.5 | 93 | 19.1 | 96 | 3 | 74-127/30 |
| 75-35-4 | 1,1-Dichloroethylene | 20 | 17.0 | 85 | 17.7 | 89 | 4 | 70-134/30 |
| 156-59-2 | cis-1,2-Dichloroethylene | 20 | 19.0 | 95 | 19.8 | 99 | 4 | 78-122/30 |
| 156-60-5 | trans-1,2-Dichloroethylene | 20 | 17.0 | 85 | 17.6 | 88 | 3 | 75-127/30 |
| 540-59-0 | 1,2-Dichloroethene (total) | 40 | 36.1 | 90 | 37.4 | 94 | 4 | 78-123/30 |
| 78-87-5 | 1,2-Dichloropropane | 20 | 18.9 | 95 | 19.4 | 97 | 3 | 82-120/30 |
| 10061-01-5 | cis-1,3-Dichloropropene | 20 | 17.4 | 87 | 17.7 | 89 | 2 | 79-122/30 |
| 10061-02-6 | trans-1,3-Dichloropropene | 20 | 17.2 | 86 | 17.7 | 89 | 3 | 78-124/30 |
| 542-75-6 | 1,3-Dichloropropene (total) | 40 | 34.6 | 87 | 35.4 | 89 | 2 | 50-150/30 ^a |
| 100-41-4 | Ethylbenzene | 20 | 19.6 | 98 | 20.1 | 101 | 3 | 84-117/30 |
| 67-72-1 | Hexachloroethane | 20 | 14.9 | 75 | 15.3 | 77 | 3 | 53-141/30 |
| 78-83-1 | Isobutyl Alcohol | 200 | 180 | 90 | 182 | 91 | 1 | 20-175/30 |
| 74-83-9 | Methyl Bromide | 20 | 19.1 | 96 | 19.4 | 97 | 2 | 37-198/30 |
| 74-87-3 | Methyl Chloride | 20 | 15.7 | 79 | 16.5 | 83 | 5 | 50-136/30 |
| 75-09-2 | Methylene Chloride | 20 | 18.7 | 94 | 19.4 | 97 | 4 | 71-130/30 |
| 78-93-3 | Methyl Ethyl Ketone | 50 | 48.7 | 97 | 50.2 | 100 | 3 | 59-149/30 |
| 108-10-1 | 4-Methyl-2-pentanone | 50 | 42.0 | 84 | 43.4 | 87 | 3 | 74-131/30 |
| 1634-04-4 | Methyl Tert Butyl Ether | 20 | 19.5 | 98 | 19.2 | 96 | 2 | 70-126/30 |
| 100-42-5 | Styrene | 20 | 18.7 | 94 | 19.1 | 96 | 2 | 79-128/30 |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 20 | 19.1 | 96 | 19.9 | 100 | 4 | 84-120/30 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 20 | 19.0 | 95 | 19.2 | 96 | 1 | 77-126/30 |

* = Outside of Control Limits.

Blank Spike/Blank Spike Duplicate Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------------|------------|----|----------|----|-----------|------------|------------------|
| V1I1896-BS1 | 1I040647.D | 1 | 10/29/18 | LS | n/a | n/a | V1I1896 |
| V1I1896-BSD1 | 1I040649.D | 1 | 10/29/18 | LS | n/a | n/a | V1I1896 |

The QC reported here applies to the following samples:

Method: SW846 8260B

LA49126-3

| CAS No. | Compound | Spike ug/l | BSP ug/l | BSP % | BSD ug/l | BSD % | RPD | Limits Rec/RPD |
|-----------|------------------------|---------------|-------------|----------|-------------|----------|-----|-------------------|
| 127-18-4 | Tetrachloroethylene | 20 | 18.3 | 92 | 19.1 | 96 | 4 | 75-133/30 |
| 108-88-3 | Toluene | 20 | 18.4 | 92 | 19.0 | 95 | 3 | 80-121/30 |
| 71-55-6 | 1,1,1-Trichloroethane | 20 | 18.1 | 91 | 18.8 | 94 | 4 | 74-126/30 |
| 79-00-5 | 1,1,2-Trichloroethane | 20 | 18.9 | 95 | 19.5 | 98 | 3 | 80-123/30 |
| 79-01-6 | Trichloroethylene | 20 | 19.3 | 97 | 19.6 | 98 | 2 | 62-125/30 |
| 75-69-4 | Trichlorofluoromethane | 20 | 15.8 | 79 | 16.6 | 83 | 5 | 62-148/30 |
| 75-01-4 | Vinyl Chloride | 20 | 17.2 | 86 | 17.9 | 90 | 4 | 67-130/30 |
| | m,p-Xylene | 40 | 41.4 | 104 | 42.6 | 107 | 3 | 82-121/30 |
| 95-47-6 | o-Xylene | 20 | 18.2 | 91 | 18.8 | 94 | 3 | 84-119/30 |
| 1330-20-7 | Xylene (total) | 60 | 59.5 | 99 | 61.4 | 102 | 3 | 81-122/30 |

| CAS No. | Surrogate Recoveries | BSP | BSD | Limits |
|------------|-----------------------|------|------|---------|
| 17060-07-0 | 1,2-Dichloroethane-D4 | 90% | 91% | 84-124% |
| 2037-26-5 | Toluene-D8 | 101% | 101% | 83-115% |
| 460-00-4 | 4-Bromofluorobenzene | 100% | 101% | 89-111% |

(a) Advisory control limits.

* = Outside of Control Limits.

4.2.1
4

Blank Spike/Blank Spike Duplicate Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------------|------------|----|----------|----|-----------|------------|------------------|
| V2G4705-BS1 | 2G090277.D | 1 | 10/31/18 | LS | n/a | n/a | V2G4705 |
| V2G4705-BSD1 | 2G090279.D | 1 | 10/31/18 | LS | n/a | n/a | V2G4705 |

The QC reported here applies to the following samples:

Method: SW846 8260B

LA49126-4

| CAS No. | Compound | Spike ug/l | BSP ug/l | BSP % | BSD ug/l | BSD % | RPD | Limits Rec/RPD |
|------------|-----------------------------|---------------|-------------|----------|-------------|----------|-----|------------------------|
| 67-64-1 | Acetone | 50 | 84.1 | 168 | 82.0 | 164 | 3 | 38-178/30 |
| 71-43-2 | Benzene | 20 | 18.7 | 94 | 19.5 | 98 | 4 | 82-119/30 |
| 75-27-4 | Bromodichloromethane | 20 | 21.7 | 109 | 22.2 | 111 | 2 | 79-120/30 |
| 75-25-2 | Bromoform | 20 | 22.9 | 115 | 22.5 | 113 | 2 | 68-128/30 |
| 75-15-0 | Carbon Disulfide | 20 | 18.9 | 95 | 19.6 | 98 | 4 | 64-133/30 |
| 56-23-5 | Carbon Tetrachloride | 20 | 20.8 | 104 | 21.4 | 107 | 3 | 69-132/30 |
| 108-90-7 | Chlorobenzene | 20 | 19.4 | 97 | 20.6 | 103 | 6 | 85-120/30 |
| 75-00-3 | Chloroethane | 20 | 19.7 | 99 | 18.6 | 93 | 6 | 33-170/30 |
| 67-66-3 | Chloroform | 20 | 19.8 | 99 | 20.2 | 101 | 2 | 80-122/30 |
| 124-48-1 | Dibromochloromethane | 20 | 22.1 | 111 | 22.6 | 113 | 2 | 73-125/30 |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | 20 | 22.3 | 112 | 22.2 | 111 | 0 | 67-131/30 |
| 541-73-1 | m-Dichlorobenzene | 20 | 19.7 | 99 | 20.1 | 101 | 2 | 84-121/30 |
| 95-50-1 | o-Dichlorobenzene | 20 | 20.2 | 101 | 20.5 | 103 | 1 | 83-120/30 |
| 106-46-7 | p-Dichlorobenzene | 20 | 19.8 | 99 | 19.8 | 99 | 0 | 83-122/30 |
| 75-34-3 | 1,1-Dichloroethane | 20 | 19.1 | 96 | 19.1 | 96 | 0 | 78-124/30 |
| 107-06-2 | 1,2-Dichloroethane | 20 | 20.2 | 101 | 19.9 | 100 | 1 | 74-127/30 |
| 75-35-4 | 1,1-Dichloroethylene | 20 | 18.3 | 92 | 18.5 | 93 | 1 | 70-134/30 |
| 156-59-2 | cis-1,2-Dichloroethylene | 20 | 18.1 | 91 | 18.1 | 91 | 0 | 78-122/30 |
| 156-60-5 | trans-1,2-Dichloroethylene | 20 | 17.7 | 89 | 18.2 | 91 | 3 | 75-127/30 |
| 540-59-0 | 1,2-Dichloroethene (total) | 40 | 35.9 | 90 | 36.2 | 91 | 1 | 78-123/30 |
| 78-87-5 | 1,2-Dichloropropane | 20 | 18.3 | 92 | 18.4 | 92 | 1 | 82-120/30 |
| 10061-01-5 | cis-1,3-Dichloropropene | 20 | 20.7 | 104 | 20.7 | 104 | 0 | 79-122/30 |
| 10061-02-6 | trans-1,3-Dichloropropene | 20 | 21.5 | 108 | 21.9 | 110 | 2 | 78-124/30 |
| 542-75-6 | 1,3-Dichloropropene (total) | 40 | 42.3 | 106 | 42.6 | 107 | 1 | 50-150/30 ^a |
| 100-41-4 | Ethylbenzene | 20 | 19.6 | 98 | 20.3 | 102 | 4 | 84-117/30 |
| 67-72-1 | Hexachloroethane | 20 | 22.6 | 113 | 23.0 | 115 | 2 | 53-141/30 |
| 78-83-1 | Isobutyl Alcohol | 200 | 196 | 98 | 189 | 95 | 4 | 20-175/30 |
| 74-83-9 | Methyl Bromide | 20 | 17.4 | 87 | 20.7 | 104 | 17 | 37-198/30 |
| 74-87-3 | Methyl Chloride | 20 | 15.3 | 77 | 15.3 | 77 | 0 | 50-136/30 |
| 75-09-2 | Methylene Chloride | 20 | 18.5 | 93 | 19.6 | 98 | 6 | 71-130/30 |
| 78-93-3 | Methyl Ethyl Ketone | 50 | 62.1 | 124 | 63.4 | 127 | 2 | 59-149/30 |
| 108-10-1 | 4-Methyl-2-pentanone | 50 | 55.5 | 111 | 55.6 | 111 | 0 | 74-131/30 |
| 1634-04-4 | Methyl Tert Butyl Ether | 20 | 19.7 | 99 | 19.6 | 98 | 1 | 70-126/30 |
| 100-42-5 | Styrene | 20 | 20.2 | 101 | 20.9 | 105 | 3 | 79-128/30 |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | 20 | 22.3 | 112 | 22.4 | 112 | 0 | 84-120/30 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | 20 | 18.8 | 94 | 19.4 | 97 | 3 | 77-126/30 |

* = Outside of Control Limits.

4.2.2
4

Blank Spike/Blank Spike Duplicate Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------------|------------|----|----------|----|-----------|------------|------------------|
| V2G4705-BS1 | 2G090277.D | 1 | 10/31/18 | LS | n/a | n/a | V2G4705 |
| V2G4705-BSD1 | 2G090279.D | 1 | 10/31/18 | LS | n/a | n/a | V2G4705 |

The QC reported here applies to the following samples:

Method: SW846 8260B

LA49126-4

| CAS No. | Compound | Spike ug/l | BSP ug/l | BSP % | BSD ug/l | BSD % | RPD | Limits Rec/RPD |
|-----------|------------------------|---------------|-------------|----------|-------------|----------|-----|-------------------|
| 127-18-4 | Tetrachloroethylene | 20 | 20.6 | 103 | 22.0 | 110 | 7 | 75-133/30 |
| 108-88-3 | Toluene | 20 | 19.1 | 96 | 19.9 | 100 | 4 | 80-121/30 |
| 71-55-6 | 1,1,1-Trichloroethane | 20 | 20.7 | 104 | 21.5 | 108 | 4 | 74-126/30 |
| 79-00-5 | 1,1,2-Trichloroethane | 20 | 20.4 | 102 | 20.0 | 100 | 2 | 80-123/30 |
| 79-01-6 | Trichloroethylene | 20 | 20.4 | 102 | 21.5 | 108 | 5 | 62-125/30 |
| 75-69-4 | Trichlorofluoromethane | 20 | 22.3 | 112 | 22.3 | 112 | 0 | 62-148/30 |
| 75-01-4 | Vinyl Chloride | 20 | 18.6 | 93 | 19.0 | 95 | 2 | 67-130/30 |
| | m,p-Xylene | 40 | 40.4 | 101 | 42.3 | 106 | 5 | 82-121/30 |
| 95-47-6 | o-Xylene | 20 | 21.2 | 106 | 21.9 | 110 | 3 | 84-119/30 |
| 1330-20-7 | Xylene (total) | 60 | 61.6 | 103 | 64.2 | 107 | 4 | 81-122/30 |

| CAS No. | Surrogate Recoveries | BSP | BSD | Limits |
|------------|-----------------------|------|------|---------|
| 17060-07-0 | 1,2-Dichloroethane-D4 | 101% | 99% | 84-124% |
| 2037-26-5 | Toluene-D8 | 100% | 98% | 83-115% |
| 460-00-4 | 4-Bromofluorobenzene | 102% | 103% | 89-111% |

(a) Advisory control limits.

* = Outside of Control Limits.

4.2.2
4

Matrix Spike/Matrix Spike Duplicate Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------------|------------|----|----------|----|-----------|------------|------------------|
| LA49176-5MS | 1I040713.D | 5 | 10/31/18 | LS | n/a | n/a | V1I1896 |
| LA49176-5MSD | 1I040715.D | 5 | 10/31/18 | LS | n/a | n/a | V1I1896 |
| LA49176-5 | 1I040659.D | 1 | 10/29/18 | LS | n/a | n/a | V1I1896 |

The QC reported here applies to the following samples:

Method: SW846 8260B

LA49126-3

| CAS No. | Compound | LA49176-5 ug/l | Spike Q | MS ug/l | MS % | Spike ug/l | MSD ug/l | MSD % | RPD | Limits Rec/RPD |
|------------|-----------------------------|-------------------|------------|------------|---------|---------------|-------------|----------|-----|------------------------|
| 67-64-1 | Acetone | ND | 250 | 172 | 69 | 250 | 179 | 72 | 4 | 39-164/27 |
| 71-43-2 | Benzene | ND | 100 | 99.6 | 100 | 100 | 104 | 104 | 4 | 31-161/15 |
| 75-27-4 | Bromodichloromethane | ND | 100 | 102 | 102 | 100 | 105 | 105 | 3 | 64-122/36 |
| 75-25-2 | Bromoform | ND | 100 | 78.0 | 78 | 100 | 82.3 | 82 | 5 | 43-125/37 |
| 75-15-0 | Carbon Disulfide | ND | 100 | 98.9 | 99 | 100 | 103 | 103 | 4 | 38-138/36 |
| 56-23-5 | Carbon Tetrachloride | ND | 100 | 97.8 | 98 | 100 | 102 | 102 | 4 | 53-133/36 |
| 108-90-7 | Chlorobenzene | ND | 100 | 97.2 | 97 | 100 | 98.8 | 99 | 2 | 74-122/34 |
| 75-00-3 | Chloroethane | ND | 100 | 118 | 118 | 100 | 119 | 119 | 1 | 14-181/43 |
| 67-66-3 | Chloroform | ND | 100 | 97.9 | 98 | 100 | 101 | 101 | 3 | 65-130/24 |
| 124-48-1 | Dibromochloromethane | ND | 100 | 88.1 | 88 | 100 | 93.1 | 93 | 6 | 57-121/36 |
| 96-12-8 | 1,2-Dibromo-3-chloropropane | ND | 100 | 63.2 | 63 | 100 | 71.1 | 71 | 12 | 46-135/25 |
| 541-73-1 | m-Dichlorobenzene | ND | 100 | 101 | 101 | 100 | 102 | 102 | 1 | 70-120/35 |
| 95-50-1 | o-Dichlorobenzene | ND | 100 | 98.6 | 99 | 100 | 102 | 102 | 3 | 72-120/35 |
| 106-46-7 | p-Dichlorobenzene | ND | 100 | 95.3 | 95 | 100 | 97.8 | 98 | 3 | 68-120/35 |
| 75-34-3 | 1,1-Dichloroethane | ND | 100 | 98.0 | 98 | 100 | 101 | 101 | 3 | 56-138/32 |
| 107-06-2 | 1,2-Dichloroethane | ND | 100 | 108 | 108 | 100 | 111 | 111 | 3 | 51-141/39 |
| 75-35-4 | 1,1-Dichloroethylene | ND | 100 | 92.9 | 93 | 100 | 95.6 | 96 | 3 | 48-139/37 |
| 156-59-2 | cis-1,2-Dichloroethylene | ND | 100 | 98.1 | 98 | 100 | 101 | 101 | 3 | 56-133/15 |
| 156-60-5 | trans-1,2-Dichloroethylene | ND | 100 | 95.5 | 96 | 100 | 96.1 | 96 | 1 | 59-128/37 |
| 540-59-0 | 1,2-Dichloroethane (total) | ND | 200 | 194 | 97 | 200 | 197 | 99 | 2 | 54-134/30 |
| 78-87-5 | 1,2-Dichloropropane | ND | 100 | 98.6 | 99 | 100 | 101 | 101 | 2 | 68-124/32 |
| 10061-01-5 | cis-1,3-Dichloropropene | ND | 100 | 84.2 | 84 | 100 | 87.1 | 87 | 3 | 62-120/35 |
| 10061-02-6 | trans-1,3-Dichloropropene | ND | 100 | 90.8 | 91 | 100 | 94.6 | 95 | 4 | 64-119/36 |
| 542-75-6 | 1,3-Dichloropropene (total) | ND | 200 | 175 | 88 | 200 | 182 | 91 | 4 | 50-150/30 ^a |
| 100-41-4 | Ethylbenzene | ND | 100 | 105 | 105 | 100 | 107 | 107 | 2 | 47-146/30 |
| 67-72-1 | Hexachloroethane | ND | 100 | 69.1 | 69 | 100 | 71.5 | 72 | 3 | 32-128/39 |
| 78-83-1 | Isobutyl Alcohol | ND | 1000 | 855 | 86 | 1000 | 885 | 89 | 3 | 33-142/54 |
| 74-83-9 | Methyl Bromide | ND | 100 | 104 | 104 | 100 | 107 | 107 | 3 | 1-150/64 |
| 74-87-3 | Methyl Chloride | ND | 100 | 90.4 | 90 | 100 | 93.4 | 93 | 3 | 16-146/29 |
| 75-09-2 | Methylene Chloride | ND | 100 | 115 | 115 | 100 | 116 | 116 | 1 | 55-134/36 |
| 78-93-3 | Methyl Ethyl Ketone | ND | 250 | 230 | 92 | 250 | 235 | 94 | 2 | 54-142/39 |
| 108-10-1 | 4-Methyl-2-pentanone | ND | 250 | 233 | 93 | 250 | 236 | 94 | 1 | 60-140/40 |
| 1634-04-4 | Methyl Tert Butyl Ether | ND | 100 | 99.0 | 99 | 100 | 99.6 | 100 | 1 | 52-146/32 |
| 100-42-5 | Styrene | ND | 100 | 96.3 | 96 | 100 | 98.8 | 99 | 3 | 67-128/35 |
| 630-20-6 | 1,1,1,2-Tetrachloroethane | ND | 100 | 104 | 104 | 100 | 106 | 106 | 2 | 67-121/35 |
| 79-34-5 | 1,1,2,2-Tetrachloroethane | ND | 100 | 104 | 104 | 100 | 103 | 103 | 1 | 64-133/38 |

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------------|------------|----|----------|----|-----------|------------|------------------|
| LA49176-5MS | 1I040713.D | 5 | 10/31/18 | LS | n/a | n/a | V1I1896 |
| LA49176-5MSD | 1I040715.D | 5 | 10/31/18 | LS | n/a | n/a | V1I1896 |
| LA49176-5 | 1I040659.D | 1 | 10/29/18 | LS | n/a | n/a | V1I1896 |

The QC reported here applies to the following samples:

Method: SW846 8260B

LA49126-3

| CAS No. | Compound | LA49176-5 ug/l | Spike Q ug/l | MS ug/l | MS % | Spike ug/l | MSD ug/l | MSD % | RPD | Limits Rec/RPD |
|-----------|------------------------|-------------------|--------------------|------------|---------|---------------|-------------|----------|-----|-------------------|
| 127-18-4 | Tetrachloroethylene | ND | 100 | 97.4 | 97 | 100 | 101 | 101 | 4 | 58-135/37 |
| 108-88-3 | Toluene | ND | 100 | 97.0 | 97 | 100 | 99.3 | 99 | 2 | 36-155/17 |
| 71-55-6 | 1,1,1-Trichloroethane | ND | 100 | 101 | 101 | 100 | 103 | 103 | 2 | 63-128/36 |
| 79-00-5 | 1,1,2-Trichloroethane | ND | 100 | 102 | 102 | 100 | 102 | 102 | 0 | 61-138/17 |
| 79-01-6 | Trichloroethylene | ND | 100 | 98.9 | 99 | 100 | 103 | 103 | 4 | 57-131/36 |
| 75-69-4 | Trichlorofluoromethane | ND | 100 | 103 | 103 | 100 | 105 | 105 | 2 | 31-156/36 |
| 75-01-4 | Vinyl Chloride | ND | 100 | 95.7 | 96 | 100 | 97.6 | 98 | 2 | 22-155/49 |
| | m,p-Xylene | ND | 200 | 224 | 112 | 200 | 230 | 115 | 3 | 35-159/31 |
| 95-47-6 | o-Xylene | ND | 100 | 93.3 | 93 | 100 | 96.6 | 97 | 3 | 50-144/35 |
| 1330-20-7 | Xylene (total) | ND | 300 | 318 | 106 | 300 | 327 | 109 | 3 | 41-154/29 |

| CAS No. | Surrogate Recoveries | MS | MSD | LA49176-5 | Limits |
|------------|-----------------------|------|------|-----------|---------|
| 17060-07-0 | 1,2-Dichloroethane-D4 | 100% | 100% | 110% | 84-124% |
| 2037-26-5 | Toluene-D8 | 98% | 99% | 99% | 83-115% |
| 460-00-4 | 4-Bromofluorobenzene | 102% | 102% | 94% | 89-111% |

(a) Advisory control limits.

* = Outside of Control Limits.

MS Semi-volatiles

QC Data Summaries

Includes the following where applicable:

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries

Method Blank Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|------------|------------|----|----------|----|-----------|------------|------------------|
| OP12657-MB | C0039827.D | 1 | 10/29/18 | IK | 10/29/18 | OP12657 | EC1678 |

The QC reported here applies to the following samples:

Method: SW846 8270D

LA49126-1, LA49126-2

| CAS No. | Compound | Result | RL | Units | Q |
|----------|------------------------------|--------|------|-------|---|
| 95-57-8 | 2-Chlorophenol | ND | 5.0 | ug/l | |
| 120-83-2 | 2,4-Dichlorophenol | ND | 5.0 | ug/l | |
| 105-67-9 | 2,4-Dimethylphenol | ND | 5.0 | ug/l | |
| 51-28-5 | 2,4-Dinitrophenol | ND | 20 | ug/l | |
| 100-02-7 | 4-Nitrophenol | ND | 25 | ug/l | |
| 87-86-5 | Pentachlorophenol | ND | 10 | ug/l | |
| 108-95-2 | Phenol | ND | 5.0 | ug/l | |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | ND | 5.0 | ug/l | |
| 95-95-4 | 2,4,5-Trichlorophenol | ND | 5.0 | ug/l | |
| 88-06-2 | 2,4,6-Trichlorophenol | ND | 5.0 | ug/l | |
| 83-32-9 | Acenaphthene | ND | 0.20 | ug/l | |
| 208-96-8 | Acenaphthylene | ND | 0.20 | ug/l | |
| 62-53-3 | Aniline | ND | 5.0 | ug/l | |
| 120-12-7 | Anthracene | ND | 0.20 | ug/l | |
| 56-55-3 | Benzo(a)anthracene | ND | 0.20 | ug/l | |
| 50-32-8 | Benzo(a)pyrene | ND | 0.20 | ug/l | |
| 205-99-2 | Benzo(b)fluoranthene | ND | 0.20 | ug/l | |
| 207-08-9 | Benzo(k)fluoranthene | ND | 0.20 | ug/l | |
| 92-52-4 | 1,1'-Biphenyl | ND | 10 | ug/l | |
| 85-68-7 | Butyl Benzyl Phthalate | ND | 5.0 | ug/l | |
| 106-47-8 | 4-Chloroaniline | ND | 5.0 | ug/l | |
| 111-44-4 | bis(2-Chloroethyl)ether | ND | 5.0 | ug/l | |
| 108-60-1 | 2,2'-Oxybis(1-chloropropane) | ND | 5.0 | ug/l | |
| 91-58-7 | 2-Chloronaphthalene | ND | 5.0 | ug/l | |
| 218-01-9 | Chrysene | ND | 0.20 | ug/l | |
| 53-70-3 | Dibenzo(a,h)anthracene | ND | 0.20 | ug/l | |
| 132-64-9 | Dibenzofuran | ND | 5.0 | ug/l | |
| 91-94-1 | 3,3'-Dichlorobenzidine | ND | 10 | ug/l | |
| 84-66-2 | Diethyl Phthalate | ND | 5.0 | ug/l | |
| 131-11-3 | Dimethyl Phthalate | ND | 5.0 | ug/l | |
| 117-84-0 | Di-n-octyl Phthalate | ND | 5.0 | ug/l | |
| 99-65-0 | 1,3-Dinitrobenzene | ND | 5.0 | ug/l | |
| 121-14-2 | 2,4-Dinitrotoluene | ND | 5.0 | ug/l | |
| 606-20-2 | 2,6-Dinitrotoluene | ND | 5.0 | ug/l | |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | ND | 5.0 | ug/l | |
| 206-44-0 | Fluoranthene | ND | 0.20 | ug/l | |

5.1.1
5

Method Blank Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|------------|------------|----|----------|----|-----------|------------|------------------|
| OP12657-MB | C0039827.D | 1 | 10/29/18 | IK | 10/29/18 | OP12657 | EC1678 |

The QC reported here applies to the following samples:

Method: SW846 8270D

LA49126-1, LA49126-2

| CAS No. | Compound | Result | RL | Units | Q |
|----------|----------------------------|--------|------|-------|---|
| 86-73-7 | Fluorene | ND | 0.20 | ug/l | |
| 118-74-1 | Hexachlorobenzene | ND | 5.0 | ug/l | |
| 87-68-3 | Hexachlorobutadiene | ND | 5.0 | ug/l | |
| 77-47-4 | Hexachlorocyclopentadiene | ND | 10 | ug/l | |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | ND | 0.20 | ug/l | |
| 78-59-1 | Isophorone | ND | 5.0 | ug/l | |
| 91-57-6 | 2-Methylnaphthalene | ND | 0.20 | ug/l | |
| 91-20-3 | Naphthalene | ND | 0.20 | ug/l | |
| 88-74-4 | 2-Nitroaniline | ND | 5.0 | ug/l | |
| 99-09-2 | 3-Nitroaniline | ND | 5.0 | ug/l | |
| 100-01-6 | 4-Nitroaniline | ND | 5.0 | ug/l | |
| 98-95-3 | Nitrobenzene | ND | 5.0 | ug/l | |
| 621-64-7 | N-Nitroso-di-n-propylamine | ND | 5.0 | ug/l | |
| 86-30-6 | N-Nitrosodiphenylamine | ND | 5.0 | ug/l | |
| 85-01-8 | Phenanthrene | ND | 0.20 | ug/l | |
| 129-00-0 | Pyrene | ND | 0.20 | ug/l | |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | ND | 5.0 | ug/l | |
| 120-82-1 | 1,2,4-Trichlorobenzene | ND | 5.0 | ug/l | |

| CAS No. | Surrogate Recoveries | Limits | |
|-----------|----------------------|--------|---------|
| 367-12-4 | 2-Fluorophenol | 49% | 23-85% |
| 4165-62-2 | Phenol-d5 | 35% | 10-69% |
| 118-79-6 | 2,4,6-Tribromophenol | 88% | 48-138% |
| 4165-60-0 | Nitrobenzene-d5 | 89% | 51-128% |
| 321-60-8 | 2-Fluorobiphenyl | 93% | 55-122% |
| 1718-51-0 | Terphenyl-d14 | 91% | 43-138% |

Blank Spike/Blank Spike Duplicate Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|-------------|------------|----|----------|----|-----------|------------|------------------|
| OP12657-BS | C0039828.D | 1 | 10/29/18 | IK | 10/29/18 | OP12657 | EC1678 |
| OP12657-BSD | C0039829.D | 1 | 10/29/18 | IK | 10/29/18 | OP12657 | EC1678 |

The QC reported here applies to the following samples:

Method: SW846 8270D

LA49126-1, LA49126-2

| CAS No. | Compound | Spike ug/l | BSP ug/l | BSP % | BSD ug/l | BSD % | RPD | Limits Rec/RPD |
|----------|------------------------------|---------------|-------------|----------|-------------|----------|-------|-------------------|
| 95-57-8 | 2-Chlorophenol | 50 | 38.7 | 77 | 33.7 | 67 | 14 | 63-104/19 |
| 120-83-2 | 2,4-Dichlorophenol | 50 | 44.0 | 88 | 37.8 | 76 | 15 | 68-112/19 |
| 105-67-9 | 2,4-Dimethylphenol | 50 | 44.4 | 89 | 35.6 | 71 | 22* a | 64-110/20 |
| 51-28-5 | 2,4-Dinitrophenol | 50 | 44.8 | 90 | 40.4 | 81 | 10 | 51-121/30 |
| 100-02-7 | 4-Nitrophenol | 50 | 18.4 | 37 | 16.0 | 32 | 14 | 20-68/23 |
| 87-86-5 | Pentachlorophenol | 50 | 43.0 | 86 | 35.0 | 70 | 21 | 52-120/29 |
| 108-95-2 | Phenol | 50 | 18.2 | 36 | 16.6 | 33 | 9 | 18-67/20 |
| 58-90-2 | 2,3,4,6-Tetrachlorophenol | 50 | 52.0 | 104 | 44.4 | 89 | 16 | 67-121/21 |
| 95-95-4 | 2,4,5-Trichlorophenol | 50 | 52.0 | 104 | 44.9 | 90 | 15 | 67-119/21 |
| 88-06-2 | 2,4,6-Trichlorophenol | 50 | 45.6 | 91 | 39.2 | 78 | 15 | 67-120/21 |
| 83-32-9 | Acenaphthene | 50 | 46.0 | 92 | 40.5 | 81 | 13 | 67-114/28 |
| 208-96-8 | Acenaphthylene | 50 | 46.9 | 94 | 38.2 | 76 | 20 | 67-119/26 |
| 62-53-3 | Aniline | 50 | 28.5 | 57 | 27.8 | 56 | 2 | 40-114/40 |
| 120-12-7 | Anthracene | 50 | 42.2 | 84 | 37.1 | 74 | 13 | 68-121/24 |
| 56-55-3 | Benzo(a)anthracene | 50 | 44.1 | 88 | 38.5 | 77 | 14 | 69-113/20 |
| 50-32-8 | Benzo(a)pyrene | 50 | 44.4 | 89 | 38.8 | 78 | 13 | 71-124/22 |
| 205-99-2 | Benzo(b)fluoranthene | 50 | 45.1 | 90 | 39.4 | 79 | 13 | 72-120/22 |
| 207-08-9 | Benzo(k)fluoranthene | 50 | 44.8 | 90 | 40.8 | 82 | 9 | 71-124/21 |
| 92-52-4 | 1,1'-Biphenyl | 50 | 47.0 | 94 | 40.8 | 82 | 14 | 65-122/29 |
| 85-68-7 | Butyl Benzyl Phthalate | 50 | 40.5 | 81 | 37.1 | 74 | 9 | 73-123/21 |
| 106-47-8 | 4-Chloroaniline | 50 | 43.9 | 88 | 40.3 | 81 | 9 | 58-113/51 |
| 111-44-4 | bis(2-Chloroethyl)ether | 50 | 44.2 | 88 | 39.2 | 78 | 12 | 50-118/28 |
| 108-60-1 | 2,2'-Oxybis(1-chloropropane) | 50 | 49.0 | 98 | 44.0 | 88 | 11 | 43-138/21 |
| 91-58-7 | 2-Chloronaphthalene | 50 | 45.9 | 92 | 40.9 | 82 | 12 | 64-114/30 |
| 218-01-9 | Chrysene | 50 | 44.4 | 89 | 39.0 | 78 | 13 | 70-115/20 |
| 53-70-3 | Dibenzo(a,h)anthracene | 50 | 46.8 | 94 | 41.6 | 83 | 12 | 70-124/21 |
| 132-64-9 | Dibenzofuran | 50 | 50.3 | 101 | 44.2 | 88 | 13 | 67-117/27 |
| 91-94-1 | 3,3'-Dichlorobenzidine | 50 | 41.2 | 82 | 36.4 | 73 | 12 | 69-122/38 |
| 84-66-2 | Diethyl Phthalate | 50 | 48.2 | 96 | 42.9 | 86 | 12 | 71-123/21 |
| 131-11-3 | Dimethyl Phthalate | 50 | 48.1 | 96 | 43.4 | 87 | 10 | 69-119/20 |
| 117-84-0 | Di-n-octyl Phthalate | 50 | 40.3 | 81 | 35.6 | 71 | 12 | 66-121/22 |
| 99-65-0 | 1,3-Dinitrobenzene | 50 | 50.4 | 101 | 45.9 | 92 | 9 | 71-122/21 |
| 121-14-2 | 2,4-Dinitrotoluene | 50 | 52.0 | 104 | 47.1 | 94 | 10 | 73-122/21 |
| 606-20-2 | 2,6-Dinitrotoluene | 50 | 49.5 | 99 | 44.2 | 88 | 11 | 72-121/21 |
| 117-81-7 | bis(2-Ethylhexyl)phthalate | 50 | 39.2 | 78 | 34.1 | 68 | 14 | 68-126/21 |
| 206-44-0 | Fluoranthene | 50 | 43.8 | 88 | 38.0 | 76 | 14 | 73-120/21 |

* = Outside of Control Limits.

5.2.1
5

Blank Spike/Blank Spike Duplicate Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|-------------|------------|----|----------|----|-----------|------------|------------------|
| OP12657-BS | C0039828.D | 1 | 10/29/18 | IK | 10/29/18 | OP12657 | EC1678 |
| OP12657-BSD | C0039829.D | 1 | 10/29/18 | IK | 10/29/18 | OP12657 | EC1678 |

The QC reported here applies to the following samples:

Method: SW846 8270D

LA49126-1, LA49126-2

| CAS No. | Compound | Spike ug/l | BSP ug/l | BSP % | BSD ug/l | BSD % | RPD | Limits Rec/RPD |
|----------|----------------------------|---------------|-------------|----------|-------------|----------|-----|-------------------|
| 86-73-7 | Fluorene | 50 | 49.8 | 100 | 43.9 | 88 | 13 | 69-118/25 |
| 118-74-1 | Hexachlorobenzene | 50 | 43.0 | 86 | 38.0 | 76 | 12 | 67-117/23 |
| 87-68-3 | Hexachlorobutadiene | 50 | 41.1 | 82 | 36.6 | 73 | 12 | 42-120/35 |
| 77-47-4 | Hexachlorocyclopentadiene | 50 | 41.0 | 82 | 38.0 | 76 | 8 | 35-123/48 |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 50 | 43.6 | 87 | 38.6 | 77 | 12 | 70-123/21 |
| 78-59-1 | Isophorone | 50 | 52.6 | 105 | 46.3 | 93 | 13 | 70-119/19 |
| 91-57-6 | 2-Methylnaphthalene | 50 | 41.6 | 83 | 36.6 | 73 | 13 | 65-113/27 |
| 91-20-3 | Naphthalene | 50 | 40.9 | 82 | 35.6 | 71 | 14 | 63-114/23 |
| 88-74-4 | 2-Nitroaniline | 50 | 46.7 | 93 | 42.0 | 84 | 11 | 68-125/21 |
| 99-09-2 | 3-Nitroaniline | 50 | 46.7 | 93 | 42.1 | 84 | 10 | 69-117/23 |
| 100-01-6 | 4-Nitroaniline | 50 | 44.7 | 89 | 40.2 | 80 | 11 | 67-122/19 |
| 98-95-3 | Nitrobenzene | 50 | 53.7 | 107 | 47.2 | 94 | 13 | 69-116/21 |
| 621-64-7 | N-Nitroso-di-n-propylamine | 50 | 49.6 | 99 | 43.8 | 88 | 12 | 67-120/20 |
| 86-30-6 | N-Nitrosodiphenylamine | 50 | 39.2 | 78 | 34.1 | 68 | 14 | 67-119/25 |
| 85-01-8 | Phenanthrene | 50 | 43.3 | 87 | 38.4 | 77 | 12 | 70-117/23 |
| 129-00-0 | Pyrene | 50 | 43.7 | 87 | 39.0 | 78 | 11 | 70-119/21 |
| 95-94-3 | 1,2,4,5-Tetrachlorobenzene | 50 | 53.9 | 108 | 46.2 | 92 | 15 | 55-117/35 |
| 120-82-1 | 1,2,4-Trichlorobenzene | 50 | 41.4 | 83 | 36.2 | 72 | 13 | 56-111/30 |

| CAS No. | Surrogate Recoveries | BSP | BSD | Limits |
|-----------|----------------------|-----|-----|---------|
| 367-12-4 | 2-Fluorophenol | 47% | 43% | 23-85% |
| 4165-62-2 | Phenol-d5 | 34% | 32% | 10-69% |
| 118-79-6 | 2,4,6-Tribromophenol | 88% | 77% | 48-138% |
| 4165-60-0 | Nitrobenzene-d5 | 93% | 84% | 51-128% |
| 321-60-8 | 2-Fluorobiphenyl | 93% | 83% | 55-122% |
| 1718-51-0 | Terphenyl-d14 | 88% | 81% | 43-138% |

(a) High RPD. Analyte not detected in associated samples.

* = Outside of Control Limits.

5.2.1
5

GC Volatiles

QC Data Summaries

Includes the following where applicable:

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries

Method Blank Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|-------------|------------|----|----------|----|-----------|------------|------------------|
| GLC1882-MB1 | LC381160.D | 1 | 10/27/18 | SV | n/a | n/a | GLC1882 |

The QC reported here applies to the following samples:

Method: MADEP VPH REV 1.1

LA49126-1, LA49126-2, LA49126-3, LA49126-4

| CAS No. | Compound | Result | RL | Units | Q |
|---------|------------------------------|--------|----|-------|---|
| | Aliphatics C6-C8 (Unadj.) | ND | 30 | ug/l | |
| | Aliphatics > C8-C10 (Unadj.) | ND | 50 | ug/l | |
| | Aromatics > C8-C10 (Unadj.) | ND | 50 | ug/l | |

| CAS No. | Surrogate Recoveries | Limits | |
|----------|----------------------|-------------------|---------|
| 615-59-8 | 2,5-Dibromotoluene | 101% ^a | 70-130% |
| 615-59-8 | 2,5-Dibromotoluene | 100% ^b | 70-130% |

- (a) Recovery from Aromatics fraction.
- (b) Recovery from Aliphatics fraction.

6.1.1
6

Blank Spike/Blank Spike Duplicate Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------------|------------|----|----------|----|-----------|------------|------------------|
| GLC1882-BS1 | LC381158.D | 1 | 10/27/18 | SV | n/a | n/a | GLC1882 |
| GLC1882-BSD1 | LC381159.D | 1 | 10/27/18 | SV | n/a | n/a | GLC1882 |

The QC reported here applies to the following samples:

Method: MADEP VPH REV 1.1

LA49126-1, LA49126-2, LA49126-3, LA49126-4

| CAS No. | Compound | Spike ug/l | BSP ug/l | BSP % | BSD ug/l | BSD % | RPD | Limits Rec/RPD |
|---------|------------------------------|---------------|-------------|----------|-------------|----------|-----|-------------------|
| | Aliphatics C6-C8 (Unadj.) | 150 | 123 | 82 | 132 | 88 | 7 | 70-130/30 |
| | Aliphatics > C8-C10 (Unadj.) | 250 | 214 | 86 | 209 | 84 | 2 | 70-130/30 |
| | Aromatics > C8-C10 (Unadj.) | 250 | 221 | 88 | 221 | 88 | 0 | 70-130/30 |

| CAS No. | Surrogate Recoveries | BSP | BSD | Limits |
|----------|----------------------|-------------------|-------------------|---------|
| 615-59-8 | 2,5-Dibromotoluene | 105% ^a | 103% ^a | 70-130% |
| 615-59-8 | 2,5-Dibromotoluene | 106% ^b | 114% ^b | 70-130% |

(a) Recovery from Aromatics fraction.

(b) Recovery from Aliphatics fraction.

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|--------------|------------|----|----------|----|-----------|------------|------------------|
| LA49116-1MS | LC381164.D | 5 | 10/27/18 | SV | n/a | n/a | GLC1882 |
| LA49116-1MSD | LC381165.D | 5 | 10/27/18 | SV | n/a | n/a | GLC1882 |
| LA49116-1 | LC381161.D | 1 | 10/27/18 | SV | n/a | n/a | GLC1882 |

The QC reported here applies to the following samples:

Method: MADEP VPH REV 1.1

LA49126-1, LA49126-2, LA49126-3, LA49126-4

| CAS No. | Compound | LA49116-1 ug/l | Spike Q ug/l | MS ug/l | MS % | Spike ug/l | MSD ug/l | MSD % | RPD | Limits Rec/RPD |
|---------|------------------------------|-------------------|--------------------|------------|---------|---------------|-------------|----------|-----|-------------------|
| | Aliphatics C6-C8 (Unadj.) | ND | 750 | 769 | 103 | 750 | 721 | 96 | 6 | 70-130/50 |
| | Aliphatics > C8-C10 (Unadj.) | ND | 1250 | 1240 | 99 | 1250 | 1270 | 102 | 2 | 70-130/50 |
| | Aromatics > C8-C10 (Unadj.) | ND | 1250 | 1300 | 104 | 1250 | 1310 | 105 | 1 | 70-130/50 |

| CAS No. | Surrogate Recoveries | MS | MSD | LA49116-1 | Limits |
|----------|----------------------|-------------------|-------------------|-------------------|---------|
| 615-59-8 | 2,5-Dibromotoluene | 105% ^a | 104% ^a | 100% ^a | 70-130% |
| 615-59-8 | 2,5-Dibromotoluene | 105% ^b | 104% ^b | 101% ^b | 70-130% |

(a) Recovery from Aromatics fraction.

(b) Recovery from Aliphatics fraction.

* = Outside of Control Limits.

GC/LC Semi-volatiles

QC Data Summaries

Includes the following where applicable:

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries

Method Blank Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|------------|------------|----|----------|----|-----------|------------|------------------|
| OP12635-MB | X0005563.D | 1 | 10/29/18 | JT | 10/26/18 | OP12635 | GLB1651 |

The QC reported here applies to the following samples:

Method: MADEP EPH REV 1.1

LA49126-1, LA49126-2

| CAS No. | Compound | Result | RL | Units | Q |
|---------|------------------------------|--------|-----|-------|---|
| | Aromatics > C10-C12 (Unadj.) | ND | 140 | ug/l | |
| | Aromatics > C12-C16 (Unadj.) | ND | 140 | ug/l | |
| | Aromatics > C16-C21 (Unadj.) | ND | 140 | ug/l | |
| | Aromatics > C21-C35 (Unadj.) | ND | 140 | ug/l | |

| CAS No. | Surrogate Recoveries | Results | Limits |
|----------|----------------------|---------|---------|
| 84-15-1 | o-Terphenyl | 73% | 40-140% |
| 321-60-8 | 2-Fluorobiphenyl | 73% | 40-140% |

7.1.1
7

Method Blank Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|------------|------------|----|----------|----|-----------|------------|------------------|
| OP12635-MB | Y0005563.D | 1 | 10/29/18 | JT | 10/26/18 | OP12635 | GLB1652 |

The QC reported here applies to the following samples:

Method: MADEP EPH REV 1.1

LA49126-1, LA49126-2

| CAS No. | Compound | Result | RL | Units | Q |
|---------|-------------------------------|--------|-----|-------|---|
| | Aliphatics > C10-C12 (Unadj.) | ND | 140 | ug/l | |
| | Aliphatics > C12-C16 (Unadj.) | ND | 140 | ug/l | |
| | Aliphatics > C16-C35 (Unadj.) | ND | 140 | ug/l | |

| CAS No. | Surrogate Recoveries | Results | Limits |
|-----------|----------------------|---------|---------|
| 3386-33-2 | 1-Chlorooctadecane | 84% | 40-140% |

7.1.2
7

Blank Spike/Blank Spike Duplicate Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|-------------|------------|----|----------|----|-----------|------------|------------------|
| OP12635-BS | X0005564.D | 1 | 10/29/18 | JT | 10/26/18 | OP12635 | GLB1651 |
| OP12635-BSD | X0005565.D | 1 | 10/29/18 | JT | 10/26/18 | OP12635 | GLB1651 |

The QC reported here applies to the following samples:

Method: MADEP EPH REV 1.1

LA49126-1, LA49126-2

| CAS No. | Compound | Spike ug/l | BSP ug/l | BSP % | BSD ug/l | BSD % | RPD | Limits Rec/RPD |
|---------|------------------------------|---------------|-------------|----------|-------------|----------|-----|-------------------|
| | Aromatics > C10-C12 (Unadj.) | 461 | 340 | 74 | 339 | 73 | 0 | 40-140/30 |
| | Aromatics > C12-C16 (Unadj.) | 1380 | 1030 | 74 | 1020 | 73 | 1 | 40-140/30 |
| | Aromatics > C16-C21 (Unadj.) | 2310 | 1860 | 81 | 1900 | 82 | 2 | 40-140/30 |
| | Aromatics > C21-C35 (Unadj.) | 3690 | 3090 | 84 | 3140 | 85 | 2 | 40-140/30 |

| CAS No. | Surrogate Recoveries | BSP | BSD | Limits |
|----------|----------------------|-----|-----|---------|
| 84-15-1 | o-Terphenyl | 75% | 75% | 40-140% |
| 321-60-8 | 2-Fluorobiphenyl | 76% | 76% | 40-140% |

* = Outside of Control Limits.

7.2.1
7

Blank Spike/Blank Spike Duplicate Summary

Job Number: LA49126
 Account: HETILAL Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

| Sample | File ID | DF | Analyzed | By | Prep Date | Prep Batch | Analytical Batch |
|-------------|------------|----|----------|----|-----------|------------|------------------|
| OP12635-BS | Y0005564.D | 1 | 10/29/18 | JT | 10/26/18 | OP12635 | GLB1652 |
| OP12635-BSD | Y0005565.D | 1 | 10/29/18 | JT | 10/26/18 | OP12635 | GLB1652 |

The QC reported here applies to the following samples:

Method: MADEP EPH REV 1.1

LA49126-1, LA49126-2

| CAS No. | Compound | Spike ug/l | BSP ug/l | BSP % | BSD ug/l | BSD % | RPD | Limits Rec/RPD |
|---------|-------------------------------|---------------|-------------|----------|-------------|----------|-----|-------------------|
| | Aliphatics > C10-C12 (Unadj.) | 461 | 351 | 76 | 369 | 80 | 5 | 40-140/30 |
| | Aliphatics > C12-C16 (Unadj.) | 923 | 699 | 76 | 741 | 80 | 6 | 40-140/30 |
| | Aliphatics > C16-C35 (Unadj.) | 4150 | 2980 | 72 | 3130 | 75 | 5 | 40-140/30 |

| CAS No. | Surrogate Recoveries | BSP | BSD | Limits |
|-----------|----------------------|-----|-----|---------|
| 3386-33-2 | 1-Chlorooctadecane | 79% | 82% | 40-140% |

* = Outside of Control Limits.

7.2.2
7

Metals Analysis

QC Data Summaries



Includes the following where applicable:

- Method Blank Summaries
- Matrix Spike and Duplicate Summaries
- Blank Spike and Lab Control Sample Summaries
- Serial Dilution Summaries

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: LA49126
Account: HETILAL - Hydro-Environmental Technology, Inc.
Project: 8060.00 Indigo-Desoto Parish, LA

QC Batch ID: MP13152
Matrix Type: AQUEOUS

Methods: SW846 6020A
Units: ug/l

Prep Date: 10/26/18

| Metal | RL | IDL | MDL | MB raw | final |
|------------|------|-------|-----|-----------|-------|
| Aluminum | 100 | 6.9 | 9.3 | -0.76 | <100 |
| Antimony | 1.0 | .043 | .34 | | |
| Arsenic | 1.0 | .062 | .26 | -0.023 | <1.0 |
| Barium | 1.0 | .033 | .46 | -0.019 | <1.0 |
| Beryllium | 1.0 | .0077 | .28 | | |
| Boron | 20 | 1.3 | 2.9 | | |
| Cadmium | 0.50 | .011 | .12 | 0.034 | <0.50 |
| Calcium | 100 | 5.7 | 20 | -14 | <100 |
| Cerium | 1.0 | .0041 | .16 | | |
| Chromium | 1.0 | .11 | .15 | -0.029 | <1.0 |
| Cobalt | 1.0 | .012 | .14 | | |
| Copper | 1.0 | .91 | .74 | | |
| Iron | 100 | 48 | 16 | -20 | <100 |
| Lanthanum | 1.0 | .0038 | .41 | | |
| Lithium | 2.0 | .1 | .61 | | |
| Lead | 1.0 | .0081 | .13 | -0.13 | <1.0 |
| Magnesium | 100 | 1.6 | 11 | -10 | <100 |
| Manganese | 2.0 | .48 | .53 | 0.011 | <2.0 |
| Molybdenum | 1.0 | .048 | .89 | | |
| Nickel | 1.0 | .037 | .2 | | |
| Potassium | 100 | 3.4 | 7.6 | -9.6 | <100 |
| Selenium | 5.0 | .38 | 3.1 | 0.048 | <5.0 |
| Silver | 1.0 | .0047 | .13 | -0.081 | <1.0 |
| Silicon | 500 | 6.6 | 130 | | |
| Sodium | 100 | 24 | 9.9 | -8.7 | <100 |
| Strontium | 2.0 | .12 | .27 | 0.025 | <2.0 |
| Thallium | 1.0 | .021 | .86 | | |
| Tin | 2.0 | .034 | .19 | | |
| Titanium | 1.0 | .15 | .77 | | |
| Uranium | 1.0 | .0048 | .17 | | |
| Vanadium | 1.0 | .027 | .1 | | |
| Zinc | 5.0 | 1.5 | 1.1 | -0.32 | <5.0 |

Associated samples MP13152: LA49126-1, LA49126-2, LA49126-1F, LA49126-2F

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: LA49126
Account: HETILAL - Hydro-Environmental Technology, Inc.
Project: 8060.00 Indigo-Desoto Parish, LA

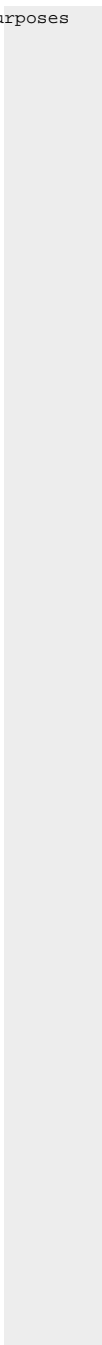
QC Batch ID: MP13152
Matrix Type: AQUEOUS

Methods: SW846 6020A
Units: ug/l

Prep Date: 10/26/18

| Metal | RL | IDL | MDL | MB raw | final |
|-------|----|-----|-----|-----------|-------|
|-------|----|-----|-----|-----------|-------|

Results < IDL are shown as zero for calculation purposes
(*) Outside of QC limits
(anr) Analyte not requested



8.1.1
8

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: LA49126
 Account: HETILAL - Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

QC Batch ID: MP13152
 Matrix Type: AQUEOUS

Methods: SW846 6020A
 Units: ug/l

Prep Date: 10/26/18

| Metal | LA49127-1 Original MS | | SpikeLot MPICPMS6 % Rec | QC Limits | |
|------------|--------------------------|--------|----------------------------|--------------|--------|
| Aluminum | 249 | 5540 | 5100 | 103.7 | 75-125 |
| Antimony | | | | | |
| Arsenic | 0.0 | 108 | 100 | 108.0 | 75-125 |
| Barium | 111 | 221 | 100 | 110.0 | 75-125 |
| Beryllium | | | | | |
| Boron | | | | | |
| Cadmium | 0.36 | 108 | 100 | 107.6 | 75-125 |
| Calcium | 9960 | 15700 | 5000 | 114.8 | 75-125 |
| Cerium | | | | | |
| Chromium | 1.1 | 111 | 100 | 109.9 | 75-125 |
| Cobalt | | | | | |
| Copper | | | | | |
| Iron | 0.0 | 5460 | 5000 | 109.2 | 75-125 |
| Lanthanum | | | | | |
| Lithium | | | | | |
| Lead | 0.0 | 103 | 100 | 103.0 | 75-125 |
| Magnesium | 3720 | 9290 | 5000 | 111.4 | 75-125 |
| Manganese | 22.5 | 132 | 100 | 109.5 | 75-125 |
| Molybdenum | | | | | |
| Nickel | | | | | |
| Potassium | 2290 | 7580 | 5000 | 105.8 | 75-125 |
| Selenium | 0.0 | 537 | 500 | 107.4 | 75-125 |
| Silver | 0.0 | 95.5 | 100 | 95.5 | 75-125 |
| Silicon | | | | | |
| Sodium | 196000 | 211000 | 5000 | 300.0(a) | 75-125 |
| Strontium | 680 | 820 | 100 | 140.0(a) | 75-125 |
| Thallium | | | | | |
| Tin | | | | | |
| Titanium | | | | | |
| Uranium | | | | | |
| Vanadium | | | | | |
| Zinc | 0.0 | 121 | 100 | 121.0 | 75-125 |

Associated samples MP13152: LA49126-1, LA49126-2, LA49126-1F, LA49126-2F

8.12
8

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: LA49126
Account: HETILAL - Hydro-Environmental Technology, Inc.
Project: 8060.00 Indigo-Desoto Parish, LA

QC Batch ID: MP13152
Matrix Type: AQUEOUS

Methods: SW846 6020A
Units: ug/l

Prep Date: 10/26/18

| Metal | LA49127-1 Original MS | SpikeLot MPICPMS6 % Rec | QC Limits |
|-------|--------------------------|----------------------------|--------------|
|-------|--------------------------|----------------------------|--------------|

Results < IDL are shown as zero for calculation purposes

- (*) Outside of QC limits
- (N) Matrix Spike Rec. outside of QC limits
- (anr) Analyte not requested
- (a) Spike amount low relative to the sample amount. Refer to lab control or spike blank for recovery information.

8.1.2
8

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: LA49126
 Account: HETILAL - Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

QC Batch ID: MP13152
 Matrix Type: AQUEOUS

Methods: SW846 6020A
 Units: ug/l

Prep Date: 10/26/18

| Metal | LA49127-1 Original MSD | | SpikeLot MPICPMS6 % Rec | | MSD RPD | QC Limit |
|------------|---------------------------|--------|----------------------------|-----------|------------|-------------|
| Aluminum | 249 | 5230 | 5100 | 97.7 | 5.8 | 20 |
| Antimony | | | | | | |
| Arsenic | 0.0 | 105 | 100 | 105.0 | 2.8 | 20 |
| Barium | 111 | 210 | 100 | 99.0 | 5.1 | 20 |
| Beryllium | | | | | | |
| Boron | | | | | | |
| Cadmium | 0.36 | 99.6 | 100 | 99.2 | 8.1 | 20 |
| Calcium | 9960 | 14900 | 5000 | 98.8 | 5.2 | 20 |
| Cerium | | | | | | |
| Chromium | 1.1 | 104 | 100 | 102.9 | 6.5 | 20 |
| Cobalt | | | | | | |
| Copper | | | | | | |
| Iron | 0.0 | 5170 | 5000 | 103.4 | 5.5 | 20 |
| Lanthanum | | | | | | |
| Lithium | | | | | | |
| Lead | 0.0 | 96.6 | 100 | 96.6 | 6.4 | 20 |
| Magnesium | 3720 | 8860 | 5000 | 102.8 | 4.7 | 20 |
| Manganese | 22.5 | 126 | 100 | 103.5 | 4.7 | 20 |
| Molybdenum | | | | | | |
| Nickel | | | | | | |
| Potassium | 2290 | 7240 | 5000 | 99.0 | 4.6 | 20 |
| Selenium | 0.0 | 496 | 500 | 99.2 | 7.9 | 20 |
| Silver | 0.0 | 91.0 | 100 | 91.0 | 4.8 | 20 |
| Silicon | | | | | | |
| Sodium | 196000 | 205000 | 5000 | 180.0(a) | 2.9 | 20 |
| Strontium | 680 | 788 | 100 | 108.0 | 4.0 | 20 |
| Thallium | | | | | | |
| Tin | | | | | | |
| Titanium | | | | | | |
| Uranium | | | | | | |
| Vanadium | | | | | | |
| Zinc | 0.0 | 128 | 100 | 128.0N(b) | 5.6 | 20 |

Associated samples MP13152: LA49126-1, LA49126-2, LA49126-1F, LA49126-2F

8.1.2
8

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: LA49126
 Account: HETILAL - Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

QC Batch ID: MP13152
 Matrix Type: AQUEOUS

Methods: SW846 6020A
 Units: ug/l

Prep Date: 10/26/18

| Metal | LA49127-1 Original MSD | Spike lot MPICPMS6 % Rec | MSD RPD | QC Limit |
|-------|---------------------------|-----------------------------|------------|-------------|
|-------|---------------------------|-----------------------------|------------|-------------|

Results < IDL are shown as zero for calculation purposes

- (*) Outside of QC limits
- (N) Matrix Spike Rec. outside of QC limits
- (anr) Analyte not requested
- (a) Spike amount low relative to the sample amount. Refer to lab control or spike blank for recovery information.
- (b) Spike recovery indicates possible matrix interference or sample non-homogeneity.

8.1.2
8

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: LA49126
 Account: HETILAL - Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

QC Batch ID: MP13152
 Matrix Type: AQUEOUS

Methods: SW846 6020A
 Units: ug/l

Prep Date: 10/26/18

| Metal | BSP Result | Spikelot MPICPMS6 | % Rec | QC Limits |
|------------|------------|-------------------|-------|-----------|
| Aluminum | 5270 | 5100 | 103.3 | 80-120 |
| Antimony | | | | |
| Arsenic | 101 | 100 | 101.0 | 80-120 |
| Barium | 107 | 100 | 107.0 | 80-120 |
| Beryllium | | | | |
| Boron | | | | |
| Cadmium | 101 | 100 | 101.0 | 80-120 |
| Calcium | 5430 | 5000 | 108.6 | 80-120 |
| Cerium | | | | |
| Chromium | 106 | 100 | 106.0 | 80-120 |
| Cobalt | | | | |
| Copper | | | | |
| Iron | 5250 | 5000 | 105.0 | 80-120 |
| Lanthanum | | | | |
| Lithium | | | | |
| Lead | 103 | 100 | 103.0 | 80-120 |
| Magnesium | 5070 | 5000 | 101.4 | 80-120 |
| Manganese | 105 | 100 | 105.0 | 80-120 |
| Molybdenum | | | | |
| Nickel | | | | |
| Potassium | 5350 | 5000 | 107.0 | 80-120 |
| Selenium | 493 | 500 | 98.6 | 80-120 |
| Silver | 101 | 100 | 101.0 | 80-120 |
| Silicon | | | | |
| Sodium | 4980 | 5000 | 99.6 | 80-120 |
| Strontium | 105 | 100 | 105.0 | 80-120 |
| Thallium | | | | |
| Tin | | | | |
| Titanium | | | | |
| Uranium | | | | |
| Vanadium | | | | |
| Zinc | 98.9 | 100 | 98.9 | 80-120 |

Associated samples MP13152: LA49126-1, LA49126-2, LA49126-1F, LA49126-2F

8.1.3
8

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: LA49126
Account: HETILAL - Hydro-Environmental Technology, Inc.
Project: 8060.00 Indigo-Desoto Parish, LA

QC Batch ID: MP13152
Matrix Type: AQUEOUS

Methods: SW846 6020A
Units: ug/l

Prep Date: 10/26/18

| Metal | BSP Result | Spikelot MPICPMS6 % Rec | QC Limits |
|-------|---------------|----------------------------|--------------|
|-------|---------------|----------------------------|--------------|

Results < IDL are shown as zero for calculation purposes
(*) Outside of QC limits
(anr) Analyte not requested

8.1.3

8

SERIAL DILUTION RESULTS SUMMARY

Login Number: LA49126
 Account: HETILAL - Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

QC Batch ID: MP13152
 Matrix Type: AQUEOUS

Methods: SW846 6020A
 Units: ug/l

Prep Date: 10/26/18

| Metal | LA49127-1 Original SDL 10:50%DIF | | QC Limits |
|------------|-------------------------------------|--------|----------------|
| Aluminum | 249 | 0.00 | 100.0 (a) 0-10 |
| Antimony | | | |
| Arsenic | 0.00 | 0.00 | NC 0-10 |
| Barium | 111 | 94.5 | 15.0* (b) 0-10 |
| Beryllium | | | |
| Boron | | | |
| Cadmium | 0.363 | 1.93 | 431.3 (a) 0-10 |
| Calcium | 9960 | 7900 | 20.7* (b) 0-10 |
| Cerium | | | |
| Chromium | 0.00 | 0.00 | NC 0-10 |
| Cobalt | | | |
| Copper | | | |
| Iron | 0.00 | 0.00 | NC 0-10 |
| Lanthanum | | | |
| Lithium | | | |
| Lead | 0.00 | 0.00 | NC 0-10 |
| Magnesium | 3720 | 3000 | 19.4* (b) 0-10 |
| Manganese | 22.5 | 0.00 | 100.0 (a) 0-10 |
| Molybdenum | | | |
| Nickel | | | |
| Potassium | 2290 | 1630 | 28.6* (b) 0-10 |
| Selenium | 0.00 | 0.00 | NC 0-10 |
| Silver | 0.00 | 0.00 | NC 0-10 |
| Silicon | | | |
| Sodium | 196000 | 183000 | 7.0 0-10 |
| Strontium | 680 | 634 | 6.7 0-10 |
| Thallium | | | |
| Tin | | | |
| Titanium | | | |
| Uranium | | | |
| Vanadium | | | |
| Zinc | 0.00 | 0.00 | NC 0-10 |

Associated samples MP13152: LA49126-1, LA49126-2, LA49126-1F, LA49126-2F

8.1.4
8

SERIAL DILUTION RESULTS SUMMARY

Login Number: LA49126
Account: HETILAL - Hydro-Environmental Technology, Inc.
Project: 8060.00 Indigo-Desoto Parish, LA

QC Batch ID: MP13152
Matrix Type: AQUEOUS

Methods: SW846 6020A
Units: ug/l

Prep Date: 10/26/18

| | | |
|-------|-------------------------------------|--------------|
| Metal | LA49127-1 Original SDL 10:50%DIF | QC Limits |
|-------|-------------------------------------|--------------|

Results < IDL are shown as zero for calculation purposes

- (*) Outside of QC limits
- (anr) Analyte not requested
- (a) Percent difference acceptable due to low initial sample concentration (< 50 times IDL).
- (b) Serial dilution indicates possible matrix interference.

POST DIGESTATE SPIKE SUMMARY

Login Number: LA49126
 Account: HETILAL - Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

QC Batch ID: MP13152
 Matrix Type: AQUEOUS

Methods: SW846 6020A
 Units: ug/l

Prep Date:

10/26/18

| Metal | Sample ml | Final ml | LA49127-1 Raw | PS Corr.** ug/l | Spike ml | Spike ug/ml | Spike ug/l | % Rec | QC Limits |
|------------|-----------|----------|---------------|-----------------|----------|-------------|------------|-------|-----------|
| Antimony | | | | | | | | | |
| Beryllium | | | | | | | | | |
| Boron | | | | | | | | | |
| Cerium | | | | | | | | | |
| Cobalt | | | | | | | | | |
| Copper | | | | | | | | | |
| Lanthanum | | | | | | | | | |
| Lithium | | | | | | | | | |
| Molybdenum | | | | | | | | | |
| Nickel | | | | | | | | | |
| Silicon | | | | | | | | | |
| Thallium | | | | | | | | | |
| Tin | | | | | | | | | |
| Titanium | | | | | | | | | |
| Uranium | | | | | | | | | |
| Vanadium | | | | | | | | | |
| Zinc | 0.2 | 10 | | 104.8 | 0.1 | 10 | 100 | 104.8 | 75-125 |

Associated samples MP13152: LA49126-1, LA49126-2, LA49126-1F, LA49126-2F

Results < IDL are shown as zero for calculation purposes
 (*) Outside of QC limits
 (**) Corr. sample result = Raw * (sample volume / final volume)
 (anr) Analyte not requested

8.1.5
8

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: LA49126
Account: HETILAL - Hydro-Environmental Technology, Inc.
Project: 8060.00 Indigo-Desoto Parish, LA

QC Batch ID: MP13160
Matrix Type: AQUEOUS

Methods: SW846 7470A
Units: ug/l

Prep Date: 10/29/18

| Metal | RL | IDL | MDL | MB raw | final |
|---------|------|-----|------|-----------|-------|
| Mercury | 0.20 | .06 | .081 | -0.0034 | <0.20 |

Associated samples MP13160: LA49126-1, LA49126-2, LA49126-1F, LA49126-2F

Results < IDL are shown as zero for calculation purposes
(*) Outside of QC limits
(anr) Analyte not requested

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: LA49126
 Account: HETILAL - Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

QC Batch ID: MP13160
 Matrix Type: AQUEOUS

Methods: SW846 7470A
 Units: ug/l

Prep Date: 10/29/18

| Metal | TD29406-3F Original MS | SpikeLot HGSPIKE1 % Rec | QC Limits |
|-------|---------------------------|----------------------------|--------------|
|-------|---------------------------|----------------------------|--------------|

Mercury 0.0 4.1 5 82.0 75-125

Associated samples MP13160: LA49126-1, LA49126-2, LA49126-1F, LA49126-2F

Results < IDL are shown as zero for calculation purposes
 (*) Outside of QC limits
 (N) Matrix Spike Rec. outside of QC limits
 (anr) Analyte not requested

8.2.2
8

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: LA49126
 Account: HETILAL - Hydro-Environmental Technology, Inc.
 Project: 8060.00 Indigo-Desoto Parish, LA

QC Batch ID: MP13160
 Matrix Type: AQUEOUS

Methods: SW846 7470A
 Units: ug/l

Prep Date: 10/29/18

| Metal | TD29406-3F Original MSD | Spikelot HGSPIKE1 | % Rec | MSD RPD | QC Limit |
|---------|----------------------------|----------------------|-------|------------|-------------|
| Mercury | 0.0 | 4.1 | 5 | 82.0 | 0.0 20 |

Associated samples MP13160: LA49126-1, LA49126-2, LA49126-1F, LA49126-2F

Results < IDL are shown as zero for calculation purposes
 (*) Outside of QC limits
 (N) Matrix Spike Rec. outside of QC limits
 (anr) Analyte not requested

8.2.2
8

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: LA49126
Account: HETILAL - Hydro-Environmental Technology, Inc.
Project: 8060.00 Indigo-Desoto Parish, LA

QC Batch ID: MP13160
Matrix Type: AQUEOUS

Methods: SW846 7470A
Units: ug/l

Prep Date: 10/29/18

| Metal | BSP Result | Spikelot HGSPIKE1 | % Rec | QC Limits |
|---------|---------------|----------------------|-------|--------------|
| Mercury | 4.3 | 5 | 86.0 | 80-120 |

Associated samples MP13160: LA49126-1, LA49126-2, LA49126-1F, LA49126-2F

Results < IDL are shown as zero for calculation purposes
(*) Outside of QC limits
(anr) Analyte not requested

8.2.3
8

SERIAL DILUTION RESULTS SUMMARY

Login Number: LA49126
Account: HETILAL - Hydro-Environmental Technology, Inc.
Project: 8060.00 Indigo-Desoto Parish, LA

QC Batch ID: MP13160
Matrix Type: AQUEOUS

Methods: SW846 7470A
Units: ug/l

Prep Date: 10/29/18

| Metal | TD29406-3F Original | SDL 1:5 | %DIF | QC Limits |
|-------|------------------------|---------|------|--------------|
|-------|------------------------|---------|------|--------------|

Mercury 0.00 0.00 NC 0-

Associated samples MP13160: LA49126-1, LA49126-2, LA49126-1F, LA49126-2F

Results < IDL are shown as zero for calculation purposes
(*) Outside of QC limits
(anr) Analyte not requested

Misc. Forms

Custody Documents and Other Forms

(SGS Houston, TX)

Includes the following where applicable:

- Chain of Custody

SGS Sample Receipt Summary

Job Number: LA49126 **Client:** SGS **Project:** 8060.00 RL INDIGO
Date / Time Received: _____ **Delivery Method:** _____ **Airbill #'s:** _____
No. Coolers: 1 **Therm ID:** IR-3; **Temp Adjustment Factor:** 0;
Cooler Temps (Initial/Adjusted): #1: (2.8/2.8);

| | | | |
|-------------------------------------|--|-------------------------------------|--|
| Cooler Security | <u>Y or N</u> | | <u>Y or N</u> |
| 1. Custody Seals Present: | <input checked="" type="checkbox"/> <input type="checkbox"/> | 3. COC Present: | <input checked="" type="checkbox"/> <input type="checkbox"/> |
| 2. Custody Seals Intact: | <input checked="" type="checkbox"/> <input type="checkbox"/> | 4. Smpl Dates/Time OK | <input checked="" type="checkbox"/> <input type="checkbox"/> |
| | | | |
| Cooler Temperature | <u>Y or N</u> | | |
| 1. Temp criteria achieved: | <input checked="" type="checkbox"/> <input type="checkbox"/> | | |
| 2. Cooler temp verification: | _____ | | |
| 3. Cooler media: | Ice (Bag) | | |
| | | | |
| Quality Control Preservation | <u>Y or N</u> | <u>N/A</u> | <u>WTB STB</u> |
| 1. Trip Blank present / cooler: | <input type="checkbox"/> <input type="checkbox"/> | <input checked="" type="checkbox"/> | <input type="checkbox"/> <input type="checkbox"/> |
| 2. Trip Blank listed on COC: | <input type="checkbox"/> <input type="checkbox"/> | <input checked="" type="checkbox"/> | |
| 3. Samples preserved properly: | <input checked="" type="checkbox"/> <input type="checkbox"/> | | |
| 4. VOCs headspace free: | <input type="checkbox"/> <input type="checkbox"/> | <input checked="" type="checkbox"/> | |

| | |
|---|---|
| Sample Integrity - Documentation | <u>Y or N</u> |
| 1. Sample labels present on bottles: | <input checked="" type="checkbox"/> <input type="checkbox"/> |
| 2. Container labeling complete: | <input checked="" type="checkbox"/> <input type="checkbox"/> |
| 3. Sample container label / COC agree: | <input checked="" type="checkbox"/> <input type="checkbox"/> |
| | |
| Sample Integrity - Condition | <u>Y or N</u> |
| 1. Sample recvd within HT: | <input checked="" type="checkbox"/> <input type="checkbox"/> |
| 2. All containers accounted for: | <input checked="" type="checkbox"/> <input type="checkbox"/> |
| 3. Condition of sample: | Intact |
| | |
| Sample Integrity - Instructions | <u>Y or N</u> <u>N/A</u> |
| 1. Analysis requested is clear: | <input checked="" type="checkbox"/> <input type="checkbox"/> |
| 2. Bottles received for unspecified tests | <input type="checkbox"/> <input checked="" type="checkbox"/> |
| 3. Sufficient volume recvd for analysis: | <input checked="" type="checkbox"/> <input type="checkbox"/> |
| 4. Compositing instructions clear: | <input type="checkbox"/> <input type="checkbox"/> <input checked="" type="checkbox"/> |
| 5. Filtering instructions clear: | <input type="checkbox"/> <input type="checkbox"/> <input checked="" type="checkbox"/> |

Comments

9.1
9

Sample Receipt Log

Job #: LA49126 _____

Date / Time Received: 10/26/2018 11:20:00 PM _____

Initials: DS _____

Client: SGS _____

| Cooler # | Sample ID: | Vol | Bot # | Location | Pres | pH | Therm ID | Initial Temp | Therm CF | Corrected Temp |
|----------|------------|-------|-------|----------|------|--|----------|--------------|----------|----------------|
| 1 | LA49126-1 | 500ml | 1 | M1A | N/P | Note #2 - Preservative check not applicable. | IR-3 | 2.8 | 0 | 2.8 |
| 1 | LA49126-2 | 500ml | 1 | M1A | N/P | Note #2 - Preservative check not applicable. | IR-3 | 2.8 | 0 | 2.8 |

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LA49126: Chain of Custody
Page 3 of 3

General Chemistry

QC Data Summaries

(SGS Houston, TX)

Includes the following where applicable:

- Method Blank and Blank Spike Summaries
- Duplicate Summaries
- Matrix Spike Summaries

METHOD BLANK AND SPIKE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: LA49126
Account: ALLA - SGS Scott, LA
Project: HETILAL: 8060.00 Indigo-Desoto Parish, LA

| Analyte | Batch ID | RL | MB Result | Units | Spike Amount | BSP Result | BSP %Recov | QC Limits |
|----------------------------|-----------------|-------|-----------|----------|--------------|------------|------------|-----------|
| Alkalinity, Bicarbonate | GN93922 | 5.0 | 2.0 | mg/l | | | | |
| Alkalinity, Carbonate | GN93923 | 5.0 | 0.0 | mg/l | | | | |
| Alkalinity, Total as CaCO3 | GN93920 | 5.0 | 0.0 | mg/l | 100 | 104 | 104.0 | 90-100% |
| Bromide | GP50318/GN94140 | 0.50 | 0.0 | mg/l | 10 | 10.4 | 104.0 | 90-110% |
| Bromide | GP50320/GN94143 | 0.50 | 0.0 | mg/l | 10 | 10.6 | 106.0 | 90-110% |
| Chloride | GP50318/GN94140 | 0.50 | 0.0 | mg/l | 10 | 10.1 | 101.0 | 90-110% |
| Chloride | GP50320/GN94143 | 0.50 | 0.0 | mg/l | 10 | 10.3 | 103.0 | 90-110% |
| Fluoride | GP50318/GN94140 | 0.50 | 0.0 | mg/l | 10 | 10.5 | 105.0 | 90-110% |
| Fluoride | GP50320/GN94143 | 0.50 | 0.0 | mg/l | 10 | 10.8 | 108.0 | 90-110% |
| Nitrogen, Nitrate | GP50318/GN94140 | 0.50 | 0.0 | mg/l | 10 | 10.0 | 100.0 | 90-110% |
| Nitrogen, Nitrate | GP50320/GN94143 | 0.50 | 0.0 | mg/l | 10 | 10.3 | 103.0 | 90-110% |
| Nitrogen, Nitrite | GP50318/GN94140 | 0.50 | 0.0 | mg/l | 10 | 10.4 | 104.0 | 90-110% |
| Nitrogen, Nitrite | GP50320/GN94143 | 0.50 | 0.0 | mg/l | 10 | 10.7 | 107.0 | 90-110% |
| Silica, Dissolved | GN93969 | 0.070 | 0.0 | mg/l | 1.07 | 0.97 | 90.7 | 80-120% |
| Solids, Total Dissolved | GN93903 | 10 | 0.0 | mg/l | 500 | 479 | 95.8 | 88-110% |
| Specific Conductivity | GN93899 | 1.0 | <1.0 | umhos/cm | | | | |
| Sulfate | GP50318/GN94140 | 0.50 | 0.0 | mg/l | 10 | 10.3 | 103.0 | 90-110% |
| Sulfate | GP50320/GN94143 | 0.50 | 0.0 | mg/l | 10 | 10.6 | 106.0 | 90-110% |

Associated Samples:

Batch GN93899: LA49126-1, LA49126-2
 Batch GN93903: LA49126-1, LA49126-2
 Batch GN93920: LA49126-1, LA49126-2
 Batch GN93922: LA49126-1, LA49126-2
 Batch GN93923: LA49126-1, LA49126-2
 Batch GN93969: LA49126-1, LA49126-2
 Batch GP50318: LA49126-2
 Batch GP50320: LA49126-1, LA49126-2
 (*) Outside of QC limits

10.1
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DUPLICATE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: LA49126
Account: ALLA - SGS Scott, LA
Project: HETILAL: 8060.00 Indigo-Desoto Parish, LA

| Analyte | Batch ID | QC Sample | Units | Original Result | DUP Result | RPD | QC Limits |
|----------------------------|-----------------|------------|----------|-----------------|------------|-----|-----------|
| Alkalinity, Bicarbonate | GN93922 | LA49146-1 | mg/l | 621 | 621 | 0.0 | 0-10% |
| Alkalinity, Carbonate | GN93923 | LA49146-1 | mg/l | 3.5 | 3.3 | 0.0 | 0-20% |
| Alkalinity, Total as CaCO3 | GN93920 | LA49146-1E | mg/l | 625 | 625 | 0.0 | 0-10% |
| Bromide | GP50318/GN94140 | LA49114-8 | mg/l | 0.0 | 0.0 | 0.0 | 0-19% |
| Bromide | GP50320/GN94143 | LA49128-1 | mg/l | 0.45 | 0.45 | 0.0 | 0-19% |
| Chloride | GP50318/GN94140 | LA49114-8 | mg/l | 0.0 | 0.0 | 0.0 | 0-13% |
| Chloride | GP50320/GN94143 | LA49128-1 | mg/l | 46.8 | 46.7 | 0.2 | 0-13% |
| Fluoride | GP50318/GN94140 | LA49114-8 | mg/l | 0.0 | 0.0 | 0.0 | 0-12% |
| Nitrogen, Nitrate | GP50318/GN94140 | LA49114-8 | mg/l | 0.0 | 0.0 | 0.0 | 0-14% |
| Nitrogen, Nitrite | GP50318/GN94140 | LA49114-8 | mg/l | 0.0 | 0.0 | 0.0 | 0-10% |
| Silica, Dissolved | GN93969 | LA49190-5 | mg/l | 5.6 | 0.0 | 0.0 | 0-20% |
| Solids, Total Dissolved | GN93903 | LA49126-2 | mg/l | 638 | 630 | 1.3 | 0-5% |
| Specific Conductivity | GN93899 | TD29212-8 | umhos/cm | 343 | 343 | 0.0 | 0-10% |
| Sulfate | GP50318/GN94140 | LA49114-8 | mg/l | 0.0 | 0.0 | 0.0 | 0-20% |
| Sulfate | GP50320/GN94143 | LA49128-1 | mg/l | 0.80 | 0.82 | 2.5 | 0-20% |

Associated Samples:

Batch GN93899: LA49126-1, LA49126-2
 Batch GN93903: LA49126-1, LA49126-2
 Batch GN93920: LA49126-1, LA49126-2
 Batch GN93922: LA49126-1, LA49126-2
 Batch GN93923: LA49126-1, LA49126-2
 Batch GN93969: LA49126-1, LA49126-2
 Batch GP50318: LA49126-2
 Batch GP50320: LA49126-1, LA49126-2
 (*) Outside of QC limits

10.2
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MATRIX SPIKE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: LA49126
Account: ALLA - SGS Scott, LA
Project: HETILAL: 8060.00 Indigo-Desoto Parish, LA

| Analyte | Batch ID | QC Sample | Units | Original Result | Spike Amount | MS Result | %Rec | QC Limits |
|----------------------------|-----------------|------------|-------|-----------------|--------------|-----------|-------|-----------|
| Alkalinity, Total as CaCO3 | GN93920 | LA49146-1E | mg/l | 625 | 25 | 650 | 100.0 | 75-117% |
| Bromide | GP50318/GN94140 | LA49114-8 | mg/l | 0.0 | 10 | 10.4 | 104.0 | 80-120% |
| Bromide | GP50320/GN94143 | LA49128-1 | mg/l | 0.45 | 10 | 10.8 | 103.5 | 80-120% |
| Chloride | GP50318/GN94140 | LA49114-8 | mg/l | 0.0 | 10 | 10.1 | 101.0 | 80-120% |
| Chloride | GP50320/GN94143 | LA49128-1 | mg/l | 46.8 | 50 | 104 | 114.4 | 80-120% |
| Fluoride | GP50318/GN94140 | LA49114-8 | mg/l | 0.0 | 10 | 10.5 | 105.0 | 80-120% |
| Nitrogen, Nitrate | GP50318/GN94140 | LA49114-8 | mg/l | 0.0 | 10 | 10.1 | 101.0 | 80-120% |
| Nitrogen, Nitrite | GP50318/GN94140 | LA49114-8 | mg/l | 0.0 | 10 | 10.5 | 105.0 | 80-120% |
| Silica, Dissolved | GN93969 | LA49190-5 | mg/l | 5.6 | 1.07 | 0.85 | 79.0 | 75-125% |
| Sulfate | GP50318/GN94140 | LA49114-8 | mg/l | 0.0 | 10 | 10.4 | 104.0 | 80-120% |
| Sulfate | GP50320/GN94143 | LA49128-1 | mg/l | 0.80 | 10 | 11.3 | 105.0 | 80-120% |

Associated Samples:

Batch GN93920: LA49126-1, LA49126-2

Batch GN93969: LA49126-1, LA49126-2

Batch GP50318: LA49126-2

Batch GP50320: LA49126-1, LA49126-2

(*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

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