

## LAB CHRONICLE

|                 |         |                   |                                    |
|-----------------|---------|-------------------|------------------------------------|
| <b>OrderID:</b> | Q2924   | <b>OrderDate:</b> | 8/20/2025 3:43:00 PM               |
| <b>Client:</b>  | AECOM   | <b>Project:</b>   | National Grid Equity - Brooklyn NY |
| <b>Contact:</b> | Peter S | <b>Location:</b>  | J23,VOA Lab                        |

| LabID                   | ClientID                | Matrix       | Test          | Method | Sample Date     | Prep Date | Anal Date | Received        |
|-------------------------|-------------------------|--------------|---------------|--------|-----------------|-----------|-----------|-----------------|
| <b>Q2924-01</b>         | <b>MW-10C-082025</b>    | <b>Water</b> | VOC-TCLVOA-10 | 8260D  | <b>08/20/25</b> |           | 08/21/25  | <b>08/20/25</b> |
| <b>Q2924-02</b>         | <b>MW-201-082025</b>    | <b>Water</b> | VOC-TCLVOA-10 | 8260D  | <b>08/20/25</b> |           | 08/21/25  | <b>08/20/25</b> |
| <b>Q2924-02DL</b>       | <b>MW-201-082025DL</b>  | <b>Water</b> | VOC-TCLVOA-10 | 8260D  | <b>08/20/25</b> |           | 08/22/25  | <b>08/20/25</b> |
| <b>Q2924-02DL<br/>2</b> | <b>MW-201-082025DL2</b> | <b>Water</b> | VOC-TCLVOA-10 | 8260D  | <b>08/20/25</b> |           | 08/22/25  | <b>08/20/25</b> |
| <b>Q2924-05</b>         | <b>MW-2B-082025</b>     | <b>Water</b> | VOC-TCLVOA-10 | 8260D  | <b>08/20/25</b> |           | 08/21/25  | <b>08/20/25</b> |
| <b>Q2924-05DL</b>       | <b>MW-2B-082025DL</b>   | <b>Water</b> | VOC-TCLVOA-10 | 8260D  | <b>08/20/25</b> |           | 08/22/25  | <b>08/20/25</b> |
| <b>Q2924-05DL<br/>2</b> | <b>MW-2B-082025DL2</b>  | <b>Water</b> | VOC-TCLVOA-10 | 8260D  | <b>08/20/25</b> |           | 08/22/25  | <b>08/20/25</b> |
| <b>Q2924-06</b>         | <b>MW-2A-082025</b>     | <b>Water</b> | VOC-TCLVOA-10 | 8260D  | <b>08/20/25</b> |           | 08/21/25  | <b>08/20/25</b> |
| <b>Q2924-07</b>         | <b>MW-18C-082025</b>    | <b>Water</b> | VOC-TCLVOA-10 | 8260D  | <b>08/20/25</b> |           | 08/21/25  | <b>08/20/25</b> |
| <b>Q2924-08</b>         | <b>MW-16A-082025</b>    | <b>Water</b> | VOC-TCLVOA-10 | 8260D  | <b>08/20/25</b> |           | 08/21/25  | <b>08/20/25</b> |
| <b>Q2924-09</b>         | <b>MW-16C-082025</b>    | <b>Water</b> | VOC-TCLVOA-10 | 8260D  | <b>08/20/25</b> |           | 08/21/25  | <b>08/20/25</b> |

## LAB CHRONICLE

|                 |                     |              |               |       |                 |          |                 |
|-----------------|---------------------|--------------|---------------|-------|-----------------|----------|-----------------|
| <b>Q2924-10</b> | <b>MW-8A-082025</b> | <b>Water</b> |               |       | <b>08/20/25</b> |          | <b>08/20/25</b> |
|                 |                     |              | VOC-TCLVOA-10 | 8260D |                 | 08/21/25 |                 |
| <b>Q2924-11</b> | <b>FB-01-082025</b> | <b>Water</b> |               |       | <b>08/20/25</b> |          | <b>08/20/25</b> |
|                 |                     |              | VOC-TCLVOA-10 | 8260D |                 | 08/21/25 |                 |
| <b>Q2924-12</b> | <b>TB</b>           | <b>Water</b> |               |       | <b>08/15/25</b> |          | <b>08/20/25</b> |
|                 |                     |              | VOC-TCLVOA-10 | 8260D |                 | 08/21/25 |                 |

A

B

C

D

E

F

G

**Hit Summary Sheet**  
**SW-846**

**SDG No.:** Q2924  
**Client:** AECOM

| Sample ID         | Client ID              | Matrix | Parameter                     | Concentration | C   | MDL   | RDL  | Units |
|-------------------|------------------------|--------|-------------------------------|---------------|-----|-------|------|-------|
| <b>Client ID:</b> | <b>MW-10C-082025</b>   |        |                               |               |     |       |      |       |
| Q2924-01          | MW-10C-082025          | Water  | Vinyl Chloride                | 2.60          | J   | 0.26  | 5.00 | ug/L  |
| Q2924-01          | MW-10C-082025          | Water  | cis-1,2-Dichloroethene        | 6.70          |     | 0.19  | 5.00 | ug/L  |
| Q2924-01          | MW-10C-082025          | Water  | Trichloroethene               | 5.40          |     | 0.090 | 5.00 | ug/L  |
| Q2924-01          | MW-10C-082025          | Water  | Tetrachloroethene             | 29.9          |     | 0.23  | 5.00 | ug/L  |
|                   |                        |        | <b>Total Voc :</b>            | <b>44.6</b>   |     |       |      |       |
|                   |                        |        | <b>Total Concentration:</b>   | <b>44.6</b>   |     |       |      |       |
| <b>Client ID:</b> | <b>MW-201-082025</b>   |        |                               |               |     |       |      |       |
| Q2924-02          | MW-201-082025          | Water  | Methyl tert-butyl Ether       | 0.92          | J   | 0.16  | 5.00 | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | trans-1,2-Dichloroethene      | 210           | E   | 0.23  | 5.00 | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | cis-1,2-Dichloroethene        | 150           | E   | 0.19  | 5.00 | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | Benzene                       | 1700          | E   | 0.15  | 5.00 | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | Trichloroethene               | 26.9          |     | 0.090 | 5.00 | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | Toluene                       | 3400          | E   | 0.14  | 5.00 | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | Ethyl Benzene                 | 650           | E   | 0.13  | 5.00 | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | m/p-Xylenes                   | 2400          | E   | 0.24  | 10.0 | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | o-Xylene                      | 2500          | E   | 0.12  | 5.00 | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | Styrene                       | 1900          | E   | 0.15  | 5.00 | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | Isopropylbenzene              | 8.40          |     | 0.12  | 5.00 | ug/L  |
|                   |                        |        | <b>Total Voc :</b>            | <b>12900</b>  |     |       |      |       |
| Q2924-02          | MW-201-082025          | Water  | Indene                        | * 4200        | J   | 0     | 0    | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | Benzene, 1-ethenyl-3-methyl-  | * 340         | J   | 0     | 0    | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | Benzo[c]thiophene             | * 280         | J   | 0     | 0    | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | Benzene, 1,2,3-trimethyl-     | * 230         | J   | 0     | 0    | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | Benzene, 1-ethenyl-4-methyl-  | * 960         | J   | 0     | 0    | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | 2-Methylindene                | * 620         | J   | 0     | 0    | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | Benzene, (1-methyl-2-cyclopro | * 750         | J   | 0     | 0    | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | n-propylbenzene               | * 30.4        | J   | 0.13  | 5.00 | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | 1,3,5-Trimethylbenzene        | * 190         | J   | 0.15  | 5.00 | ug/L  |
| Q2924-02          | MW-201-082025          | Water  | 1,2,4-Trimethylbenzene        | * 660         | J   | 0.14  | 5.00 | ug/L  |
|                   |                        |        | <b>Total Tics :</b>           | <b>8260</b>   |     |       |      |       |
|                   |                        |        | <b>Total Concentration:</b>   | <b>21200</b>  |     |       |      |       |
| <b>Client ID:</b> | <b>MW-201-082025DL</b> |        |                               |               |     |       |      |       |
| Q2924-02DL        | MW-201-082025DI        | Water  | trans-1,2-Dichloroethene      | 200           | JDQ | 23.0  | 500  | ug/L  |
| Q2924-02DL        | MW-201-082025DI        | Water  | cis-1,2-Dichloroethene        | 130           | JD  | 19.0  | 500  | ug/L  |
| Q2924-02DL        | MW-201-082025DI        | Water  | Benzene                       | 5200          | D   | 15.0  | 500  | ug/L  |
| Q2924-02DL        | MW-201-082025DI        | Water  | Toluene                       | 19900         | ED  | 14.0  | 500  | ug/L  |

**Hit Summary Sheet**  
**SW-846**

**SDG No.:** Q2924

**Client:** AECOM

| Sample ID            | Client ID        | Matrix | Parameter                      | Concentration | C  | MDL   | RDL  | Units |
|----------------------|------------------|--------|--------------------------------|---------------|----|-------|------|-------|
| Q2924-02DL           | MW-201-082025DI  | Water  | Ethyl Benzene                  | 570           | D  | 13.0  | 500  | ug/L  |
| Q2924-02DL           | MW-201-082025DI  | Water  | m/p-Xylenes                    | 3300          | D  | 24.0  | 1000 | ug/L  |
| Q2924-02DL           | MW-201-082025DI  | Water  | o-Xylene                       | 2000          | D  | 12.0  | 500  | ug/L  |
| Q2924-02DL           | MW-201-082025DI  | Water  | Styrene                        | 5100          | D  | 15.0  | 500  | ug/L  |
| Total Voc :          |                  |        |                                | 36400         |    |       |      |       |
| Total Concentration: |                  |        |                                | 36400         |    |       |      |       |
| Client ID:           | MW-201-082025DL2 |        |                                |               |    |       |      |       |
| Q2924-02DL2          | MW-201-082025DI  | Water  | Benzene                        | 3700          | D  | 60.0  | 2000 | ug/L  |
| Q2924-02DL2          | MW-201-082025DI  | Water  | Toluene                        | 14000         | D  | 56.0  | 2000 | ug/L  |
| Q2924-02DL2          | MW-201-082025DI  | Water  | Ethyl Benzene                  | 410           | JD | 52.0  | 2000 | ug/L  |
| Q2924-02DL2          | MW-201-082025DI  | Water  | m/p-Xylenes                    | 2200          | JD | 96.0  | 4000 | ug/L  |
| Q2924-02DL2          | MW-201-082025DI  | Water  | o-Xylene                       | 1300          | JD | 48.0  | 2000 | ug/L  |
| Q2924-02DL2          | MW-201-082025DI  | Water  | Styrene                        | 3200          | D  | 60.0  | 2000 | ug/L  |
| Total Voc :          |                  |        |                                | 24800         |    |       |      |       |
| Total Concentration: |                  |        |                                | 24800         |    |       |      |       |
| Client ID:           | MW-2B-082025     |        |                                |               |    |       |      |       |
| Q2924-05             | MW-2B-082025     | Water  | Acetone                        | 2.40          | J  | 1.50  | 25.0 | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | Methyl tert-butyl Ether        | 9.60          |    | 0.16  | 5.00 | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | trans-1,2-Dichloroethene       | 30.5          |    | 0.23  | 5.00 | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | cis-1,2-Dichloroethene         | 58.4          |    | 0.19  | 5.00 | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | Methylcyclohexane              | 1.30          | J  | 0.16  | 5.00 | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | Benzene                        | 1600          | E  | 0.15  | 5.00 | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | Trichloroethene                | 1.90          | J  | 0.090 | 5.00 | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | Toluene                        | 13.6          |    | 0.14  | 5.00 | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | Ethyl Benzene                  | 610           | E  | 0.13  | 5.00 | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | m/p-Xylenes                    | 55.8          |    | 0.24  | 10.0 | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | o-Xylene                       | 50.4          |    | 0.12  | 5.00 | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | Isopropylbenzene               | 78.7          |    | 0.12  | 5.00 | ug/L  |
| Total Voc :          |                  |        |                                | 2510          |    |       |      |       |
| Q2924-05             | MW-2B-082025     | Water  | Indene                         | * 93.1        | J  | 0     | 0    | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | Indane                         | * 2600        | J  | 0     | 0    | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | Benzene, 1,2,3-trimethyl-      | * 130         | J  | 0     | 0    | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | Benzene, 2-ethenyl-1,4-dimethy | * 150         | J  | 0     | 0    | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | 2-Methylindene                 | * 74.7        | J  | 0     | 0    | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | Benzene, 1-ethenyl-4-ethyl-    | * 180         | J  | 0     | 0    | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | Benzene, (1-methyl-2-cyclopro  | * 220         | J  | 0     | 0    | ug/L  |
| Q2924-05             | MW-2B-082025     | Water  | n-propylbenzene                | * 22.1        | J  | 0.13  | 5.00 | ug/L  |

### Hit Summary Sheet SW-846

SDG No.: Q2924

Client: AECOM

| Sample ID            | Client ID       | Matrix | Parameter                | Concentration | C   | MDL   | RDL  | Units |
|----------------------|-----------------|--------|--------------------------|---------------|-----|-------|------|-------|
| Q2924-05             | MW-2B-082025    | Water  | 1,3,5-Trimethylbenzene   | * 19.7        | J   | 0.15  | 5.00 | ug/L  |
| Q2924-05             | MW-2B-082025    | Water  | 1,2,4-Trimethylbenzene   | * 57.4        | J   | 0.14  | 5.00 | ug/L  |
| Total Tics :         |                 |        |                          | 3550          |     |       |      |       |
| Total Concentration: |                 |        |                          | 6060          |     |       |      |       |
| Client ID:           | MW-2B-082025DL  |        |                          |               |     |       |      |       |
| Q2924-05DL           | MW-2B-082025DL  | Water  | trans-1,2-Dichloroethene | 30.5          | JDQ | 4.60  | 100  | ug/L  |
| Q2924-05DL           | MW-2B-082025DL  | Water  | cis-1,2-Dichloroethene   | 47.2          | JD  | 3.80  | 100  | ug/L  |
| Q2924-05DL           | MW-2B-082025DL  | Water  | Benzene                  | 5100          | ED  | 3.00  | 100  | ug/L  |
| Q2924-05DL           | MW-2B-082025DL  | Water  | Ethyl Benzene            | 680           | D   | 2.60  | 100  | ug/L  |
| Q2924-05DL           | MW-2B-082025DL  | Water  | m/p-Xylenes              | 49.5          | JD  | 4.80  | 200  | ug/L  |
| Q2924-05DL           | MW-2B-082025DL  | Water  | o-Xylene                 | 42.5          | JD  | 2.40  | 100  | ug/L  |
| Q2924-05DL           | MW-2B-082025DL  | Water  | Isopropylbenzene         | 66.0          | JD  | 2.40  | 100  | ug/L  |
| Total Voc :          |                 |        |                          | 6020          |     |       |      |       |
| Total Concentration: |                 |        |                          | 6020          |     |       |      |       |
| Client ID:           | MW-2B-082025DL2 |        |                          |               |     |       |      |       |
| Q2924-05DL2          | MW-2B-082025DL  | Water  | cis-1,2-Dichloroethene   | 59.4          | JD  | 19.0  | 500  | ug/L  |
| Q2924-05DL2          | MW-2B-082025DL  | Water  | Benzene                  | 4900          | D   | 15.0  | 500  | ug/L  |
| Q2924-05DL2          | MW-2B-082025DL  | Water  | Ethyl Benzene            | 570           | D   | 13.0  | 500  | ug/L  |
| Total Voc :          |                 |        |                          | 5530          |     |       |      |       |
| Total Concentration: |                 |        |                          | 5530          |     |       |      |       |
| Client ID:           | MW-2A-082025    |        |                          |               |     |       |      |       |
| Q2924-06             | MW-2A-082025    | Water  | Acetone                  | 190           |     | 1.50  | 25.0 | ug/L  |
| Q2924-06             | MW-2A-082025    | Water  | Chloroform               | 2.10          | J   | 0.25  | 5.00 | ug/L  |
| Q2924-06             | MW-2A-082025    | Water  | Benzene                  | 30.7          |     | 0.15  | 5.00 | ug/L  |
| Q2924-06             | MW-2A-082025    | Water  | Trichloroethene          | 0.48          | J   | 0.090 | 5.00 | ug/L  |
| Q2924-06             | MW-2A-082025    | Water  | Tetrachloroethene        | 5.40          |     | 0.23  | 5.00 | ug/L  |
| Q2924-06             | MW-2A-082025    | Water  | Ethyl Benzene            | 0.46          | J   | 0.13  | 5.00 | ug/L  |
| Q2924-06             | MW-2A-082025    | Water  | Isopropylbenzene         | 0.26          | J   | 0.12  | 5.00 | ug/L  |
| Total Voc :          |                 |        |                          | 229           |     |       |      |       |
| Q2924-06             | MW-2A-082025    | Water  | Ethanol                  | * 61.3        | J   | 0     | 0    | ug/L  |
| Q2924-06             | MW-2A-082025    | Water  | Sulfur dioxide           | * 150         | J   | 0     | 0    | ug/L  |
| Q2924-06             | MW-2A-082025    | Water  | Isopropyl Alcohol        | * 14.4        | J   | 0     | 0    | ug/L  |
| Total Tics :         |                 |        |                          | 226           |     |       |      |       |
| Total Concentration: |                 |        |                          | 455           |     |       |      |       |
| Client ID:           | MW-18C-082025   |        |                          |               |     |       |      |       |
| Q2924-07             | MW-18C-082025   | Water  | Vinyl Chloride           | 37.3          |     | 0.26  | 5.00 | ug/L  |
| Q2924-07             | MW-18C-082025   | Water  | Acetone                  | 1.70          | J   | 1.50  | 25.0 | ug/L  |
| Q2924-07             | MW-18C-082025   | Water  | Methyl tert-butyl Ether  | 5.30          |     | 0.16  | 5.00 | ug/L  |

**Hit Summary Sheet**  
**SW-846**

**SDG No.:** Q2924

**Client:** AECOM

| Sample ID         | Client ID            | Matrix | Parameter                   | Concentration | C | MDL   | RDL  | Units |
|-------------------|----------------------|--------|-----------------------------|---------------|---|-------|------|-------|
| Q2924-07          | MW-18C-082025        | Water  | trans-1,2-Dichloroethene    | 5.90          |   | 0.23  | 5.00 | ug/L  |
| Q2924-07          | MW-18C-082025        | Water  | cis-1,2-Dichloroethene      | 6.20          |   | 0.19  | 5.00 | ug/L  |
| Q2924-07          | MW-18C-082025        | Water  | Benzene                     | 0.65          | J | 0.15  | 5.00 | ug/L  |
| Q2924-07          | MW-18C-082025        | Water  | Trichloroethene             | 2.20          | J | 0.090 | 5.00 | ug/L  |
| Q2924-07          | MW-18C-082025        | Water  | Toluene                     | 4.30          | J | 0.14  | 5.00 | ug/L  |
| Q2924-07          | MW-18C-082025        | Water  | Tetrachloroethene           | 2.10          | J | 0.23  | 5.00 | ug/L  |
|                   |                      |        | <b>Total Voc :</b>          |               |   | 65.7  |      |       |
| Q2924-07          | MW-18C-082025        | Water  | Sulfur dioxide              | * 38.7        | J | 0     | 0    | ug/L  |
|                   |                      |        | <b>Total Tics :</b>         |               |   | 38.7  |      |       |
|                   |                      |        | <b>Total Concentration:</b> |               |   | 104   |      |       |
| <b>Client ID:</b> | <b>MW-16A-082025</b> |        |                             |               |   |       |      |       |
| Q2924-08          | MW-16A-082025        | Water  | Sulfur dioxide              | * 100         | J | 0     | 0    | ug/L  |
|                   |                      |        | <b>Total Tics :</b>         |               |   | 100   |      |       |
|                   |                      |        | <b>Total Concentration:</b> |               |   | 100   |      |       |
| <b>Client ID:</b> | <b>MW-16C-082025</b> |        |                             |               |   |       |      |       |
| Q2924-09          | MW-16C-082025        | Water  | Vinyl Chloride              | 53.8          |   | 0.26  | 5.00 | ug/L  |
| Q2924-09          | MW-16C-082025        | Water  | 1,1-Dichloroethene          | 0.65          | J | 0.23  | 5.00 | ug/L  |
| Q2924-09          | MW-16C-082025        | Water  | Methyl tert-butyl Ether     | 17.7          |   | 0.16  | 5.00 | ug/L  |
| Q2924-09          | MW-16C-082025        | Water  | trans-1,2-Dichloroethene    | 85.5          |   | 0.23  | 5.00 | ug/L  |
| Q2924-09          | MW-16C-082025        | Water  | cis-1,2-Dichloroethene      | 32.7          |   | 0.19  | 5.00 | ug/L  |
| Q2924-09          | MW-16C-082025        | Water  | Benzene                     | 5.50          |   | 0.15  | 5.00 | ug/L  |
| Q2924-09          | MW-16C-082025        | Water  | Trichloroethene             | 14.7          |   | 0.090 | 5.00 | ug/L  |
| Q2924-09          | MW-16C-082025        | Water  | Tetrachloroethene           | 1.30          | J | 0.23  | 5.00 | ug/L  |
| Q2924-09          | MW-16C-082025        | Water  | Ethyl Benzene               | 0.61          | J | 0.13  | 5.00 | ug/L  |
| Q2924-09          | MW-16C-082025        | Water  | Isopropylbenzene            | 0.54          | J | 0.12  | 5.00 | ug/L  |
|                   |                      |        | <b>Total Voc :</b>          |               |   | 213   |      |       |
| Q2924-09          | MW-16C-082025        | Water  | 2-Methylindene              | * 7.10        | J | 0     | 0    | ug/L  |
| Q2924-09          | MW-16C-082025        | Water  | Sulfur dioxide              | * 10.9        | J | 0     | 0    | ug/L  |
| Q2924-09          | MW-16C-082025        | Water  | Benzene, 1-ethenyl-3-ethyl- | * 46.6        | J | 0     | 0    | ug/L  |
|                   |                      |        | <b>Total Tics :</b>         |               |   | 64.6  |      |       |
|                   |                      |        | <b>Total Concentration:</b> |               |   | 278   |      |       |
| <b>Client ID:</b> | <b>MW-8A-082025</b>  |        |                             |               |   |       |      |       |
| Q2924-10          | MW-8A-082025         | Water  | Acetone                     | 2.80          | J | 1.50  | 25.0 | ug/L  |
| Q2924-10          | MW-8A-082025         | Water  | Trichloroethene             | 0.49          | J | 0.090 | 5.00 | ug/L  |
|                   |                      |        | <b>Total Voc :</b>          |               |   | 3.29  |      |       |
| Q2924-10          | MW-8A-082025         | Water  | Sulfur dioxide              | * 20.8        | J | 0     | 0    | ug/L  |
|                   |                      |        | <b>Total Tics :</b>         |               |   | 20.8  |      |       |
|                   |                      |        | <b>Total Concentration:</b> |               |   | 24.1  |      |       |
| <b>Client ID:</b> | <b>FB-01-082025</b>  |        |                             |               |   |       |      |       |

**Hit Summary Sheet**  
**SW-846**

**SDG No.:** Q2924

**Client:** AECOM

| Sample ID | Client ID    | Matrix | Parameter                   | Concentration | C | MDL  | RDL  | Units |
|-----------|--------------|--------|-----------------------------|---------------|---|------|------|-------|
| Q2924-11  | FB-01-082025 | Water  | Acetone                     | 14.1          | J | 1.50 | 25.0 | ug/L  |
|           |              |        | <b>Total Voc :</b>          | 14.1          |   |      |      |       |
|           |              |        | <b>Total Concentration:</b> | 14.1          |   |      |      |       |



# SAMPLE DATA

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-10C-082025                      |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-01                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038997.D        | 1         | 08/21/25 21:28 | VV082125      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 5.00  | U         | 0.32  | 5.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 2.60  | J         | 0.26  | 5.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 5.00  | U         | 1.40  | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 5.00  | U         | 0.47  | 5.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 5.00  | U         | 0.33  | 5.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 67-64-1        | Acetone                        | 25.0  | U         | 1.50  | 25.0       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 5.00  | U         | 0.21  | 5.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 5.00  | U         | 0.16  | 5.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 5.00  | U         | 0.27  | 5.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 5.00  | U         | 0.28  | 5.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 5.00  | U         | 1.50  | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 25.0  | U         | 0.98  | 25.0       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 6.70  |           | 0.19  | 5.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 67-66-3        | Chloroform                     | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 5.00  | U         | 0.16  | 5.00       | ug/L  |
| 71-43-2        | Benzene                        | 5.00  | U         | 0.15  | 5.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 5.40  |           | 0.090 | 5.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 25.0  | U         | 0.68  | 25.0       | ug/L  |
| 108-88-3       | Toluene                        | 5.00  | U         | 0.14  | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-10C-082025                      |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-01                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038997.D        | 1         | 08/21/25 21:28 | VV082125      |

| CAS Number                | Parameter                   | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|---------|-----------|----------|------------|---------|
| 10061-02-6                | t-1,3-Dichloropropene       | 5.00    | U         | 0.17     | 5.00       | ug/L    |
| 10061-01-5                | cis-1,3-Dichloropropene     | 5.00    | U         | 0.16     | 5.00       | ug/L    |
| 79-00-5                   | 1,1,2-Trichloroethane       | 5.00    | U         | 0.21     | 5.00       | ug/L    |
| 591-78-6                  | 2-Hexanone                  | 25.0    | U         | 0.89     | 25.0       | ug/L    |
| 124-48-1                  | Dibromochloromethane        | 5.00    | U         | 0.18     | 5.00       | ug/L    |
| 106-93-4                  | 1,2-Dibromoethane           | 5.00    | U         | 0.15     | 5.00       | ug/L    |
| 127-18-4                  | Tetrachloroethene           | 29.9    |           | 0.23     | 5.00       | ug/L    |
| 108-90-7                  | Chlorobenzene               | 5.00    | U         | 0.12     | 5.00       | ug/L    |
| 100-41-4                  | Ethyl Benzene               | 5.00    | U         | 0.13     | 5.00       | ug/L    |
| 179601-23-1               | m/p-Xylenes                 | 10.0    | U         | 0.24     | 10.0       | ug/L    |
| 95-47-6                   | o-Xylene                    | 5.00    | U         | 0.12     | 5.00       | ug/L    |
| 100-42-5                  | Styrene                     | 5.00    | U         | 0.15     | 5.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 5.00    | U         | 0.19     | 5.00       | ug/L    |
| 98-82-8                   | Isopropylbenzene            | 5.00    | U         | 0.12     | 5.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 5.00    | U         | 0.26     | 5.00       | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene         | 5.00    | U         | 0.16     | 5.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 5.00    | U         | 0.19     | 5.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 5.00    | U         | 0.16     | 5.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 5.00    | U         | 0.53     | 5.00       | ug/L    |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 5.00    | U         | 0.20     | 5.00       | ug/L    |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 5.00    | U         | 0.20     | 5.00       | ug/L    |
| <b>SURROGATES</b>         |                             |         |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 54.8    |           | 74 - 125 | 110%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 46.5    |           | 75 - 124 | 93%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 45.0    |           | 86 - 113 | 90%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 41.7    |           | 77 - 121 | 83%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |         |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 585000  | 4.6       |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 1060000 | 5.532     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 975000  | 8.776     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 355000  | 11.175    |          |            |         |

## Report of Analysis

|                    |                                    |                 |               |
|--------------------|------------------------------------|-----------------|---------------|
| Client:            | AECOM                              | Date Collected: | 08/20/25      |
| Project:           | National Grid Equity - Brooklyn NY | Date Received:  | 08/20/25      |
| Client Sample ID:  | MW-10C-082025                      | SDG No.:        | Q2924         |
| Lab Sample ID:     | Q2924-01                           | Matrix:         | Water         |
| Analytical Method: | 8260D                              | % Solid:        | 0             |
| Sample Wt/Vol:     | 5                                  | Units:          | mL            |
| Soil Aliquot Vol:  |                                    |                 | uL            |
| GC Column:         | DB-624UI                           | ID :            | 0.18          |
| Prep Method :      |                                    | Level :         | LOW           |
|                    |                                    | Final Vol:      | 5000 uL       |
|                    |                                    | Test:           | VOC-TCLVOA-10 |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038997.D        | 1         | 08/21/25 21:28 | VV082125      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

U = Not Detected  
 LOQ = Limit of Quantitation  
 MDL = Method Detection Limit  
 LOD = Limit of Detection  
 E = Value Exceeds Calibration Range  
 Q = indicates LCS control criteria did not meet requirements  
 M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value  
 B = Analyte Found in Associated Method Blank  
 N = Presumptive Evidence of a Compound  
 \* = Values outside of QC limits  
 D = Dilution  
 () = Laboratory InHouse Limit  
 A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-201-082025                      |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-02                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039004.D        | 1         | 08/21/25 23:59 | VV082125      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 5.00  | U         | 0.32  | 5.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 5.00  | U         | 0.26  | 5.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 5.00  | U         | 1.40  | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 5.00  | U         | 0.47  | 5.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 5.00  | U         | 0.33  | 5.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 67-64-1        | Acetone                        | 25.0  | U         | 1.50  | 25.0       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 5.00  | U         | 0.21  | 5.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 0.92  | J         | 0.16  | 5.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 5.00  | U         | 0.27  | 5.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 5.00  | U         | 0.28  | 5.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 210   | E         | 0.23  | 5.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 5.00  | U         | 1.50  | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 25.0  | U         | 0.98  | 25.0       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 150   | E         | 0.19  | 5.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 67-66-3        | Chloroform                     | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 5.00  | U         | 0.16  | 5.00       | ug/L  |
| 71-43-2        | Benzene                        | 1700  | E         | 0.15  | 5.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 26.9  |           | 0.090 | 5.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 25.0  | U         | 0.68  | 25.0       | ug/L  |
| 108-88-3       | Toluene                        | 3400  | E         | 0.14  | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |                 |          |
|--------------------|------------------------------------|-----------------|----------|
| Client:            | AECOM                              | Date Collected: | 08/20/25 |
| Project:           | National Grid Equity - Brooklyn NY | Date Received:  | 08/20/25 |
| Client Sample ID:  | MW-201-082025                      | SDG No.:        | Q2924    |
| Lab Sample ID:     | Q2924-02                           | Matrix:         | Water    |
| Analytical Method: | 8260D                              | % Solid:        | 0        |
| Sample Wt/Vol:     | 5                                  | Units:          | mL       |
| Soil Aliquot Vol:  |                                    |                 | uL       |
| GC Column:         | DB-624UI                           | ID :            | 0.18     |
| Prep Method :      |                                    | Level :         | LOW      |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039004.D        | 1         | 08/21/25 23:59 | VV082125      |

| CAS Number                | Parameter                   | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|--------|-----------|----------|------------|---------|
| 10061-02-6                | t-1,3-Dichloropropene       | 5.00   | U         | 0.17     | 5.00       | ug/L    |
| 10061-01-5                | cis-1,3-Dichloropropene     | 5.00   | U         | 0.16     | 5.00       | ug/L    |
| 79-00-5                   | 1,1,2-Trichloroethane       | 5.00   | U         | 0.21     | 5.00       | ug/L    |
| 591-78-6                  | 2-Hexanone                  | 25.0   | U         | 0.89     | 25.0       | ug/L    |
| 124-48-1                  | Dibromochloromethane        | 5.00   | U         | 0.18     | 5.00       | ug/L    |
| 106-93-4                  | 1,2-Dibromoethane           | 5.00   | U         | 0.15     | 5.00       | ug/L    |
| 127-18-4                  | Tetrachloroethene           | 5.00   | U         | 0.23     | 5.00       | ug/L    |
| 108-90-7                  | Chlorobenzene               | 5.00   | U         | 0.12     | 5.00       | ug/L    |
| 100-41-4                  | Ethyl Benzene               | 650    | E         | 0.13     | 5.00       | ug/L    |
| 179601-23-1               | m/p-Xylenes                 | 2400   | E         | 0.24     | 10.0       | ug/L    |
| 95-47-6                   | o-Xylene                    | 2500   | E         | 0.12     | 5.00       | ug/L    |
| 100-42-5                  | Styrene                     | 1900   | E         | 0.15     | 5.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 5.00   | U         | 0.19     | 5.00       | ug/L    |
| 98-82-8                   | Isopropylbenzene            | 8.40   |           | 0.12     | 5.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 5.00   | U         | 0.26     | 5.00       | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene         | 5.00   | U         | 0.16     | 5.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 5.00   | U         | 0.19     | 5.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 5.00   | U         | 0.16     | 5.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 5.00   | U         | 0.53     | 5.00       | ug/L    |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 5.00   | U         | 0.20     | 5.00       | ug/L    |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 5.00   | U         | 0.20     | 5.00       | ug/L    |
| <b>SURROGATES</b>         |                             |        |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 57.7   |           | 74 - 125 | 115%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 44.5   |           | 75 - 124 | 89%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 44.9   |           | 86 - 113 | 90%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 55.7   |           | 77 - 121 | 111%       | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |        |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 456000 | 4.606     |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 817000 | 5.535     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 704000 | 8.779     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 443000 | 11.178    |          |            |         |

### TENTATIVE IDENTIFIED COMPOUNDS

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-201-082025                      |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-02                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039004.D        | 1         | 08/21/25 23:59 | VV082125      |

| CAS Number  | Parameter                          | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|-------------|------------------------------------|-------|-----------|-----|------------|-------|
| 103-65-1    | n-propylbenzene                    | 30.4  | J         |     | 10.3       | ug/L  |
| 108-67-8    | 1,3,5-Trimethylbenzene             | 190   | J         |     | 10.5       | ug/L  |
| 95-63-6     | 1,2,4-Trimethylbenzene             | 660   | J         |     | 10.8       | ug/L  |
| 000622-97-9 | Benzene, 1-ethenyl-4-methyl-       | 960   | J         |     | 10.9       | ug/L  |
| 000100-80-1 | Benzene, 1-ethenyl-3-methyl-       | 340   | J         |     | 11.0       | ug/L  |
| 000526-73-8 | Benzene, 1,2,3-trimethyl-          | 230   | J         |     | 11.3       | ug/L  |
| 000095-13-6 | Indene                             | 4200  | J         |     | 11.7       | ug/L  |
| 002177-47-1 | 2-Methylindene                     | 620   | J         |     | 12.9       | ug/L  |
| 065051-83-4 | Benzene, (1-methyl-2-cyclopropen-1 | 750   | J         |     | 13.0       | ug/L  |
| 000270-82-6 | Benzo[c]thiophene                  | 280   | J         |     | 13.6       | ug/L  |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-201-082025DL                    |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-02DL                         |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039015.D        | 100       | 08/22/25 14:10 | VV082225      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |      |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 500   | UD        | 22.0 | 500        | ug/L  |
| 74-87-3        | Chloromethane                  | 500   | UD        | 32.0 | 500        | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 500   | UD        | 26.0 | 500        | ug/L  |
| 74-83-9        | Bromomethane                   | 500   | UD        | 140  | 500        | ug/L  |
| 75-00-3        | Chloroethane                   | 500   | UD        | 47.0 | 500        | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 500   | UD        | 33.0 | 500        | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 500   | UD        | 25.0 | 500        | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 500   | UD        | 23.0 | 500        | ug/L  |
| 67-64-1        | Acetone                        | 2500  | UD        | 150  | 2500       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 500   | UD        | 21.0 | 500        | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 500   | UD        | 16.0 | 500        | ug/L  |
| 79-20-9        | Methyl Acetate                 | 500   | UD        | 27.0 | 500        | ug/L  |
| 75-09-2        | Methylene Chloride             | 500   | UD        | 28.0 | 500        | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 200   | JDQ       | 23.0 | 500        | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 500   | UD        | 23.0 | 500        | ug/L  |
| 110-82-7       | Cyclohexane                    | 500   | UD        | 150  | 500        | ug/L  |
| 78-93-3        | 2-Butanone                     | 2500  | UD        | 98.0 | 2500       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 500   | UD        | 25.0 | 500        | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 130   | JD        | 19.0 | 500        | ug/L  |
| 74-97-5        | Bromochloromethane             | 500   | UD        | 22.0 | 500        | ug/L  |
| 67-66-3        | Chloroform                     | 500   | UD        | 25.0 | 500        | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 500   | UD        | 20.0 | 500        | ug/L  |
| 108-87-2       | Methylcyclohexane              | 500   | UD        | 16.0 | 500        | ug/L  |
| 71-43-2        | Benzene                        | 5200  | D         | 15.0 | 500        | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 500   | UD        | 22.0 | 500        | ug/L  |
| 79-01-6        | Trichloroethene                | 500   | UD        | 9.30 | 500        | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 500   | UD        | 20.0 | 500        | ug/L  |
| 75-27-4        | Bromodichloromethane           | 500   | UD        | 22.0 | 500        | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 2500  | UD        | 68.0 | 2500       | ug/L  |
| 108-88-3       | Toluene                        | 19900 | ED        | 14.0 | 500        | ug/L  |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-201-082025DL                    |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-02DL                         |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039015.D        | 100       | 08/22/25 14:10 | VV082225      |

| CAS Number                | Parameter                   | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|---------|-----------|----------|------------|---------|
| 10061-02-6                | t-1,3-Dichloropropene       | 500     | UD        | 17.0     | 500        | ug/L    |
| 10061-01-5                | cis-1,3-Dichloropropene     | 500     | UD        | 16.0     | 500        | ug/L    |
| 79-00-5                   | 1,1,2-Trichloroethane       | 500     | UD        | 21.0     | 500        | ug/L    |
| 591-78-6                  | 2-Hexanone                  | 2500    | UD        | 89.0     | 2500       | ug/L    |
| 124-48-1                  | Dibromochloromethane        | 500     | UD        | 18.0     | 500        | ug/L    |
| 106-93-4                  | 1,2-Dibromoethane           | 500     | UD        | 15.0     | 500        | ug/L    |
| 127-18-4                  | Tetrachloroethene           | 500     | UD        | 23.0     | 500        | ug/L    |
| 108-90-7                  | Chlorobenzene               | 500     | UD        | 12.0     | 500        | ug/L    |
| 100-41-4                  | Ethyl Benzene               | 570     | D         | 13.0     | 500        | ug/L    |
| 179601-23-1               | m/p-Xylenes                 | 3300    | D         | 24.0     | 1000       | ug/L    |
| 95-47-6                   | o-Xylene                    | 2000    | D         | 12.0     | 500        | ug/L    |
| 100-42-5                  | Styrene                     | 5100    | D         | 15.0     | 500        | ug/L    |
| 75-25-2                   | Bromoform                   | 500     | UD        | 19.0     | 500        | ug/L    |
| 98-82-8                   | Isopropylbenzene            | 500     | UD        | 12.0     | 500        | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 500     | UD        | 26.0     | 500        | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene         | 500     | UD        | 16.0     | 500        | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 500     | UD        | 19.0     | 500        | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 500     | UD        | 16.0     | 500        | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 500     | UD        | 53.0     | 500        | ug/L    |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 500     | UD        | 20.0     | 500        | ug/L    |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 500     | UD        | 20.0     | 500        | ug/L    |
| <b>SURROGATES</b>         |                             |         |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 59.8    |           | 74 - 125 | 120%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 44.9    |           | 75 - 124 | 90%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 45.7    |           | 86 - 113 | 91%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 46.9    |           | 77 - 121 | 94%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |         |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 582000  | 4.6       |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 1170000 | 5.532     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 1040000 | 8.776     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 510000  | 11.172    |          |            |         |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-201-082025DL                    |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-02DL                         |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039015.D        | 100       | 08/22/25 14:10 | VV082225      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-201-082025DL2                   |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-02DL2                        |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039018.D        | 400       | 08/22/25 15:27 | VV082225      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |      |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 2000  | UD        | 88.0 | 2000       | ug/L  |
| 74-87-3        | Chloromethane                  | 2000  | UD        | 130  | 2000       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 2000  | UD        | 100  | 2000       | ug/L  |
| 74-83-9        | Bromomethane                   | 2000  | UD        | 580  | 2000       | ug/L  |
| 75-00-3        | Chloroethane                   | 2000  | UD        | 190  | 2000       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 2000  | UD        | 130  | 2000       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 2000  | UD        | 100  | 2000       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 2000  | UD        | 92.0 | 2000       | ug/L  |
| 67-64-1        | Acetone                        | 10000 | UD        | 600  | 10000      | ug/L  |
| 75-15-0        | Carbon Disulfide               | 2000  | UD        | 84.0 | 2000       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 2000  | UDQ       | 64.0 | 2000       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 2000  | UDQ       | 110  | 2000       | ug/L  |
| 75-09-2        | Methylene Chloride             | 2000  | UD        | 110  | 2000       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 2000  | UDQ       | 92.0 | 2000       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 2000  | UD        | 92.0 | 2000       | ug/L  |
| 110-82-7       | Cyclohexane                    | 2000  | UD        | 580  | 2000       | ug/L  |
| 78-93-3        | 2-Butanone                     | 10000 | UD        | 390  | 10000      | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 2000  | UD        | 100  | 2000       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 2000  | UD        | 76.0 | 2000       | ug/L  |
| 74-97-5        | Bromochloromethane             | 2000  | UDQ       | 88.0 | 2000       | ug/L  |
| 67-66-3        | Chloroform                     | 2000  | UD        | 100  | 2000       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 2000  | UD        | 80.0 | 2000       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 2000  | UD        | 64.0 | 2000       | ug/L  |
| 71-43-2        | Benzene                        | 3700  | D         | 60.0 | 2000       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 2000  | UD        | 88.0 | 2000       | ug/L  |
| 79-01-6        | Trichloroethene                | 2000  | UD        | 37.2 | 2000       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 2000  | UD        | 80.0 | 2000       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 2000  | UD        | 88.0 | 2000       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 10000 | UD        | 270  | 10000      | ug/L  |
| 108-88-3       | Toluene                        | 14000 | D         | 56.0 | 2000       | ug/L  |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-201-082025DL2                   |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-02DL2                        |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039018.D        | 400       | 08/22/25 15:27 | VV082225      |

| CAS Number                | Parameter                   | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|---------|-----------|----------|------------|---------|
| 10061-02-6                | t-1,3-Dichloropropene       | 2000    | UD        | 68.0     | 2000       | ug/L    |
| 10061-01-5                | cis-1,3-Dichloropropene     | 2000    | UD        | 64.0     | 2000       | ug/L    |
| 79-00-5                   | 1,1,2-Trichloroethane       | 2000    | UD        | 84.0     | 2000       | ug/L    |
| 591-78-6                  | 2-Hexanone                  | 10000   | UD        | 360      | 10000      | ug/L    |
| 124-48-1                  | Dibromochloromethane        | 2000    | UD        | 72.0     | 2000       | ug/L    |
| 106-93-4                  | 1,2-Dibromoethane           | 2000    | UD        | 60.0     | 2000       | ug/L    |
| 127-18-4                  | Tetrachloroethene           | 2000    | UD        | 92.0     | 2000       | ug/L    |
| 108-90-7                  | Chlorobenzene               | 2000    | UD        | 48.0     | 2000       | ug/L    |
| 100-41-4                  | Ethyl Benzene               | 410     | JD        | 52.0     | 2000       | ug/L    |
| 179601-23-1               | m/p-Xylenes                 | 2200    | JD        | 96.0     | 4000       | ug/L    |
| 95-47-6                   | o-Xylene                    | 1300    | JD        | 48.0     | 2000       | ug/L    |
| 100-42-5                  | Styrene                     | 3200    | D         | 60.0     | 2000       | ug/L    |
| 75-25-2                   | Bromoform                   | 2000    | UD        | 76.0     | 2000       | ug/L    |
| 98-82-8                   | Isopropylbenzene            | 2000    | UD        | 48.0     | 2000       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 2000    | UD        | 100      | 2000       | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene         | 2000    | UD        | 64.0     | 2000       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 2000    | UD        | 76.0     | 2000       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 2000    | UD        | 64.0     | 2000       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 2000    | UD        | 210      | 2000       | ug/L    |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 2000    | UD        | 80.0     | 2000       | ug/L    |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 2000    | UD        | 80.0     | 2000       | ug/L    |
| <b>SURROGATES</b>         |                             |         |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 51.2    |           | 74 - 125 | 102%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 44.3    |           | 75 - 124 | 89%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 44.2    |           | 86 - 113 | 88%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 46.6    |           | 77 - 121 | 93%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |         |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 766000  | 4.6       |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 1440000 | 5.532     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 1310000 | 8.773     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 617000  | 11.172    |          |            |         |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-201-082025DL2                   |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-02DL2                        |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039018.D        | 400       | 08/22/25 15:27 | VV082225      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-2B-082025                       |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-05                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038998.D        | 1         | 08/21/25 21:49 | VV082125      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 5.00  | U         | 0.32  | 5.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 5.00  | U         | 0.26  | 5.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 5.00  | U         | 1.40  | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 5.00  | U         | 0.47  | 5.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 5.00  | U         | 0.33  | 5.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 67-64-1        | Acetone                        | 2.40  | J         | 1.50  | 25.0       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 5.00  | U         | 0.21  | 5.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 9.60  |           | 0.16  | 5.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 5.00  | U         | 0.27  | 5.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 5.00  | U         | 0.28  | 5.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 30.5  |           | 0.23  | 5.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 5.00  | U         | 1.50  | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 25.0  | U         | 0.98  | 25.0       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 58.4  |           | 0.19  | 5.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 67-66-3        | Chloroform                     | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 1.30  | J         | 0.16  | 5.00       | ug/L  |
| 71-43-2        | Benzene                        | 1600  | E         | 0.15  | 5.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 1.90  | J         | 0.090 | 5.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 25.0  | U         | 0.68  | 25.0       | ug/L  |
| 108-88-3       | Toluene                        | 13.6  |           | 0.14  | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |                 |          |
|--------------------|------------------------------------|-----------------|----------|
| Client:            | AECOM                              | Date Collected: | 08/20/25 |
| Project:           | National Grid Equity - Brooklyn NY | Date Received:  | 08/20/25 |
| Client Sample ID:  | MW-2B-082025                       | SDG No.:        | Q2924    |
| Lab Sample ID:     | Q2924-05                           | Matrix:         | Water    |
| Analytical Method: | 8260D                              | % Solid:        | 0        |
| Sample Wt/Vol:     | 5                                  | Units:          | mL       |
| Soil Aliquot Vol:  |                                    |                 | uL       |
| GC Column:         | DB-624UI                           | ID :            | 0.18     |
| Prep Method :      |                                    | Level :         | LOW      |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038998.D        | 1         | 08/21/25 21:49 | VV082125      |

[illegible]

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-2B-082025                       |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-05                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038998.D        | 1         | 08/21/25 21:49 | VV082125      |

| CAS Number  | Parameter                          | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|-------------|------------------------------------|-------|-----------|-----|------------|-------|
| 103-65-1    | n-propylbenzene                    | 22.1  | J         |     | 10.3       | ug/L  |
| 108-67-8    | 1,3,5-Trimethylbenzene             | 19.7  | J         |     | 10.5       | ug/L  |
| 95-63-6     | 1,2,4-Trimethylbenzene             | 57.4  | J         |     | 10.8       | ug/L  |
| 000526-73-8 | Benzene, 1,2,3-trimethyl-          | 130   | J         |     | 11.3       | ug/L  |
| 000496-11-7 | Indane                             | 2600  | J         |     | 11.5       | ug/L  |
| 000095-13-6 | Indene                             | 93.1  | J         |     | 11.7       | ug/L  |
| 002039-89-6 | Benzene, 2-ethenyl-1,4-dimethyl-   | 150   | J         |     | 12.6       | ug/L  |
| 003454-07-7 | Benzene, 1-ethenyl-4-ethyl-        | 180   | J         |     | 12.8       | ug/L  |
| 002177-47-1 | 2-Methylindene                     | 74.7  | J         |     | 12.9       | ug/L  |
| 065051-83-4 | Benzene, (1-methyl-2-cyclopropen-1 | 220   | J         |     | 13.0       | ug/L  |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

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E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-2B-082025DL                     |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-05DL                         |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039014.D        | 20        | 08/22/25 13:48 | VV082225      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |      |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 100   | UD        | 4.40 | 100        | ug/L  |
| 74-87-3        | Chloromethane                  | 100   | UD        | 6.40 | 100        | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 100   | UD        | 5.20 | 100        | ug/L  |
| 74-83-9        | Bromomethane                   | 100   | UD        | 28.8 | 100        | ug/L  |
| 75-00-3        | Chloroethane                   | 100   | UD        | 9.40 | 100        | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 100   | UD        | 6.60 | 100        | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 100   | UD        | 5.00 | 100        | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 100   | UD        | 4.60 | 100        | ug/L  |
| 67-64-1        | Acetone                        | 500   | UD        | 30.2 | 500        | ug/L  |
| 75-15-0        | Carbon Disulfide               | 100   | UD        | 4.20 | 100        | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 100   | UDQ       | 3.20 | 100        | ug/L  |
| 79-20-9        | Methyl Acetate                 | 100   | UDQ       | 5.40 | 100        | ug/L  |
| 75-09-2        | Methylene Chloride             | 100   | UD        | 5.60 | 100        | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 30.5  | JDQ       | 4.60 | 100        | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 100   | UD        | 4.60 | 100        | ug/L  |
| 110-82-7       | Cyclohexane                    | 100   | UD        | 29.0 | 100        | ug/L  |
| 78-93-3        | 2-Butanone                     | 500   | UD        | 19.6 | 500        | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 100   | UD        | 5.00 | 100        | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 47.2  | JD        | 3.80 | 100        | ug/L  |
| 74-97-5        | Bromochloromethane             | 100   | UDQ       | 4.40 | 100        | ug/L  |
| 67-66-3        | Chloroform                     | 100   | UD        | 5.00 | 100        | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 100   | UD        | 4.00 | 100        | ug/L  |
| 108-87-2       | Methylcyclohexane              | 100   | UD        | 3.20 | 100        | ug/L  |
| 71-43-2        | Benzene                        | 5100  | ED        | 3.00 | 100        | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 100   | UD        | 4.40 | 100        | ug/L  |
| 79-01-6        | Trichloroethene                | 100   | UD        | 1.90 | 100        | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 100   | UD        | 4.00 | 100        | ug/L  |
| 75-27-4        | Bromodichloromethane           | 100   | UD        | 4.40 | 100        | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 500   | UD        | 13.6 | 500        | ug/L  |
| 108-88-3       | Toluene                        | 100   | UD        | 2.80 | 100        | ug/L  |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-2B-082025DL                     |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-05DL                         |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039014.D        | 20        | 08/22/25 13:48 | VV082225      |

| CAS Number                | Parameter                   | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|---------|-----------|----------|------------|---------|
| 10061-02-6                | t-1,3-Dichloropropene       | 100     | UD        | 3.40     | 100        | ug/L    |
| 10061-01-5                | cis-1,3-Dichloropropene     | 100     | UD        | 3.20     | 100        | ug/L    |
| 79-00-5                   | 1,1,2-Trichloroethane       | 100     | UD        | 4.20     | 100        | ug/L    |
| 591-78-6                  | 2-Hexanone                  | 500     | UD        | 17.8     | 500        | ug/L    |
| 124-48-1                  | Dibromochloromethane        | 100     | UD        | 3.60     | 100        | ug/L    |
| 106-93-4                  | 1,2-Dibromoethane           | 100     | UD        | 3.00     | 100        | ug/L    |
| 127-18-4                  | Tetrachloroethene           | 100     | UD        | 4.60     | 100        | ug/L    |
| 108-90-7                  | Chlorobenzene               | 100     | UD        | 2.40     | 100        | ug/L    |
| 100-41-4                  | Ethyl Benzene               | 680     | D         | 2.60     | 100        | ug/L    |
| 179601-23-1               | m/p-Xylenes                 | 49.5    | JD        | 4.80     | 200        | ug/L    |
| 95-47-6                   | o-Xylene                    | 42.5    | JD        | 2.40     | 100        | ug/L    |
| 100-42-5                  | Styrene                     | 100     | UD        | 3.00     | 100        | ug/L    |
| 75-25-2                   | Bromoform                   | 100     | UD        | 3.80     | 100        | ug/L    |
| 98-82-8                   | Isopropylbenzene            | 66.0    | JD        | 2.40     | 100        | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 100     | UD        | 5.20     | 100        | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene         | 100     | UD        | 3.20     | 100        | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 100     | UD        | 3.80     | 100        | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 100     | UD        | 3.20     | 100        | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 100     | UD        | 10.6     | 100        | ug/L    |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 100     | UD        | 4.00     | 100        | ug/L    |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 100     | UD        | 4.00     | 100        | ug/L    |
| <b>SURROGATES</b>         |                             |         |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 51.7    |           | 74 - 125 | 103%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 43.5    |           | 75 - 124 | 87%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 43.7    |           | 86 - 113 | 87%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 47.5    |           | 77 - 121 | 95%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |         |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 617000  | 4.603     |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 1160000 | 5.532     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 1030000 | 8.776     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 477000  | 11.172    |          |            |         |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-2B-082025DL                     |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-05DL                         |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039014.D        | 20        | 08/22/25 13:48 | VV082225      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-2B-082025DL2                    |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-05DL2                        |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039017.D        | 100       | 08/22/25 15:06 | VV082225      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                                |       |           |      |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 500   | UD        | 22.0 | 500        | ug/L  |
| 74-87-3        | Chloromethane                  | 500   | UD        | 32.0 | 500        | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 500   | UD        | 26.0 | 500        | ug/L  |
| 74-83-9        | Bromomethane                   | 500   | UD        | 140  | 500        | ug/L  |
| 75-00-3        | Chloroethane                   | 500   | UD        | 47.0 | 500        | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 500   | UD        | 33.0 | 500        | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 500   | UD        | 25.0 | 500        | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 500   | UD        | 23.0 | 500        | ug/L  |
| 67-64-1        | Acetone                        | 2500  | UD        | 150  | 2500       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 500   | UD        | 21.0 | 500        | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 500   | UDQ       | 16.0 | 500        | ug/L  |
| 79-20-9        | Methyl Acetate                 | 500   | UDQ       | 27.0 | 500        | ug/L  |
| 75-09-2        | Methylene Chloride             | 500   | UD        | 28.0 | 500        | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 500   | UDQ       | 23.0 | 500        | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 500   | UD        | 23.0 | 500        | ug/L  |
| 110-82-7       | Cyclohexane                    | 500   | UD        | 150  | 500        | ug/L  |
| 78-93-3        | 2-Butanone                     | 2500  | UD        | 98.0 | 2500       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 500   | UD        | 25.0 | 500        | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 59.4  | JD        | 19.0 | 500        | ug/L  |
| 74-97-5        | Bromochloromethane             | 500   | UDQ       | 22.0 | 500        | ug/L  |
| 67-66-3        | Chloroform                     | 500   | UD        | 25.0 | 500        | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 500   | UD        | 20.0 | 500        | ug/L  |
| 108-87-2       | Methylcyclohexane              | 500   | UD        | 16.0 | 500        | ug/L  |
| 71-43-2        | Benzene                        | 4900  | D         | 15.0 | 500        | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 500   | UD        | 22.0 | 500        | ug/L  |
| 79-01-6        | Trichloroethene                | 500   | UD        | 9.30 | 500        | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 500   | UD        | 20.0 | 500        | ug/L  |
| 75-27-4        | Bromodichloromethane           | 500   | UD        | 22.0 | 500        | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 2500  | UD        | 68.0 | 2500       | ug/L  |
| 108-88-3       | Toluene                        | 500   | UD        | 14.0 | 500        | ug/L  |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-2B-082025DL2                    |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-05DL2                        |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039017.D        | 100       | 08/22/25 15:06 | VV082225      |

| CAS Number                | Parameter                   | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|---------|-----------|----------|------------|---------|
| 10061-02-6                | t-1,3-Dichloropropene       | 500     | UD        | 17.0     | 500        | ug/L    |
| 10061-01-5                | cis-1,3-Dichloropropene     | 500     | UD        | 16.0     | 500        | ug/L    |
| 79-00-5                   | 1,1,2-Trichloroethane       | 500     | UD        | 21.0     | 500        | ug/L    |
| 591-78-6                  | 2-Hexanone                  | 2500    | UD        | 89.0     | 2500       | ug/L    |
| 124-48-1                  | Dibromochloromethane        | 500     | UD        | 18.0     | 500        | ug/L    |
| 106-93-4                  | 1,2-Dibromoethane           | 500     | UD        | 15.0     | 500        | ug/L    |
| 127-18-4                  | Tetrachloroethene           | 500     | UD        | 23.0     | 500        | ug/L    |
| 108-90-7                  | Chlorobenzene               | 500     | UD        | 12.0     | 500        | ug/L    |
| 100-41-4                  | Ethyl Benzene               | 570     | D         | 13.0     | 500        | ug/L    |
| 179601-23-1               | m/p-Xylenes                 | 1000    | UD        | 24.0     | 1000       | ug/L    |
| 95-47-6                   | o-Xylene                    | 500     | UD        | 12.0     | 500        | ug/L    |
| 100-42-5                  | Styrene                     | 500     | UD        | 15.0     | 500        | ug/L    |
| 75-25-2                   | Bromoform                   | 500     | UD        | 19.0     | 500        | ug/L    |
| 98-82-8                   | Isopropylbenzene            | 500     | UD        | 12.0     | 500        | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 500     | UD        | 26.0     | 500        | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene         | 500     | UD        | 16.0     | 500        | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 500     | UD        | 19.0     | 500        | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 500     | UD        | 16.0     | 500        | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 500     | UD        | 53.0     | 500        | ug/L    |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 500     | UD        | 20.0     | 500        | ug/L    |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 500     | UD        | 20.0     | 500        | ug/L    |
| <b>SURROGATES</b>         |                             |         |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 59.7    |           | 74 - 125 | 119%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 45.2    |           | 75 - 124 | 90%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 44.9    |           | 86 - 113 | 90%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 45.2    |           | 77 - 121 | 90%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |         |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 583000  | 4.596     |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 1170000 | 5.529     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 1040000 | 8.773     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 457000  | 11.172    |          |            |         |



## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-2A-082025                       |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-06                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038999.D        | 1         | 08/21/25 22:11 | VV082125      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 5.00  | U         | 0.32  | 5.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 5.00  | U         | 0.26  | 5.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 5.00  | U         | 1.40  | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 5.00  | U         | 0.47  | 5.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 5.00  | U         | 0.33  | 5.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 67-64-1        | Acetone                        | 190   |           | 1.50  | 25.0       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 5.00  | U         | 0.21  | 5.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 5.00  | U         | 0.16  | 5.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 5.00  | U         | 0.27  | 5.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 5.00  | U         | 0.28  | 5.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 5.00  | U         | 1.50  | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 25.0  | U         | 0.98  | 25.0       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 5.00  | U         | 0.19  | 5.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 67-66-3        | Chloroform                     | 2.10  | J         | 0.25  | 5.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 5.00  | U         | 0.16  | 5.00       | ug/L  |
| 71-43-2        | Benzene                        | 30.7  |           | 0.15  | 5.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 0.48  | J         | 0.090 | 5.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 25.0  | U         | 0.68  | 25.0       | ug/L  |
| 108-88-3       | Toluene                        | 5.00  | U         | 0.14  | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |                 |          |
|--------------------|------------------------------------|-----------------|----------|
| Client:            | AECOM                              | Date Collected: | 08/20/25 |
| Project:           | National Grid Equity - Brooklyn NY | Date Received:  | 08/20/25 |
| Client Sample ID:  | MW-2A-082025                       | SDG No.:        | Q2924    |
| Lab Sample ID:     | Q2924-06                           | Matrix:         | Water    |
| Analytical Method: | 8260D                              | % Solid:        | 0        |
| Sample Wt/Vol:     | 5                                  | Units:          | mL       |
| Soil Aliquot Vol:  |                                    |                 | uL       |
| GC Column:         | DB-624UI                           | ID :            | 0.18     |
| Prep Method :      |                                    | Level :         | LOW      |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038999.D        | 1         | 08/21/25 22:11 | VV082125      |

[illegible]

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-2A-082025                       |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-06                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038999.D        | 1         | 08/21/25 22:11 | VV082125      |

| CAS Number  | Parameter         | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|-------------|-------------------|-------|-----------|-----|------------|-------|
| 007446-09-5 | Sulfur dioxide    | 150   | J         |     | 1.20       | ug/L  |
| 000064-17-5 | Ethanol           | 61.3  | J         |     | 1.84       | ug/L  |
| 67-63-0     | Isopropyl Alcohol | 14.4  | J         |     | 2.22       | ug/L  |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-18C-082025                      |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-07                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039000.D        | 1         | 08/21/25 22:33 | VV082125      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 5.00  | U         | 0.32  | 5.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 37.3  |           | 0.26  | 5.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 5.00  | U         | 1.40  | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 5.00  | U         | 0.47  | 5.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 5.00  | U         | 0.33  | 5.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 67-64-1        | Acetone                        | 1.70  | J         | 1.50  | 25.0       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 5.00  | U         | 0.21  | 5.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 5.30  |           | 0.16  | 5.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 5.00  | U         | 0.27  | 5.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 5.00  | U         | 0.28  | 5.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 5.90  |           | 0.23  | 5.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 5.00  | U         | 1.50  | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 25.0  | U         | 0.98  | 25.0       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 6.20  |           | 0.19  | 5.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 67-66-3        | Chloroform                     | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 5.00  | U         | 0.16  | 5.00       | ug/L  |
| 71-43-2        | Benzene                        | 0.65  | J         | 0.15  | 5.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 2.20  | J         | 0.090 | 5.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 25.0  | U         | 0.68  | 25.0       | ug/L  |
| 108-88-3       | Toluene                        | 4.30  | J         | 0.14  | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |                 |          |
|--------------------|------------------------------------|-----------------|----------|
| Client:            | AECOM                              | Date Collected: | 08/20/25 |
| Project:           | National Grid Equity - Brooklyn NY | Date Received:  | 08/20/25 |
| Client Sample ID:  | MW-18C-082025                      | SDG No.:        | Q2924    |
| Lab Sample ID:     | Q2924-07                           | Matrix:         | Water    |
| Analytical Method: | 8260D                              | % Solid:        | 0        |
| Sample Wt/Vol:     | 5                                  | Units:          | mL       |
| Soil Aliquot Vol:  |                                    |                 | uL       |
| GC Column:         | DB-624UI                           | ID :            | 0.18     |
| Prep Method :      |                                    | Level :         | LOW      |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039000.D        | 1         | 08/21/25 22:33 | VV082125      |

[illegible]

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-18C-082025                      |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-07                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039000.D        | 1         | 08/21/25 22:33 | VV082125      |

| CAS Number  | Parameter      | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|-------------|----------------|-------|-----------|-----|------------|-------|
| 007446-09-5 | Sulfur dioxide | 38.7  | J         |     | 1.20       | ug/L  |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-16A-082025                      |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-08                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039001.D        | 1         | 08/21/25 22:54 | VV082125      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 5.00  | U         | 0.32  | 5.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 5.00  | U         | 0.26  | 5.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 5.00  | U         | 1.40  | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 5.00  | U         | 0.47  | 5.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 5.00  | U         | 0.33  | 5.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 67-64-1        | Acetone                        | 25.0  | U         | 1.50  | 25.0       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 5.00  | U         | 0.21  | 5.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 5.00  | U         | 0.16  | 5.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 5.00  | U         | 0.27  | 5.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 5.00  | U         | 0.28  | 5.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 5.00  | U         | 1.50  | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 25.0  | U         | 0.98  | 25.0       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 5.00  | U         | 0.19  | 5.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 67-66-3        | Chloroform                     | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 5.00  | U         | 0.16  | 5.00       | ug/L  |
| 71-43-2        | Benzene                        | 5.00  | U         | 0.15  | 5.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 5.00  | U         | 0.090 | 5.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 25.0  | U         | 0.68  | 25.0       | ug/L  |
| 108-88-3       | Toluene                        | 5.00  | U         | 0.14  | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |                 |               |
|--------------------|------------------------------------|-----------------|---------------|
| Client:            | AECOM                              | Date Collected: | 08/20/25      |
| Project:           | National Grid Equity - Brooklyn NY | Date Received:  | 08/20/25      |
| Client Sample ID:  | MW-16A-082025                      | SDG No.:        | Q2924         |
| Lab Sample ID:     | Q2924-08                           | Matrix:         | Water         |
| Analytical Method: | 8260D                              | % Solid:        | 0             |
| Sample Wt/Vol:     | 5 Units: mL                        | Final Vol:      | 5000 uL       |
| Soil Aliquot Vol:  | uL                                 | Test:           | VOC-TCLVOA-10 |
| GC Column:         | DB-624UI ID : 0.18                 | Level :         | LOW           |
| Prep Method :      |                                    |                 |               |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039001.D        | 1         | 08/21/25 22:54 | VV082125      |

[illegible]

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-16A-082025                      |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-08                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039001.D        | 1         | 08/21/25 22:54 | VV082125      |

| CAS Number  | Parameter      | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|-------------|----------------|-------|-----------|-----|------------|-------|
| 007446-09-5 | Sulfur dioxide | 100   | J         |     | 1.20       | ug/L  |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-16C-082025                      |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-09                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039002.D        | 1         | 08/21/25 23:16 | VV082125      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 5.00  | U         | 0.32  | 5.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 53.8  |           | 0.26  | 5.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 5.00  | U         | 1.40  | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 5.00  | U         | 0.47  | 5.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 5.00  | U         | 0.33  | 5.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 0.65  | J         | 0.23  | 5.00       | ug/L  |
| 67-64-1        | Acetone                        | 25.0  | U         | 1.50  | 25.0       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 5.00  | U         | 0.21  | 5.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 17.7  |           | 0.16  | 5.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 5.00  | U         | 0.27  | 5.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 5.00  | U         | 0.28  | 5.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 85.5  |           | 0.23  | 5.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 5.00  | U         | 1.50  | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 25.0  | U         | 0.98  | 25.0       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 32.7  |           | 0.19  | 5.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 67-66-3        | Chloroform                     | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 5.00  | U         | 0.16  | 5.00       | ug/L  |
| 71-43-2        | Benzene                        | 5.50  |           | 0.15  | 5.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 14.7  |           | 0.090 | 5.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 25.0  | U         | 0.68  | 25.0       | ug/L  |
| 108-88-3       | Toluene                        | 5.00  | U         | 0.14  | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |                 |               |
|--------------------|------------------------------------|-----------------|---------------|
| Client:            | AECOM                              | Date Collected: | 08/20/25      |
| Project:           | National Grid Equity - Brooklyn NY | Date Received:  | 08/20/25      |
| Client Sample ID:  | MW-16C-082025                      | SDG No.:        | Q2924         |
| Lab Sample ID:     | Q2924-09                           | Matrix:         | Water         |
| Analytical Method: | 8260D                              | % Solid:        | 0             |
| Sample Wt/Vol:     | 5 Units: mL                        | Final Vol:      | 5000 uL       |
| Soil Aliquot Vol:  | uL                                 | Test:           | VOC-TCLVOA-10 |
| GC Column:         | DB-624UI ID : 0.18                 | Level :         | LOW           |
| Prep Method :      |                                    |                 |               |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039002.D        | 1         | 08/21/25 23:16 | VV082125      |

[illegible]

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-16C-082025                      |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-09                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039002.D        | 1         | 08/21/25 23:16 | VV082125      |

| CAS Number  | Parameter                   | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|-------------|-----------------------------|-------|-----------|-----|------------|-------|
| 007446-09-5 | Sulfur dioxide              | 10.9  | J         |     | 1.20       | ug/L  |
| 007525-62-4 | Benzene, 1-ethenyl-3-ethyl- | 46.6  | J         |     | 12.0       | ug/L  |
| 002177-47-1 | 2-Methylindene              | 7.10  | J         |     | 12.9       | ug/L  |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-8A-082025                       |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-10                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039003.D        | 1         | 08/21/25 23:37 | VV082125      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 5.00  | U         | 0.32  | 5.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 5.00  | U         | 0.26  | 5.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 5.00  | U         | 1.40  | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 5.00  | U         | 0.47  | 5.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 5.00  | U         | 0.33  | 5.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 67-64-1        | Acetone                        | 2.80  | J         | 1.50  | 25.0       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 5.00  | U         | 0.21  | 5.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 5.00  | U         | 0.16  | 5.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 5.00  | U         | 0.27  | 5.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 5.00  | U         | 0.28  | 5.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 5.00  | U         | 1.50  | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 25.0  | U         | 0.98  | 25.0       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 5.00  | U         | 0.19  | 5.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 67-66-3        | Chloroform                     | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 5.00  | U         | 0.16  | 5.00       | ug/L  |
| 71-43-2        | Benzene                        | 5.00  | U         | 0.15  | 5.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 0.49  | J         | 0.090 | 5.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 25.0  | U         | 0.68  | 25.0       | ug/L  |
| 108-88-3       | Toluene                        | 5.00  | U         | 0.14  | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |                 |          |
|--------------------|------------------------------------|-----------------|----------|
| Client:            | AECOM                              | Date Collected: | 08/20/25 |
| Project:           | National Grid Equity - Brooklyn NY | Date Received:  | 08/20/25 |
| Client Sample ID:  | MW-8A-082025                       | SDG No.:        | Q2924    |
| Lab Sample ID:     | Q2924-10                           | Matrix:         | Water    |
| Analytical Method: | 8260D                              | % Solid:        | 0        |
| Sample Wt/Vol:     | 5                                  | Units:          | mL       |
| Soil Aliquot Vol:  |                                    |                 | uL       |
| GC Column:         | DB-624UI                           | ID :            | 0.18     |
| Prep Method :      |                                    | Level :         | LOW      |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039003.D        | 1         | 08/21/25 23:37 | VV082125      |

[illegible]

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-8A-082025                       |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-10                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039003.D        | 1         | 08/21/25 23:37 | VV082125      |

| CAS Number  | Parameter      | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|-------------|----------------|-------|-----------|-----|------------|-------|
| 007446-09-5 | Sulfur dioxide | 20.8  | J         |     | 1.20       | ug/L  |

U = Not Detected  
 LOQ = Limit of Quantitation  
 MDL = Method Detection Limit  
 LOD = Limit of Detection  
 E = Value Exceeds Calibration Range  
 Q = indicates LCS control criteria did not meet requirements  
 M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value  
 B = Analyte Found in Associated Method Blank  
 N = Presumptive Evidence of a Compound  
 \* = Values outside of QC limits  
 D = Dilution  
 () = Laboratory InHouse Limit  
 A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | FB-01-082025                       |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-11                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038996.D        | 1         | 08/21/25 21:06 | VV082125      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 5.00  | U         | 0.32  | 5.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 5.00  | U         | 0.26  | 5.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 5.00  | U         | 1.40  | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 5.00  | U         | 0.47  | 5.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 5.00  | U         | 0.33  | 5.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 67-64-1        | Acetone                        | 14.1  | J         | 1.50  | 25.0       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 5.00  | U         | 0.21  | 5.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 5.00  | U         | 0.16  | 5.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 5.00  | U         | 0.27  | 5.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 5.00  | U         | 0.28  | 5.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 5.00  | U         | 1.50  | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 25.0  | U         | 0.98  | 25.0       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 5.00  | U         | 0.19  | 5.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 67-66-3        | Chloroform                     | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 5.00  | U         | 0.16  | 5.00       | ug/L  |
| 71-43-2        | Benzene                        | 5.00  | U         | 0.15  | 5.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 5.00  | U         | 0.090 | 5.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 25.0  | U         | 0.68  | 25.0       | ug/L  |
| 108-88-3       | Toluene                        | 5.00  | U         | 0.14  | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | FB-01-082025                       |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-11                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038996.D        | 1         | 08/21/25 21:06 | VV082125      |

| CAS Number                | Parameter                   | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|---------|-----------|----------|------------|---------|
| 10061-02-6                | t-1,3-Dichloropropene       | 5.00    | U         | 0.17     | 5.00       | ug/L    |
| 10061-01-5                | cis-1,3-Dichloropropene     | 5.00    | U         | 0.16     | 5.00       | ug/L    |
| 79-00-5                   | 1,1,2-Trichloroethane       | 5.00    | U         | 0.21     | 5.00       | ug/L    |
| 591-78-6                  | 2-Hexanone                  | 25.0    | U         | 0.89     | 25.0       | ug/L    |
| 124-48-1                  | Dibromochloromethane        | 5.00    | U         | 0.18     | 5.00       | ug/L    |
| 106-93-4                  | 1,2-Dibromoethane           | 5.00    | U         | 0.15     | 5.00       | ug/L    |
| 127-18-4                  | Tetrachloroethene           | 5.00    | U         | 0.23     | 5.00       | ug/L    |
| 108-90-7                  | Chlorobenzene               | 5.00    | U         | 0.12     | 5.00       | ug/L    |
| 100-41-4                  | Ethyl Benzene               | 5.00    | U         | 0.13     | 5.00       | ug/L    |
| 179601-23-1               | m/p-Xylenes                 | 10.0    | U         | 0.24     | 10.0       | ug/L    |
| 95-47-6                   | o-Xylene                    | 5.00    | U         | 0.12     | 5.00       | ug/L    |
| 100-42-5                  | Styrene                     | 5.00    | U         | 0.15     | 5.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 5.00    | U         | 0.19     | 5.00       | ug/L    |
| 98-82-8                   | Isopropylbenzene            | 5.00    | U         | 0.12     | 5.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 5.00    | U         | 0.26     | 5.00       | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene         | 5.00    | U         | 0.16     | 5.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 5.00    | U         | 0.19     | 5.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 5.00    | U         | 0.16     | 5.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 5.00    | U         | 0.53     | 5.00       | ug/L    |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 5.00    | U         | 0.20     | 5.00       | ug/L    |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 5.00    | U         | 0.20     | 5.00       | ug/L    |
| <b>SURROGATES</b>         |                             |         |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 53.6    |           | 74 - 125 | 107%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 46.7    |           | 75 - 124 | 93%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 44.6    |           | 86 - 113 | 89%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 41.7    |           | 77 - 121 | 83%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |         |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 585000  | 4.6       |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 1080000 | 5.532     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 978000  | 8.776     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 425000  | 11.175    |          |            |         |

## Report of Analysis

|                    |                                    |                 |               |
|--------------------|------------------------------------|-----------------|---------------|
| Client:            | AECOM                              | Date Collected: | 08/20/25      |
| Project:           | National Grid Equity - Brooklyn NY | Date Received:  | 08/20/25      |
| Client Sample ID:  | FB-01-082025                       | SDG No.:        | Q2924         |
| Lab Sample ID:     | Q2924-11                           | Matrix:         | Water         |
| Analytical Method: | 8260D                              | % Solid:        | 0             |
| Sample Wt/Vol:     | 5                                  | Units:          | mL            |
| Soil Aliquot Vol:  |                                    |                 | uL            |
| GC Column:         | DB-624UI                           | ID :            | 0.18          |
| Prep Method :      |                                    | Level :         | LOW           |
|                    |                                    | Final Vol:      | 5000 uL       |
|                    |                                    | Test:           | VOC-TCLVOA-10 |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038996.D        | 1         | 08/21/25 21:06 | VV082125      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

U = Not Detected  
 LOQ = Limit of Quantitation  
 MDL = Method Detection Limit  
 LOD = Limit of Detection  
 E = Value Exceeds Calibration Range  
 Q = indicates LCS control criteria did not meet requirements  
 M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value  
 B = Analyte Found in Associated Method Blank  
 N = Presumptive Evidence of a Compound  
 \* = Values outside of QC limits  
 D = Dilution  
 () = Laboratory InHouse Limit  
 A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/15/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | TB                                 |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-12                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038995.D        | 1         | 08/21/25 20:45 | VV082125      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 5.00  | U         | 0.32  | 5.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 5.00  | U         | 0.26  | 5.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 5.00  | U         | 1.40  | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 5.00  | U         | 0.47  | 5.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 5.00  | U         | 0.33  | 5.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 67-64-1        | Acetone                        | 25.0  | U         | 1.50  | 25.0       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 5.00  | U         | 0.21  | 5.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 5.00  | U         | 0.16  | 5.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 5.00  | U         | 0.27  | 5.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 5.00  | U         | 0.28  | 5.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 5.00  | U         | 0.23  | 5.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 5.00  | U         | 1.50  | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 25.0  | U         | 0.98  | 25.0       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 5.00  | U         | 0.19  | 5.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 67-66-3        | Chloroform                     | 5.00  | U         | 0.25  | 5.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 5.00  | U         | 0.16  | 5.00       | ug/L  |
| 71-43-2        | Benzene                        | 5.00  | U         | 0.15  | 5.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 5.00  | U         | 0.090 | 5.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 5.00  | U         | 0.20  | 5.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 5.00  | U         | 0.22  | 5.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 25.0  | U         | 0.68  | 25.0       | ug/L  |
| 108-88-3       | Toluene                        | 5.00  | U         | 0.14  | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/15/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | TB                                 |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-12                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038995.D        | 1         | 08/21/25 20:45 | VV082125      |

| CAS Number                | Parameter                   | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|---------|-----------|----------|------------|---------|
| 10061-02-6                | t-1,3-Dichloropropene       | 5.00    | U         | 0.17     | 5.00       | ug/L    |
| 10061-01-5                | cis-1,3-Dichloropropene     | 5.00    | U         | 0.16     | 5.00       | ug/L    |
| 79-00-5                   | 1,1,2-Trichloroethane       | 5.00    | U         | 0.21     | 5.00       | ug/L    |
| 591-78-6                  | 2-Hexanone                  | 25.0    | U         | 0.89     | 25.0       | ug/L    |
| 124-48-1                  | Dibromochloromethane        | 5.00    | U         | 0.18     | 5.00       | ug/L    |
| 106-93-4                  | 1,2-Dibromoethane           | 5.00    | U         | 0.15     | 5.00       | ug/L    |
| 127-18-4                  | Tetrachloroethene           | 5.00    | U         | 0.23     | 5.00       | ug/L    |
| 108-90-7                  | Chlorobenzene               | 5.00    | U         | 0.12     | 5.00       | ug/L    |
| 100-41-4                  | Ethyl Benzene               | 5.00    | U         | 0.13     | 5.00       | ug/L    |
| 179601-23-1               | m/p-Xylenes                 | 10.0    | U         | 0.24     | 10.0       | ug/L    |
| 95-47-6                   | o-Xylene                    | 5.00    | U         | 0.12     | 5.00       | ug/L    |
| 100-42-5                  | Styrene                     | 5.00    | U         | 0.15     | 5.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 5.00    | U         | 0.19     | 5.00       | ug/L    |
| 98-82-8                   | Isopropylbenzene            | 5.00    | U         | 0.12     | 5.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 5.00    | U         | 0.26     | 5.00       | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene         | 5.00    | U         | 0.16     | 5.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 5.00    | U         | 0.19     | 5.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 5.00    | U         | 0.16     | 5.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 5.00    | U         | 0.53     | 5.00       | ug/L    |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 5.00    | U         | 0.20     | 5.00       | ug/L    |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 5.00    | U         | 0.20     | 5.00       | ug/L    |
| <b>SURROGATES</b>         |                             |         |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 52.9    |           | 74 - 125 | 106%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 47.1    |           | 75 - 124 | 94%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 45.1    |           | 86 - 113 | 90%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 40.6    |           | 77 - 121 | 81%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |         |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 588000  | 4.6       |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 1060000 | 5.532     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 950000  | 8.776     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 416000  | 11.175    |          |            |         |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/15/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | TB                                 |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-12                           |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038995.D        | 1         | 08/21/25 20:45 | VV082125      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

U = Not Detected  
 LOQ = Limit of Quantitation  
 MDL = Method Detection Limit  
 LOD = Limit of Detection  
 E = Value Exceeds Calibration Range  
 Q = indicates LCS control criteria did not meet requirements  
 M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value  
 B = Analyte Found in Associated Method Blank  
 N = Presumptive Evidence of a Compound  
 \* = Values outside of QC limits  
 D = Dilution  
 () = Laboratory InHouse Limit  
 A = Aldol-Condensation Reaction Products



# QC SUMMARY

## Surrogate Summary

SDG No.: **Q2924**

Client: **AECOM**

Analytical Method: **SW8260D**

| Lab Sample ID | Client ID        | Parameter             | Spike | Result | Recovery (%) | Qual | Limits (%) |      |
|---------------|------------------|-----------------------|-------|--------|--------------|------|------------|------|
|               |                  |                       |       |        |              |      | Low        | High |
| Q2924-01      | MW-10C-082025    | 1,2-Dichloroethane-d4 | 50    | 54.8   | 110          |      | 74         | 125  |
|               |                  | Dibromofluoromethane  | 50    | 46.5   | 93           |      | 75         | 124  |
|               |                  | Toluene-d8            | 50    | 45.0   | 90           |      | 86         | 113  |
|               |                  | 4-Bromofluorobenzene  | 50    | 41.7   | 83           |      | 77         | 121  |
| Q2924-02      | MW-201-082025    | 1,2-Dichloroethane-d4 | 50    | 57.6   | 115          |      | 74         | 125  |
|               |                  | Dibromofluoromethane  | 50    | 44.5   | 89           |      | 75         | 124  |
|               |                  | Toluene-d8            | 50    | 44.9   | 90           |      | 86         | 113  |
|               |                  | 4-Bromofluorobenzene  | 50    | 55.7   | 111          |      | 77         | 121  |
| Q2924-02DL    | MW-201-082025DL  | 1,2-Dichloroethane-d4 | 50    | 59.8   | 120          |      | 74         | 125  |
|               |                  | Dibromofluoromethane  | 50    | 44.9   | 90           |      | 75         | 124  |
|               |                  | Toluene-d8            | 50    | 45.7   | 91           |      | 86         | 113  |
|               |                  | 4-Bromofluorobenzene  | 50    | 46.9   | 94           |      | 77         | 121  |
| Q2924-02DL2   | MW-201-082025DL2 | 1,2-Dichloroethane-d4 | 50    | 51.2   | 102          |      | 74         | 125  |
|               |                  | Dibromofluoromethane  | 50    | 44.3   | 89           |      | 75         | 124  |
|               |                  | Toluene-d8            | 50    | 44.2   | 88           |      | 86         | 113  |
|               |                  | 4-Bromofluorobenzene  | 50    | 46.6   | 93           |      | 77         | 121  |
| Q2924-03MS    | MW-201-082025MS  | 1,2-Dichloroethane-d4 | 50    | 50.4   | 101          |      | 74         | 125  |
|               |                  | Dibromofluoromethane  | 50    | 44.4   | 89           |      | 75         | 124  |
|               |                  | Toluene-d8            | 50    | 43.9   | 88           |      | 86         | 113  |
|               |                  | 4-Bromofluorobenzene  | 50    | 45.0   | 90           |      | 77         | 121  |
| Q2924-04MSD   | MW-201-082025MSD | 1,2-Dichloroethane-d4 | 50    | 52.6   | 105          |      | 74         | 125  |
|               |                  | Dibromofluoromethane  | 50    | 47.2   | 94           |      | 75         | 124  |
|               |                  | Toluene-d8            | 50    | 44.3   | 89           |      | 86         | 113  |
|               |                  | 4-Bromofluorobenzene  | 50    | 60.7   | 121          |      | 77         | 121  |
| Q2924-05      | MW-2B-082025     | 1,2-Dichloroethane-d4 | 50    | 57.5   | 115          |      | 74         | 125  |
|               |                  | Dibromofluoromethane  | 50    | 44.9   | 90           |      | 75         | 124  |
|               |                  | Toluene-d8            | 50    | 43.4   | 87           |      | 86         | 113  |
|               |                  | 4-Bromofluorobenzene  | 50    | 54.0   | 108          |      | 77         | 121  |
| Q2924-05DL    | MW-2B-082025DL   | 1,2-Dichloroethane-d4 | 50    | 51.7   | 103          |      | 74         | 125  |
|               |                  | Dibromofluoromethane  | 50    | 43.5   | 87           |      | 75         | 124  |
|               |                  | Toluene-d8            | 50    | 43.7   | 87           |      | 86         | 113  |
|               |                  | 4-Bromofluorobenzene  | 50    | 47.5   | 95           |      | 77         | 121  |
| Q2924-05DL2   | MW-2B-082025DL2  | 1,2-Dichloroethane-d4 | 50    | 59.7   | 119          |      | 74         | 125  |
|               |                  | Dibromofluoromethane  | 50    | 45.2   | 90           |      | 75         | 124  |
|               |                  | Toluene-d8            | 50    | 44.9   | 90           |      | 86         | 113  |
|               |                  | 4-Bromofluorobenzene  | 50    | 45.2   | 90           |      | 77         | 121  |
| Q2924-06      | MW-2A-082025     | 1,2-Dichloroethane-d4 | 50    | 49.3   | 99           |      | 74         | 125  |
|               |                  | Dibromofluoromethane  | 50    | 44.6   | 89           |      | 75         | 124  |
|               |                  | Toluene-d8            | 50    | 44.0   | 88           |      | 86         | 113  |
|               |                  | 4-Bromofluorobenzene  | 50    | 41.9   | 84           |      | 77         | 121  |
| Q2924-07      | MW-18C-082025    | 1,2-Dichloroethane-d4 | 50    | 48.7   | 97           |      | 74         | 125  |
|               |                  | Dibromofluoromethane  | 50    | 44.7   | 89           |      | 75         | 124  |
|               |                  | Toluene-d8            | 50    | 44.0   | 88           |      | 86         | 113  |
|               |                  | 4-Bromofluorobenzene  | 50    | 44.2   | 88           |      | 77         | 121  |
| Q2924-08      | MW-16A-082025    | 1,2-Dichloroethane-d4 | 50    | 51.1   | 102          |      | 74         | 125  |
|               |                  | Dibromofluoromethane  | 50    | 45.8   | 91           |      | 75         | 124  |
|               |                  | Toluene-d8            | 50    | 44.4   | 89           |      | 86         | 113  |
|               |                  | 4-Bromofluorobenzene  | 50    | 44.8   | 89           |      | 77         | 121  |
| Q2924-09      | MW-16C-082025    | 1,2-Dichloroethane-d4 | 50    | 50.9   | 102          |      | 74         | 125  |
|               |                  | Dibromofluoromethane  | 50    | 44.7   | 89           |      | 75         | 124  |

## Surrogate Summary

SDG No.: Q2924

Client: AECOM

Analytical Method: SW8260D

| Lab Sample ID | Client ID     | Parameter             | Spike | Result | Recovery (%) | Qual | Limits (%) |      |
|---------------|---------------|-----------------------|-------|--------|--------------|------|------------|------|
|               |               |                       |       |        |              |      | Low        | High |
| Q2924-09      | MW-16C-082025 | Toluene-d8            | 50    | 44.2   | 88           |      | 86         | 113  |
|               |               | 4-Bromofluorobenzene  | 50    | 45.7   | 91           |      | 77         | 121  |
| Q2924-10      | MW-8A-082025  | 1,2-Dichloroethane-d4 | 50    | 51.0   | 102          |      | 74         | 125  |
|               |               | Dibromofluoromethane  | 50    | 45.0   | 90           |      | 75         | 124  |
|               |               | Toluene-d8            | 50    | 55.1   | 110          |      | 86         | 113  |
|               |               | 4-Bromofluorobenzene  | 50    | 58.8   | 118          |      | 77         | 121  |
| Q2924-11      | FB-01-082025  | 1,2-Dichloroethane-d4 | 50    | 53.6   | 107          |      | 74         | 125  |
|               |               | Dibromofluoromethane  | 50    | 46.8   | 93           |      | 75         | 124  |
|               |               | Toluene-d8            | 50    | 44.6   | 89           |      | 86         | 113  |
|               |               | 4-Bromofluorobenzene  | 50    | 41.7   | 83           |      | 77         | 121  |
| Q2924-12      | TB            | 1,2-Dichloroethane-d4 | 50    | 52.9   | 106          |      | 74         | 125  |
|               |               | Dibromofluoromethane  | 50    | 47.1   | 94           |      | 75         | 124  |
|               |               | Toluene-d8            | 50    | 45.1   | 90           |      | 86         | 113  |
|               |               | 4-Bromofluorobenzene  | 50    | 40.6   | 81           |      | 77         | 121  |

## Surrogate Summary

SDG No.: Q2924

Client: AECOM

Analytical Method: SW8260-Low

| Lab Sample ID | Client ID   | Parameter             | Spike | Result | Recovery (%) | Qual | Limits (%) |      |
|---------------|-------------|-----------------------|-------|--------|--------------|------|------------|------|
|               |             |                       |       |        |              |      | Low        | High |
| VV0821WBL01   | VV0821WBL01 | 1,2-Dichloroethane-d4 | 50    | 48.2   | 96           |      | 74         | 125  |
|               |             | Dibromofluoromethane  | 50    | 43.8   | 88           |      | 75         | 124  |
|               |             | Toluene-d8            | 50    | 43.4   | 87           |      | 86         | 113  |
|               |             | 4-Bromofluorobenzene  | 50    | 42.3   | 85           |      | 77         | 121  |
| VV0821WBS01   | VV0821WBS01 | 1,2-Dichloroethane-d4 | 50    | 55.0   | 110          |      | 74         | 125  |
|               |             | Dibromofluoromethane  | 50    | 44.8   | 90           |      | 75         | 124  |
|               |             | Toluene-d8            | 50    | 49.6   | 99           |      | 86         | 113  |
|               |             | 4-Bromofluorobenzene  | 50    | 48.8   | 98           |      | 77         | 121  |
| VV0822WBL01   | VV0822WBL01 | 1,2-Dichloroethane-d4 | 50    | 60.0   | 120          |      | 74         | 125  |
|               |             | Dibromofluoromethane  | 50    | 47.4   | 95           |      | 75         | 124  |
|               |             | Toluene-d8            | 50    | 47.7   | 95           |      | 86         | 113  |
|               |             | 4-Bromofluorobenzene  | 50    | 49.3   | 99           |      | 77         | 121  |
| VV0822WBS01   | VV0822WBS01 | 1,2-Dichloroethane-d4 | 50    | 48.5   | 97           |      | 74         | 125  |
|               |             | Dibromofluoromethane  | 50    | 46.0   | 92           |      | 75         | 124  |
|               |             | Toluene-d8            | 50    | 46.3   | 93           |      | 86         | 113  |
|               |             | 4-Bromofluorobenzene  | 50    | 48.5   | 97           |      | 77         | 121  |

### Matrix Spike/Matrix Spike Duplicate Summary

SW-846

SDG No.: Q2924

Analytical Method:

SW8260D

Client: AECOM

Datafile :

VV039005.D

|                                |            | Sample             |                 |       |      | Rec  | RPD |      | Limits |      |     |
|--------------------------------|------------|--------------------|-----------------|-------|------|------|-----|------|--------|------|-----|
| Parameter                      | Spike      | Result             | Result          | Units | Rec  | Qual | RPD | Qual | Low    | High | RPD |
| Lab Sample ID :                | Q2924-03MS | Client Sample ID : | MW-201-082025MS |       |      |      |     |      |        |      |     |
| Dichlorodifluoromethane        | 50         | 0                  | 38.4            | ug/L  | 77   |      |     |      | 73     | 120  |     |
| Chloromethane                  | 50         | 0                  | 32.4            | ug/L  | 65   |      |     |      | 58     | 133  |     |
| Vinyl chloride                 | 50         | 0                  | 39.7            | ug/L  | 79   |      |     |      | 69     | 125  |     |
| Bromomethane                   | 50         | 0                  | 24.9            | ug/L  | 50   |      |     |      | 28     | 165  |     |
| Chloroethane                   | 50         | 0                  | 52.0            | ug/L  | 104  |      |     |      | 70     | 141  |     |
| Trichlorofluoromethane         | 50         | 0                  | 45.7            | ug/L  | 91   |      |     |      | 72     | 124  |     |
| 1,1,2-Trichlorotrifluoroethane | 50         | 0                  | 44.6            | ug/L  | 89   |      |     |      | 75     | 117  |     |
| 1,1-Dichloroethene             | 50         | 0                  | 49.2            | ug/L  | 98   |      |     |      | 53     | 162  |     |
| Acetone                        | 250        | 0                  | 210             | ug/L  | 84   |      |     |      | 44     | 150  |     |
| Carbon disulfide               | 50         | 0                  | 47.5            | ug/L  | 95   |      |     |      | 44     | 135  |     |
| Methyl tert-butyl Ether        | 50         | 0.92               | 55.4            | ug/L  | 109  |      |     |      | 82     | 133  |     |
| Methyl Acetate                 | 50         | 0                  | 47.7            | ug/L  | 95   |      |     |      | 76     | 138  |     |
| Methylene Chloride             | 50         | 0                  | 43.3            | ug/L  | 87   |      |     |      | 79     | 115  |     |
| trans-1,2-Dichloroethene       | 50         | 210                | 220             | ug/L  | 20   | *    |     |      | 76     | 118  |     |
| 1,1-Dichloroethane             | 50         | 0                  | 42.5            | ug/L  | 85   |      |     |      | 78     | 122  |     |
| Cyclohexane                    | 50         | 0                  | 41.1            | ug/L  | 82   |      |     |      | 71     | 119  |     |
| 2-Butanone                     | 250        | 0                  | 230             | ug/L  | 92   |      |     |      | 67     | 137  |     |
| Carbon Tetrachloride           | 50         | 0                  | 42.2            | ug/L  | 84   |      |     |      | 66     | 133  |     |
| cis-1,2-Dichloroethene         | 50         | 150                | 170             | ug/L  | 40   | *    |     |      | 82     | 124  |     |
| Bromochloromethane             | 50         | 0                  | 45.2            | ug/L  | 90   |      |     |      | 72     | 130  |     |
| Chloroform                     | 50         | 0                  | 41.8            | ug/L  | 84   |      |     |      | 83     | 119  |     |
| 1,1,1-Trichloroethane          | 50         | 0                  | 42.5            | ug/L  | 85   |      |     |      | 83     | 117  |     |
| Methylcyclohexane              | 50         | 0                  | 45.2            | ug/L  | 90   |      |     |      | 64     | 120  |     |
| Benzene                        | 50         | 1700               | 1600            | ug/L  | -200 | *    |     |      | 81     | 128  |     |
| 1,2-Dichloroethane             | 50         | 0                  | 180             | ug/L  | 360  | *    |     |      | 76     | 120  |     |
| Trichloroethene                | 50         | 26.9               | 74.0            | ug/L  | 94   |      |     |      | 28     | 175  |     |
| 1,2-Dichloropropane            | 50         | 0                  | 40.8            | ug/L  | 82   | *    |     |      | 85     | 116  |     |
| Bromodichloromethane           | 50         | 0                  | 40.6            | ug/L  | 81   |      |     |      | 54     | 157  |     |
| 4-Methyl-2-Pentanone           | 250        | 0                  | 220             | ug/L  | 88   |      |     |      | 72     | 137  |     |
| Toluene                        | 50         | 3400               | 3300            | ug/L  | -200 | *    |     |      | 85     | 115  |     |
| t-1,3-Dichloropropene          | 50         | 0                  | 41.0            | ug/L  | 82   |      |     |      | 60     | 141  |     |
| cis-1,3-Dichloropropene        | 50         | 0                  | 42.6            | ug/L  | 85   |      |     |      | 36     | 161  |     |
| 1,1,2-Trichloroethane          | 50         | 0                  | 37.2            | ug/L  | 74   |      |     |      | 27     | 175  |     |
| 2-Hexanone                     | 250        | 0                  | 160             | ug/L  | 64   | *    |     |      | 75     | 131  |     |
| Dibromochloromethane           | 50         | 0                  | 38.7            | ug/L  | 77   |      |     |      | 59     | 164  |     |
| 1,2-Dibromoethane              | 50         | 0                  | 39.9            | ug/L  | 80   | *    |     |      | 85     | 119  |     |
| Tetrachloroethene              | 50         | 0                  | 43.7            | ug/L  | 87   |      |     |      | 48     | 153  |     |
| Chlorobenzene                  | 50         | 0                  | 44.2            | ug/L  | 88   |      |     |      | 85     | 114  |     |
| Ethyl Benzene                  | 50         | 650                | 650             | ug/L  | 0    | *    |     |      | 81     | 128  |     |
| m/p-Xylenes                    | 100        | 2400               | 2300            | ug/L  | -100 | *    |     |      | 69     | 129  |     |
| o-Xylene                       | 50         | 2500               | 2200            | ug/L  | -600 | *    |     |      | 75     | 127  |     |
| Styrene                        | 50         | 1900               | 1800            | ug/L  | -200 | *    |     |      | 84     | 128  |     |
| Bromoform                      | 50         | 0                  | 45.4            | ug/L  | 91   |      |     |      | 73     | 147  |     |
| Isopropylbenzene               | 50         | 8.40               | 43.6            | ug/L  | 70   | *    |     |      | 76     | 121  |     |
| 1,1,2,2-Tetrachloroethane      | 50         | 0                  | 28.9            | ug/L  | 58   | *    |     |      | 81     | 131  |     |

### Matrix Spike/Matrix Spike Duplicate Summary

SW-846

SDG No.: Q2924

Analytical Method: SW8260D

Client: AECOM

Datafile : VV039005.D

| Parameter                   | Spike | Sample | Result | Units | Rec |      |     | RPD  | Limits |      |     |
|-----------------------------|-------|--------|--------|-------|-----|------|-----|------|--------|------|-----|
|                             |       | Result |        |       | Rec | Qual | RPD | Qual | Low    | High | RPD |
| 1,3-Dichlorobenzene         | 50    | 0      | 45.0   | ug/L  | 90  |      |     |      | 84     | 110  |     |
| 1,4-Dichlorobenzene         | 50    | 0      | 43.8   | ug/L  | 88  |      |     |      | 81     | 111  |     |
| 1,2-Dichlorobenzene         | 50    | 0      | 46.1   | ug/L  | 92  |      |     |      | 82     | 113  |     |
| 1,2-Dibromo-3-Chloropropane | 50    | 0      | 46.0   | ug/L  | 92  |      |     |      | 79     | 137  |     |
| 1,2,4-Trichlorobenzene      | 50    | 0      | 51.9   | ug/L  | 104 |      |     |      | 73     | 120  |     |
| 1,2,3-Trichlorobenzene      | 50    | 0      | 51.6   | ug/L  | 103 |      |     |      | 75     | 119  |     |

### Matrix Spike/Matrix Spike Duplicate Summary

SW-846

SDG No.: Q2924

Analytical Method:

SW8260D

Client: AECOM

Datafile :

VV039006.D

| Parameter                      | Spike       | Sample             | Result           | Units | Rec  |     | RPD  |     | Limits |     |      |
|--------------------------------|-------------|--------------------|------------------|-------|------|-----|------|-----|--------|-----|------|
|                                |             | Result             |                  |       |      | Rec | Qual | RPD | Qual   | Low | High |
| Lab Sample ID :                | Q2924-04MSD | Client Sample ID : | MW-201-082025MSD |       |      |     |      |     |        |     |      |
| Dichlorodifluoromethane        | 50          | 0                  | 44.7             | ug/L  | 89   |     | 15   |     | 73     | 120 | 20   |
| Chloromethane                  | 50          | 0                  | 49.3             | ug/L  | 99   |     | 41   | *   | 58     | 133 | 20   |
| Vinyl chloride                 | 50          | 0                  | 57.5             | ug/L  | 115  |     | 37   | *   | 69     | 125 | 20   |
| Bromomethane                   | 50          | 0                  | 31.1             | ug/L  | 62   |     | 22   | *   | 28     | 165 | 20   |
| Chloroethane                   | 50          | 0                  | 60.7             | ug/L  | 121  |     | 15   |     | 70     | 141 | 20   |
| Trichlorofluoromethane         | 50          | 0                  | 47.5             | ug/L  | 95   |     | 4    |     | 72     | 124 | 20   |
| 1,1,2-Trichlorotrifluoroethane | 50          | 0                  | 44.0             | ug/L  | 88   |     | 1    |     | 75     | 117 | 20   |
| 1,1-Dichloroethene             | 50          | 0                  | 54.6             | ug/L  | 109  |     | 10   |     | 53     | 162 | 20   |
| Acetone                        | 250         | 0                  | 260              | ug/L  | 104  |     | 21   | *   | 44     | 150 | 20   |
| Carbon disulfide               | 50          | 0                  | 55.4             | ug/L  | 111  |     | 15   |     | 44     | 135 | 20   |
| Methyl tert-butyl Ether        | 50          | 0.92               | 67.5             | ug/L  | 133  |     | 20   |     | 82     | 133 | 20   |
| Methyl Acetate                 | 50          | 0                  | 61.9             | ug/L  | 124  |     | 26   | *   | 76     | 138 | 20   |
| Methylene Chloride             | 50          | 0                  | 53.6             | ug/L  | 107  |     | 21   | *   | 79     | 115 | 20   |
| trans-1,2-Dichloroethene       | 50          | 210                | 250              | ug/L  | 80   |     | 120  | *   | 76     | 118 | 20   |
| 1,1-Dichloroethane             | 50          | 0                  | 49.1             | ug/L  | 98   |     | 14   |     | 78     | 122 | 20   |
| Cyclohexane                    | 50          | 0                  | 38.3             | ug/L  | 77   |     | 7    |     | 71     | 119 | 20   |
| 2-Butanone                     | 250         | 0                  | 260              | ug/L  | 104  |     | 12   |     | 67     | 137 | 20   |
| Carbon Tetrachloride           | 50          | 0                  | 41.8             | ug/L  | 84   |     | 1    |     | 66     | 133 | 20   |
| cis-1,2-Dichloroethene         | 50          | 150                | 180              | ug/L  | 60   | *   | 40   | *   | 82     | 124 | 20   |
| Bromochloromethane             | 50          | 0                  | 47.4             | ug/L  | 95   |     | 5    |     | 72     | 130 | 20   |
| Chloroform                     | 50          | 0                  | 44.9             | ug/L  | 90   |     | 7    |     | 83     | 119 | 20   |
| 1,1,1-Trichloroethane          | 50          | 0                  | 43.5             | ug/L  | 87   |     | 2    |     | 83     | 117 | 20   |
| Methylcyclohexane              | 50          | 0                  | 38.9             | ug/L  | 78   |     | 15   |     | 64     | 120 | 20   |
| Benzene                        | 50          | 1700               | 1600             | ug/L  | -200 | *   | 0    |     | 81     | 128 | 20   |
| 1,2-Dichloroethane             | 50          | 0                  | 180              | ug/L  | 360  | *   | 0    |     | 76     | 120 | 20   |
| Trichloroethene                | 50          | 26.9               | 72.9             | ug/L  | 92   |     | 2    |     | 28     | 175 | 20   |
| 1,2-Dichloropropane            | 50          | 0                  | 43.1             | ug/L  | 86   |     | 5    |     | 85     | 116 | 20   |
| Bromodichloromethane           | 50          | 0                  | 43.1             | ug/L  | 86   |     | 6    |     | 54     | 157 | 20   |
| 4-Methyl-2-Pentanone           | 250         | 0                  | 230              | ug/L  | 92   |     | 4    |     | 72     | 137 | 20   |
| Toluene                        | 50          | 3400               | 3200             | ug/L  | -400 | *   | 67   | *   | 85     | 115 | 20   |
| t-1,3-Dichloropropene          | 50          | 0                  | 43.6             | ug/L  | 87   |     | 6    |     | 60     | 141 | 20   |
| cis-1,3-Dichloropropene        | 50          | 0                  | 45.0             | ug/L  | 90   |     | 5    |     | 36     | 161 | 20   |
| 1,1,2-Trichloroethane          | 50          | 0                  | 39.4             | ug/L  | 79   |     | 6    |     | 27     | 175 | 20   |
| 2-Hexanone                     | 250         | 0                  | 170              | ug/L  | 68   | *   | 6    |     | 75     | 131 | 20   |
| Dibromochloromethane           | 50          | 0                  | 41.3             | ug/L  | 83   |     | 6    |     | 59     | 164 | 20   |
| 1,2-Dibromoethane              | 50          | 0                  | 42.9             | ug/L  | 86   |     | 7    |     | 85     | 119 | 20   |
| Tetrachloroethene              | 50          | 0                  | 44.9             | ug/L  | 90   |     | 3    |     | 48     | 153 | 20   |
| Chlorobenzene                  | 50          | 0                  | 45.7             | ug/L  | 91   |     | 3    |     | 85     | 114 | 20   |
| Ethyl Benzene                  | 50          | 650                | 650              | ug/L  | 0    | *   | 200  | *   | 81     | 128 | 20   |
| m/p-Xylenes                    | 100         | 2400               | 2300             | ug/L  | -100 | *   | 0    |     | 69     | 129 | 20   |
| o-Xylene                       | 50          | 2500               | 2500             | ug/L  | 0    | *   | 200  | *   | 75     | 127 | 20   |
| Styrene                        | 50          | 1900               | 1900             | ug/L  | 0    | *   | 200  | *   | 84     | 128 | 20   |
| Bromoform                      | 50          | 0                  | 62.5             | ug/L  | 125  |     | 32   | *   | 73     | 147 | 20   |
| Isopropylbenzene               | 50          | 8.40               | 61.7             | ug/L  | 107  |     | 42   | *   | 76     | 121 | 20   |
| 1,1,2,2-Tetrachloroethane      | 50          | 0                  | 48.4             | ug/L  | 97   |     | 50   | *   | 81     | 131 | 20   |

### Matrix Spike/Matrix Spike Duplicate Summary

SW-846

SDG No.: Q2924

Analytical Method: SW8260D

Client: AECOM

Datafile : VV039006.D

| Parameter                   | Spike | Sample<br>Result | Result | Units | Rec |      | RPD | RPD  |     | Limits |     |
|-----------------------------|-------|------------------|--------|-------|-----|------|-----|------|-----|--------|-----|
|                             |       |                  |        |       | Rec | Qual |     | Qual | Low | High   | RPD |
| 1,3-Dichlorobenzene         | 50    | 0                | 49.8   | ug/L  | 100 |      | 10  |      | 84  | 110    | 20  |
| 1,4-Dichlorobenzene         | 50    | 0                | 46.8   | ug/L  | 94  |      | 7   |      | 81  | 111    | 20  |
| 1,2-Dichlorobenzene         | 50    | 0                | 51.5   | ug/L  | 103 |      | 11  |      | 82  | 113    | 20  |
| 1,2-Dibromo-3-Chloropropane | 50    | 0                | 61.3   | ug/L  | 123 |      | 29  | *    | 79  | 137    | 20  |
| 1,2,4-Trichlorobenzene      | 50    | 0                | 55.9   | ug/L  | 112 |      | 7   |      | 73  | 120    | 20  |
| 1,2,3-Trichlorobenzene      | 50    | 0                | 55.4   | ug/L  | 111 |      | 7   |      | 75  | 119    | 20  |

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

|          |       |        |                    |            |
|----------|-------|--------|--------------------|------------|
| SDG No.: | Q2924 | SW-846 | Analytical Method: | SW8260-Low |
| Client:  | AECOM |        | Datafile :         | VV038985.D |

| Lab Sample ID | Parameter                      | Spike | Result | Unit | Rec | RPD | Qual | Low | Limits High | RPD |
|---------------|--------------------------------|-------|--------|------|-----|-----|------|-----|-------------|-----|
| VV0821WBS01   | Dichlorodifluoromethane        | 20    | 17.3   | ug/L | 86  |     |      | 69  | 116         |     |
|               | Chloromethane                  | 20    | 15.8   | ug/L | 79  |     |      | 65  | 116         |     |
|               | Vinyl chloride                 | 20    | 16.0   | ug/L | 80  |     |      | 65  | 117         |     |
|               | Bromomethane                   | 20    | 19.0   | ug/L | 95  |     |      | 58  | 125         |     |
|               | Chloroethane                   | 20    | 19.4   | ug/L | 97  |     |      | 56  | 128         |     |
|               | Trichlorofluoromethane         | 20    | 19.0   | ug/L | 95  |     |      | 73  | 115         |     |
|               | 1,1,2-Trichlorotrifluoroethane | 20    | 18.6   | ug/L | 93  |     |      | 80  | 112         |     |
|               | 1,1-Dichloroethene             | 20    | 19.2   | ug/L | 96  |     |      | 74  | 110         |     |
|               | Acetone                        | 100   | 94.1   | ug/L | 94  |     |      | 60  | 125         |     |
|               | Carbon disulfide               | 20    | 19.2   | ug/L | 96  |     |      | 64  | 112         |     |
|               | Methyl tert-butyl Ether        | 20    | 20.3   | ug/L | 102 |     |      | 78  | 114         |     |
|               | Methyl Acetate                 | 20    | 20.6   | ug/L | 103 |     |      | 67  | 125         |     |
|               | Methylene Chloride             | 20    | 18.7   | ug/L | 94  |     |      | 72  | 114         |     |
|               | trans-1,2-Dichloroethene       | 20    | 18.9   | ug/L | 95  |     |      | 75  | 108         |     |
|               | 1,1-Dichloroethane             | 20    | 17.8   | ug/L | 89  |     |      | 78  | 112         |     |
|               | Cyclohexane                    | 20    | 18.2   | ug/L | 91  |     |      | 75  | 110         |     |
|               | 2-Butanone                     | 100   | 94.3   | ug/L | 94  |     |      | 65  | 122         |     |
|               | Carbon Tetrachloride           | 20    | 16.5   | ug/L | 83  |     |      | 77  | 113         |     |
|               | cis-1,2-Dichloroethene         | 20    | 17.9   | ug/L | 90  |     |      | 77  | 110         |     |
|               | Bromochloromethane             | 20    | 16.0   | ug/L | 80  |     |      | 70  | 124         |     |
|               | Chloroform                     | 20    | 18.1   | ug/L | 91  |     |      | 79  | 113         |     |
|               | 1,1,1-Trichloroethane          | 20    | 17.6   | ug/L | 88  |     |      | 80  | 108         |     |
|               | Methylcyclohexane              | 20    | 18.7   | ug/L | 94  |     |      | 72  | 115         |     |
|               | Benzene                        | 20    | 19.2   | ug/L | 96  |     |      | 82  | 109         |     |
|               | 1,2-Dichloroethane             | 20    | 18.4   | ug/L | 92  |     |      | 80  | 115         |     |
|               | Trichloroethene                | 20    | 18.0   | ug/L | 90  |     |      | 77  | 113         |     |
|               | 1,2-Dichloropropane            | 20    | 19.4   | ug/L | 97  |     |      | 83  | 111         |     |
|               | Bromodichloromethane           | 20    | 18.1   | ug/L | 91  |     |      | 83  | 110         |     |
|               | 4-Methyl-2-Pentanone           | 100   | 110    | ug/L | 110 |     |      | 74  | 118         |     |
|               | Toluene                        | 20    | 19.4   | ug/L | 97  |     |      | 82  | 110         |     |
|               | t-1,3-Dichloropropene          | 20    | 17.9   | ug/L | 90  |     |      | 79  | 110         |     |
|               | cis-1,3-Dichloropropene        | 20    | 19.5   | ug/L | 98  |     |      | 82  | 110         |     |
|               | 1,1,2-Trichloroethane          | 20    | 17.6   | ug/L | 88  |     |      | 83  | 112         |     |
|               | 2-Hexanone                     | 100   | 82.7   | ug/L | 83  |     |      | 73  | 117         |     |
|               | Dibromochloromethane           | 20    | 17.2   | ug/L | 86  |     |      | 82  | 110         |     |
|               | 1,2-Dibromoethane              | 20    | 18.3   | ug/L | 92  |     |      | 81  | 110         |     |
|               | Tetrachloroethene              | 20    | 18.6   | ug/L | 93  |     |      | 67  | 123         |     |
|               | Chlorobenzene                  | 20    | 18.0   | ug/L | 90  |     |      | 82  | 109         |     |
|               | Ethyl Benzene                  | 20    | 19.1   | ug/L | 96  |     |      | 83  | 109         |     |
|               | m/p-Xylenes                    | 40    | 36.4   | ug/L | 91  |     |      | 82  | 110         |     |
|               | o-Xylene                       | 20    | 19.6   | ug/L | 98  |     |      | 83  | 109         |     |
|               | Styrene                        | 20    | 20.4   | ug/L | 102 |     |      | 80  | 111         |     |
|               | Bromoform                      | 20    | 18.6   | ug/L | 93  |     |      | 79  | 109         |     |
|               | Isopropylbenzene               | 20    | 19.7   | ug/L | 99  |     |      | 83  | 112         |     |
|               | 1,1,2,2-Tetrachloroethane      | 20    | 17.8   | ug/L | 89  |     |      | 76  | 118         |     |
|               | 1,3-Dichlorobenzene            | 20    | 18.6   | ug/L | 93  |     |      | 82  | 108         |     |
|               | 1,4-Dichlorobenzene            | 20    | 18.8   | ug/L | 94  |     |      | 82  | 107         |     |
|               | 1,2-Dichlorobenzene            | 20    | 18.6   | ug/L | 93  |     |      | 82  | 109         |     |

**Laboratory Control Sample/Laboratory Control Sample Duplicate Summary**

**SW-846**

|                 |              |                           |                   |
|-----------------|--------------|---------------------------|-------------------|
| <b>SDG No.:</b> | <u>Q2924</u> | <b>Analytical Method:</b> | <u>SW8260-Low</u> |
| <b>Client:</b>  | <u>AECOM</u> | <b>Datafile :</b>         | <u>VV038985.D</u> |

| Lab Sample ID | Parameter                   | Spike | Result | Unit | Rec | RPD | Qual | Low | Limits | RPD |
|---------------|-----------------------------|-------|--------|------|-----|-----|------|-----|--------|-----|
|               |                             |       |        |      |     |     |      |     | High   |     |
| VV0821WBS01   | 1,2-Dibromo-3-Chloropropane | 20    | 19.1   | ug/L | 96  |     |      | 68  | 112    |     |
|               | 1,2,4-Trichlorobenzene      | 20    | 17.0   | ug/L | 85  |     |      | 75  | 113    |     |
|               | 1,2,3-Trichlorobenzene      | 20    | 18.4   | ug/L | 92  |     |      | 76  | 114    |     |

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

|          |       |        |                    |            |
|----------|-------|--------|--------------------|------------|
| SDG No.: | Q2924 | SW-846 | Analytical Method: | SW8260-Low |
| Client:  | AECOM |        | Datafile :         | VV039011.D |

| Lab Sample ID | Parameter                      | Spike | Result | Unit | Rec | RPD | Qual | Low | Limits High | RPD |
|---------------|--------------------------------|-------|--------|------|-----|-----|------|-----|-------------|-----|
| VV0822WBS01   | Dichlorodifluoromethane        | 20    | 16.0   | ug/L | 80  |     |      | 69  | 116         |     |
|               | Chloromethane                  | 20    | 15.7   | ug/L | 79  |     |      | 65  | 116         |     |
|               | Vinyl chloride                 | 20    | 16.7   | ug/L | 84  |     |      | 65  | 117         |     |
|               | Bromomethane                   | 20    | 19.2   | ug/L | 96  |     |      | 58  | 125         |     |
|               | Chloroethane                   | 20    | 21.3   | ug/L | 106 |     |      | 56  | 128         |     |
|               | Trichlorofluoromethane         | 20    | 18.7   | ug/L | 94  |     |      | 73  | 115         |     |
|               | 1,1,2-Trichlorotrifluoroethane | 20    | 20.1   | ug/L | 101 |     |      | 80  | 112         |     |
|               | 1,1-Dichloroethene             | 20    | 21.7   | ug/L | 109 |     |      | 74  | 110         |     |
|               | Acetone                        | 100   | 120    | ug/L | 120 |     |      | 60  | 125         |     |
|               | Carbon disulfide               | 20    | 22.2   | ug/L | 111 |     |      | 64  | 112         |     |
|               | Methyl tert-butyl Ether        | 20    | 23.6   | ug/L | 118 |     | *    | 78  | 114         |     |
|               | Methyl Acetate                 | 20    | 26.9   | ug/L | 135 |     | *    | 67  | 125         |     |
|               | Methylene Chloride             | 20    | 20.7   | ug/L | 104 |     |      | 72  | 114         |     |
|               | trans-1,2-Dichloroethene       | 20    | 21.7   | ug/L | 109 |     | *    | 75  | 108         |     |
|               | 1,1-Dichloroethane             | 20    | 20.8   | ug/L | 104 |     |      | 78  | 112         |     |
|               | Cyclohexane                    | 20    | 21.4   | ug/L | 107 |     |      | 75  | 110         |     |
|               | 2-Butanone                     | 100   | 120    | ug/L | 120 |     |      | 65  | 122         |     |
|               | Carbon Tetrachloride           | 20    | 16.1   | ug/L | 81  |     |      | 77  | 113         |     |
|               | cis-1,2-Dichloroethene         | 20    | 20.9   | ug/L | 104 |     |      | 77  | 110         |     |
|               | Bromochloromethane             | 20    | 25.1   | ug/L | 126 |     | *    | 70  | 124         |     |
|               | Chloroform                     | 20    | 19.1   | ug/L | 96  |     |      | 79  | 113         |     |
|               | 1,1,1-Trichloroethane          | 20    | 17.7   | ug/L | 89  |     |      | 80  | 108         |     |
|               | Methylcyclohexane              | 20    | 19.8   | ug/L | 99  |     |      | 72  | 115         |     |
|               | Benzene                        | 20    | 19.4   | ug/L | 97  |     |      | 82  | 109         |     |
|               | 1,2-Dichloroethane             | 20    | 17.5   | ug/L | 88  |     |      | 80  | 115         |     |
|               | Trichloroethene                | 20    | 17.6   | ug/L | 88  |     |      | 77  | 113         |     |
|               | 1,2-Dichloropropane            | 20    | 19.5   | ug/L | 98  |     |      | 83  | 111         |     |
|               | Bromodichloromethane           | 20    | 17.6   | ug/L | 88  |     |      | 83  | 110         |     |
|               | 4-Methyl-2-Pentanone           | 100   | 110    | ug/L | 110 |     |      | 74  | 118         |     |
|               | Toluene                        | 20    | 19.6   | ug/L | 98  |     |      | 82  | 110         |     |
|               | t-1,3-Dichloropropene          | 20    | 18.8   | ug/L | 94  |     |      | 79  | 110         |     |
|               | cis-1,3-Dichloropropene        | 20    | 20.1   | ug/L | 101 |     |      | 82  | 110         |     |
|               | 1,1,2-Trichloroethane          | 20    | 17.6   | ug/L | 88  |     |      | 83  | 112         |     |
|               | 2-Hexanone                     | 100   | 85.6   | ug/L | 86  |     |      | 73  | 117         |     |
|               | Dibromochloromethane           | 20    | 16.8   | ug/L | 84  |     |      | 82  | 110         |     |
|               | 1,2-Dibromoethane              | 20    | 18.3   | ug/L | 92  |     |      | 81  | 110         |     |
|               | Tetrachloroethene              | 20    | 18.3   | ug/L | 92  |     |      | 67  | 123         |     |
|               | Chlorobenzene                  | 20    | 18.8   | ug/L | 94  |     |      | 82  | 109         |     |
|               | Ethyl Benzene                  | 20    | 20.7   | ug/L | 104 |     |      | 83  | 109         |     |
|               | m/p-Xylenes                    | 40    | 41.0   | ug/L | 103 |     |      | 82  | 110         |     |
|               | o-Xylene                       | 20    | 20.8   | ug/L | 104 |     |      | 83  | 109         |     |
|               | Styrene                        | 20    | 21.4   | ug/L | 107 |     |      | 80  | 111         |     |
|               | Bromoform                      | 20    | 17.7   | ug/L | 89  |     |      | 79  | 109         |     |
|               | Isopropylbenzene               | 20    | 21.5   | ug/L | 108 |     |      | 83  | 112         |     |
|               | 1,1,2,2-Tetrachloroethane      | 20    | 18.5   | ug/L | 93  |     |      | 76  | 118         |     |
|               | 1,3-Dichlorobenzene            | 20    | 19.6   | ug/L | 98  |     |      | 82  | 108         |     |
|               | 1,4-Dichlorobenzene            | 20    | 18.7   | ug/L | 94  |     |      | 82  | 107         |     |
|               | 1,2-Dichlorobenzene            | 20    | 19.4   | ug/L | 97  |     |      | 82  | 109         |     |



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,  
Fax : 908 789 8922

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

|          |              |                    |                   |
|----------|--------------|--------------------|-------------------|
| SDG No.: | <u>Q2924</u> | Analytical Method: | <u>SW8260-Low</u> |
| Client:  | <u>AECOM</u> | Datafile :         | <u>VV039011.D</u> |

| Lab Sample ID | Parameter                   | Spike | Result | Unit | Rec | RPD | Qual | Low | Limits | RPD |
|---------------|-----------------------------|-------|--------|------|-----|-----|------|-----|--------|-----|
|               |                             |       |        |      |     |     |      |     | High   |     |
| VV0822WBS01   | 1,2-Dibromo-3-Chloropropane | 20    | 21.2   | ug/L | 106 |     |      | 68  | 112    |     |
|               | 1,2,4-Trichlorobenzene      | 20    | 19.9   | ug/L | 100 |     |      | 75  | 113    |     |
|               | 1,2,3-Trichlorobenzene      | 20    | 19.9   | ug/L | 100 |     |      | 76  | 114    |     |

VOLATILE METHOD BLANK SUMMARY

Client ID

VV0821WBL01

Lab Name: Alliance

Contract: AECO02

Lab Code: ACE

SDG NO.: Q2924

Lab File ID: VV038984.D

Lab Sample ID: VV0821WBL01

Date Analyzed: 08/21/2025

Time Analyzed: 16:34

GC Column: DB-624UI ID: 0.18 (mm)

Heated Purge: (Y/N) N

Instrument ID: MSVOA\_V

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

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED |
|-------------------|------------------|----------------|------------------|
| VV0821WBS01       | VV0821WBS01      | VV038985.D     | 08/21/2025       |
| TB                | Q2924-12         | VV038995.D     | 08/21/2025       |
| FB-01-082025      | Q2924-11         | VV038996.D     | 08/21/2025       |
| MW-10C-082025     | Q2924-01         | VV038997.D     | 08/21/2025       |
| MW-2B-082025      | Q2924-05         | VV038998.D     | 08/21/2025       |
| MW-2A-082025      | Q2924-06         | VV038999.D     | 08/21/2025       |
| MW-18C-082025     | Q2924-07         | VV039000.D     | 08/21/2025       |
| MW-16A-082025     | Q2924-08         | VV039001.D     | 08/21/2025       |
| MW-16C-082025     | Q2924-09         | VV039002.D     | 08/21/2025       |
| MW-8A-082025      | Q2924-10         | VV039003.D     | 08/21/2025       |
| MW-201-082025     | Q2924-02         | VV039004.D     | 08/21/2025       |
| MW-201-082025MS   | Q2924-03MS       | VV039005.D     | 08/22/2025       |
| MW-201-082025MSD  | Q2924-04MSD      | VV039006.D     | 08/22/2025       |

COMMENTS: \_\_\_\_\_

\_\_\_\_\_

VOLATILE METHOD BLANK SUMMARY

Client ID

VV0822WBL01

Lab Name: Alliance

Contract: AECO02

Lab Code: ACE

SDG NO.: Q2924

Lab File ID: VV039010.D

Lab Sample ID: VV0822WBL01

Date Analyzed: 08/22/2025

Time Analyzed: 10:55

GC Column: DB-624UI ID: 0.18 (mm)

Heated Purge: (Y/N) N

Instrument ID: MSVOA\_V

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

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED |
|-------------------|------------------|----------------|------------------|
| VV0822WBS01       | VV0822WBS01      | VV039011.D     | 08/22/2025       |
| MW-2B-082025DL    | Q2924-05DL       | VV039014.D     | 08/22/2025       |
| MW-201-082025DL   | Q2924-02DL       | VV039015.D     | 08/22/2025       |
| MW-2B-082025DL2   | Q2924-05DL2      | VV039017.D     | 08/22/2025       |
| MW-201-082025DL2  | Q2924-02DL2      | VV039018.D     | 08/22/2025       |

COMMENTS: \_\_\_\_\_

\_\_\_\_\_

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
BROMOFLUOROBENZENE (BFB)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: VV038972.D  
Instrument ID: MSVOA\_V  
GC Column: DB-624UI ID: 0.18 (mm)

Contract: AECO02  
SDG NO.: Q2924  
BFB Injection Date: 08/21/2025  
BFB Injection Time: 08:34  
Heated Purge: Y/N N

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 50  | 15.0 - 40.0% of mass 95            | 15.6                 |
| 75  | 30.0 - 60.0% of mass 95            | 48.8                 |
| 95  | Base Peak, 100% relative abundance | 100                  |
| 96  | 5.0 - 9.0% of mass 95              | 7                    |
| 173 | Less than 2.0% of mass 174         | 1.5 ( 1.9 ) 1        |
| 174 | 50.0 - 100.0% of mass 95           | 80.3                 |
| 175 | 5.0 - 9.0% of mass 174             | 6.1 ( 7.6 ) 1        |
| 176 | 95.0 - 101.0% of mass 174          | 79.3 ( 98.8 ) 1      |
| 177 | 5.0 - 9.0% of mass 176             | 5.3 ( 6.6 ) 2        |

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| CLIENT ID | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|-----------|---------------|-------------|---------------|---------------|
| VSTDIC001 | VSTDIC001     | VV038973.D  | 08/21/2025    | 09:05         |
| VSTDIC005 | VSTDIC005     | VV038974.D  | 08/21/2025    | 09:48         |
| VSTDIC020 | VSTDIC020     | VV038975.D  | 08/21/2025    | 10:17         |
| VSTDIC050 | VSTDIC050     | VV038976.D  | 08/21/2025    | 10:38         |
| VSTDIC100 | VSTDIC100     | VV038977.D  | 08/21/2025    | 11:02         |
| VSTDIC010 | VSTDIC010     | VV038979.D  | 08/21/2025    | 11:45         |

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
BROMOFLUOROBENZENE (BFB)

Lab Name: Alliance

Contract: AECO02

Lab Code: ACE

SDG NO.: Q2924

Lab File ID: VV038981.D

BFB Injection Date: 08/21/2025

Instrument ID: MSVOA\_V

BFB Injection Time: 15:11

GC Column: DB-624UI ID: 0.18 (mm)

Heated Purge: Y/N N

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 50  | 15.0 - 40.0% of mass 95            | 16                   |
| 75  | 30.0 - 60.0% of mass 95            | 50.9                 |
| 95  | Base Peak, 100% relative abundance | 100                  |
| 96  | 5.0 - 9.0% of mass 95              | 7.4                  |
| 173 | Less than 2.0% of mass 174         | 1.5 ( 1.8 ) 1        |
| 174 | 50.0 - 100.0% of mass 95           | 83.7                 |
| 175 | 5.0 - 9.0% of mass 174             | 6.2 ( 7.4 ) 1        |
| 176 | 95.0 - 101.0% of mass 174          | 81.8 ( 97.8 ) 1      |
| 177 | 5.0 - 9.0% of mass 176             | 5.2 ( 6.4 ) 2        |

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| CLIENT ID        | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|------------------|---------------|-------------|---------------|---------------|
| VSTDCCC050       | VSTDCCC050    | VV038982.D  | 08/21/2025    | 15:42         |
| VV0821WBL01      | VV0821WBL01   | VV038984.D  | 08/21/2025    | 16:34         |
| VV0821WBS01      | VV0821WBS01   | VV038985.D  | 08/21/2025    | 17:09         |
| TB               | Q2924-12      | VV038995.D  | 08/21/2025    | 20:45         |
| FB-01-082025     | Q2924-11      | VV038996.D  | 08/21/2025    | 21:06         |
| MW-10C-082025    | Q2924-01      | VV038997.D  | 08/21/2025    | 21:28         |
| MW-2B-082025     | Q2924-05      | VV038998.D  | 08/21/2025    | 21:49         |
| MW-2A-082025     | Q2924-06      | VV038999.D  | 08/21/2025    | 22:11         |
| MW-18C-082025    | Q2924-07      | VV039000.D  | 08/21/2025    | 22:33         |
| MW-16A-082025    | Q2924-08      | VV039001.D  | 08/21/2025    | 22:54         |
| MW-16C-082025    | Q2924-09      | VV039002.D  | 08/21/2025    | 23:16         |
| MW-8A-082025     | Q2924-10      | VV039003.D  | 08/21/2025    | 23:37         |
| MW-201-082025    | Q2924-02      | VV039004.D  | 08/21/2025    | 23:59         |
| MW-201-082025MS  | Q2924-03MS    | VV039005.D  | 08/22/2025    | 00:20         |
| MW-201-082025MSD | Q2924-04MSD   | VV039006.D  | 08/22/2025    | 00:42         |

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
BROMOFLUOROBENZENE (BFB)

Lab Name: Alliance

Contract: AECO02

Lab Code: ACE

SDG NO.: Q2924

Lab File ID: VV039008.D

BFB Injection Date: 08/22/2025

Instrument ID: MSVOA\_V

BFB Injection Time: 07:56

GC Column: DB-624UI ID: 0.18 (mm)

Heated Purge: Y/N N

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 50  | 15.0 - 40.0% of mass 95            | 17.3                 |
| 75  | 30.0 - 60.0% of mass 95            | 48.3                 |
| 95  | Base Peak, 100% relative abundance | 100                  |
| 96  | 5.0 - 9.0% of mass 95              | 6.6                  |
| 173 | Less than 2.0% of mass 174         | 0.8 ( 1.1 ) 1        |
| 174 | 50.0 - 100.0% of mass 95           | 72                   |
| 175 | 5.0 - 9.0% of mass 174             | 5.8 ( 8 ) 1          |
| 176 | 95.0 - 101.0% of mass 174          | 68.9 ( 95.8 ) 1      |
| 177 | 5.0 - 9.0% of mass 176             | 4.2 ( 6.1 ) 2        |

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| CLIENT ID        | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|------------------|---------------|-------------|---------------|---------------|
| VSTDCCC050       | VSTDCCC050    | VV039009.D  | 08/22/2025    | 09:03         |
| VV0822WBL01      | VV0822WBL01   | VV039010.D  | 08/22/2025    | 10:55         |
| VV0822WBS01      | VV0822WBS01   | VV039011.D  | 08/22/2025    | 11:22         |
| MW-2B-082025DL   | Q2924-05DL    | VV039014.D  | 08/22/2025    | 13:48         |
| MW-201-082025DL  | Q2924-02DL    | VV039015.D  | 08/22/2025    | 14:10         |
| MW-2B-082025DL2  | Q2924-05DL2   | VV039017.D  | 08/22/2025    | 15:06         |
| MW-201-082025DL2 | Q2924-02DL2   | VV039018.D  | 08/22/2025    | 15:27         |

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

|                |                   |                |                   |
|----------------|-------------------|----------------|-------------------|
| Lab Name:      | <u>Alliance</u>   | Contract:      | <u>AECO02</u>     |
| Lab Code:      | <u>ACE</u>        | SDG NO.:       | <u>Q2924</u>      |
| Lab File ID:   | <u>VV038982.D</u> | Date Analyzed: | <u>08/21/2025</u> |
| Instrument ID: | <u>MSVOA_V</u>    | Time Analyzed: | <u>15:42</u>      |
| GC Column:     | <u>DB-624UI</u>   | ID:            | <u>0.18</u> (mm)  |
|                |                   | Heated Purge:  | (Y/N) <u>N</u>    |

|                  | IS1<br>AREA # | RT #  | IS2<br>AREA # | RT #  | IS3<br>AREA # | RT #  |
|------------------|---------------|-------|---------------|-------|---------------|-------|
| 12 HOUR STD      | 531490        | 4.60  | 833989        | 5.53  | 788137        | 8.77  |
| UPPER LIMIT      | 1062980       | 5.096 | 1667980       | 6.029 | 1576270       | 9.273 |
| LOWER LIMIT      | 265745        | 4.096 | 416995        | 5.029 | 394069        | 8.273 |
| EPA SAMPLE NO.   |               |       |               |       |               |       |
| MW-10C-082025    | 584771        | 4.60  | 1064789       | 5.53  | 974990        | 8.78  |
| MW-201-082025    | 456156        | 4.61  | 816696        | 5.54  | 704493        | 8.78  |
| MW-201-082025MS  | 528363        | 4.60  | 848358        | 5.54  | 744435        | 8.78  |
| MW-201-082025MSD | 516721        | 4.60  | 845786        | 5.53  | 728229        | 8.78  |
| MW-2B-082025     | 444034        | 4.60  | 845792        | 5.53  | 758397        | 8.78  |
| MW-2A-082025     | 650572        | 4.60  | 1153105       | 5.53  | 1016294       | 8.78  |
| MW-18C-082025    | 671823        | 4.60  | 1168529       | 5.53  | 1007735       | 8.78  |
| MW-16A-082025    | 676464        | 4.60  | 1214643       | 5.53  | 1082468       | 8.78  |
| MW-16C-082025    | 675326        | 4.60  | 1219427       | 5.53  | 1080769       | 8.78  |
| MW-8A-082025     | 403960        | 4.60  | 696194        | 5.53  | 801104        | 8.78  |
| FB-01-082025     | 585262        | 4.60  | 1084020       | 5.53  | 978319        | 8.78  |
| TB               | 587684        | 4.60  | 1058155       | 5.53  | 949705        | 8.78  |
| VV0821WBL01      | 502511        | 4.60  | 836910        | 5.53  | 730493        | 8.78  |
| VV0821WBS01      | 490499        | 4.60  | 863998        | 5.53  | 809489        | 8.77  |

IS1 = Pentafluorobenzene  
IS2 = 1,4-Difluorobenzene  
IS3 = Chlorobenzene-d5

AREA UPPER LIMIT = +100% of internal standard area  
AREA LOWER LIMIT = -50% of internal standard area  
RT UPPER LIMIT = +0.50 minutes of internal standard RT  
RT LOWER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.  
\* Values outside of QC limits.

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

|                |                   |                |                   |
|----------------|-------------------|----------------|-------------------|
| Lab Name:      | <u>Alliance</u>   | Contract:      | <u>AECO02</u>     |
| Lab Code:      | <u>ACE</u>        | SDG NO.:       | <u>Q2924</u>      |
| Lab File ID:   | <u>VV038982.D</u> | Date Analyzed: | <u>08/21/2025</u> |
| Instrument ID: | <u>MSVOA_V</u>    | Time Analyzed: | <u>15:42</u>      |
| GC Column:     | <u>DB-624UI</u>   | ID:            | <u>0.18</u> (mm)  |
|                |                   | Heated Purge:  | (Y/N) <u>N</u>    |

|                  | IS4<br>AREA # | RT #   |  |  |  |  |
|------------------|---------------|--------|--|--|--|--|
| 12 HOUR STD      | 452177        | 11.172 |  |  |  |  |
| UPPER LIMIT      | 904354        | 11.672 |  |  |  |  |
| LOWER LIMIT      | 226089        | 10.672 |  |  |  |  |
| EPA SAMPLE NO.   |               |        |  |  |  |  |
| MW-10C-082025    | 354625        | 11.18  |  |  |  |  |
| MW-201-082025    | 442841        | 11.18  |  |  |  |  |
| MW-201-082025MS  | 531973        | 11.18  |  |  |  |  |
| MW-201-082025MSD | 478184        | 11.18  |  |  |  |  |
| MW-2B-082025     | 411561        | 11.18  |  |  |  |  |
| MW-2A-082025     | 448609        | 11.18  |  |  |  |  |
| MW-18C-082025    | 449215        | 11.18  |  |  |  |  |
| MW-16A-082025    | 493806        | 11.18  |  |  |  |  |
| MW-16C-082025    | 439427        | 11.17  |  |  |  |  |
| MW-8A-082025     | 375249        | 11.18  |  |  |  |  |
| FB-01-082025     | 424569        | 11.18  |  |  |  |  |
| TB               | 415869        | 11.18  |  |  |  |  |
| VV0821WBL01      | 318475        | 11.18  |  |  |  |  |
| VV0821WBS01      | 439751        | 11.17  |  |  |  |  |

IS4 = 1,4-Dichlorobenzene-d4

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT LOWER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

|                |                   |                |                   |
|----------------|-------------------|----------------|-------------------|
| Lab Name:      | <u>Alliance</u>   | Contract:      | <u>AECO02</u>     |
| Lab Code:      | <u>ACE</u>        | SDG NO.:       | <u>Q2924</u>      |
| Lab File ID:   | <u>VV039009.D</u> | Date Analyzed: | <u>08/22/2025</u> |
| Instrument ID: | <u>MSVOA_V</u>    | Time Analyzed: | <u>09:03</u>      |
| GC Column:     | <u>DB-624UI</u>   | ID:            | <u>0.18</u> (mm)  |
|                |                   | Heated Purge:  | (Y/N) <u>N</u>    |

|                  | IS1<br>AREA # | RT #  | IS2<br>AREA # | RT #  | IS3<br>AREA # | RT #  |
|------------------|---------------|-------|---------------|-------|---------------|-------|
| 12 HOUR STD      | 735484        | 4.60  | 1263800       | 5.53  | 1133330       | 8.77  |
| UPPER LIMIT      | 1470970       | 5.097 | 2527600       | 6.029 | 2266660       | 9.273 |
| LOWER LIMIT      | 367742        | 4.097 | 631899        | 5.029 | 566665        | 8.273 |
| EPA SAMPLE NO.   |               |       |               |       |               |       |
| MW-201-082025DL  | 581936        | 4.60  | 1169048       | 5.53  | 1041283       | 8.78  |
| MW-201-082025DL2 | 766342        | 4.60  | 1437458       | 5.53  | 1305700       | 8.77  |
| MW-2B-082025DL   | 617177        | 4.60  | 1163830       | 5.53  | 1033790       | 8.78  |
| MW-2B-082025DL2  | 582933        | 4.60  | 1168109       | 5.53  | 1042248       | 8.77  |
| VV0822WBL01      | 530331        | 4.60  | 1082642       | 5.53  | 991839        | 8.78  |
| VV0822WBS01      | 711068        | 4.60  | 1223613       | 5.53  | 1111109       | 8.77  |

IS1 = Pentafluorobenzene  
IS2 = 1,4-Difluorobenzene  
IS3 = Chlorobenzene-d5

AREA UPPER LIMIT = +100% of internal standard area  
AREA LOWER LIMIT = -50% of internal standard area  
RT UPPER LIMIT = +0.50 minutes of internal standard RT  
RT LOWER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.  
\* Values outside of QC limits.

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

|                |                   |                |                   |
|----------------|-------------------|----------------|-------------------|
| Lab Name:      | <u>Alliance</u>   | Contract:      | <u>AECO02</u>     |
| Lab Code:      | <u>ACE</u>        | SDG NO.:       | <u>Q2924</u>      |
| Lab File ID:   | <u>VV039009.D</u> | Date Analyzed: | <u>08/22/2025</u> |
| Instrument ID: | <u>MSVOA_V</u>    | Time Analyzed: | <u>09:03</u>      |
| GC Column:     | <u>DB-624UI</u>   | ID:            | <u>0.18</u> (mm)  |
|                |                   | Heated Purge:  | (Y/N) <u>N</u>    |

|                  | IS4<br>AREA # | RT #   |  |  |  |  |
|------------------|---------------|--------|--|--|--|--|
| 12 HOUR STD      | 587247        | 11.172 |  |  |  |  |
| UPPER LIMIT      | 1174490       | 11.672 |  |  |  |  |
| LOWER LIMIT      | 293624        | 10.672 |  |  |  |  |
| EPA SAMPLE NO.   |               |        |  |  |  |  |
| MW-201-082025DL  | 510275        | 11.17  |  |  |  |  |
| MW-201-082025DL2 | 617094        | 11.17  |  |  |  |  |
| MW-2B-082025DL   | 477379        | 11.17  |  |  |  |  |
| MW-2B-082025DL2  | 457197        | 11.17  |  |  |  |  |
| VV0822WBL01      | 416909        | 11.18  |  |  |  |  |
| VV0822WBS01      | 587275        | 11.17  |  |  |  |  |

IS4 = 1,4-Dichlorobenzene-d4

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT LOWER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.



# QC SAMPLE DATA

## Report of Analysis

|                    |                                    |            |                 |    |
|--------------------|------------------------------------|------------|-----------------|----|
| Client:            | AECOM                              |            | Date Collected: |    |
| Project:           | National Grid Equity - Brooklyn NY |            | Date Received:  |    |
| Client Sample ID:  | VV0821WBL01                        | SDG No.:   | Q2924           |    |
| Lab Sample ID:     | VV0821WBL01                        | Matrix:    | Water           |    |
| Analytical Method: | 8260D                              | % Solid:   | 0               |    |
| Sample Wt/Vol:     | 5                                  | Units:     | mL              |    |
| Soil Aliquot Vol:  |                                    |            | uL              |    |
| GC Column:         | DB-624UI                           | ID :       | 0.18            |    |
| Prep Method :      |                                    | Level :    | LOW             |    |
|                    |                                    | Final Vol: | 5000            | uL |
|                    |                                    | Test:      | VOC-TCLVOA-10   |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038984.D        | 1         | 08/21/25 16:34 | VV082125      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 1.00  | U         | 0.22  | 1.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 1.00  | U         | 0.32  | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 1.00  | U         | 0.26  | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 5.00  | U         | 1.40  | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 1.00  | U         | 0.47  | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 1.00  | U         | 0.33  | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 1.00  | U         | 0.25  | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 1.00  | U         | 0.23  | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 5.00  | U         | 1.50  | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 1.00  | U         | 0.21  | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 1.00  | U         | 0.16  | 1.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 1.00  | U         | 0.27  | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 1.00  | U         | 0.28  | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 1.00  | U         | 0.23  | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 1.00  | U         | 0.23  | 1.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 5.00  | U         | 1.50  | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 5.00  | U         | 0.98  | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 1.00  | U         | 0.25  | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 1.00  | U         | 0.19  | 1.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 1.00  | U         | 0.22  | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 1.00  | U         | 0.25  | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 1.00  | U         | 0.20  | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 1.00  | U         | 0.16  | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 1.00  | U         | 0.15  | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 1.00  | U         | 0.22  | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 1.00  | U         | 0.090 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 1.00  | U         | 0.20  | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 1.00  | U         | 0.22  | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 5.00  | U         | 0.68  | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 1.00  | U         | 0.14  | 1.00       | ug/L  |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: |               |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  |               |    |
| Client Sample ID:  | VV0821WBL01                        |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | VV0821WBL01                        |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038984.D        | 1         | 08/21/25 16:34 | VV082125      |

| CAS Number                | Parameter                   | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|--------|-----------|----------|------------|---------|
| 10061-02-6                | t-1,3-Dichloropropene       | 1.00   | U         | 0.17     | 1.00       | ug/L    |
| 10061-01-5                | cis-1,3-Dichloropropene     | 1.00   | U         | 0.16     | 1.00       | ug/L    |
| 79-00-5                   | 1,1,2-Trichloroethane       | 1.00   | U         | 0.21     | 1.00       | ug/L    |
| 591-78-6                  | 2-Hexanone                  | 5.00   | U         | 0.89     | 5.00       | ug/L    |
| 124-48-1                  | Dibromochloromethane        | 1.00   | U         | 0.18     | 1.00       | ug/L    |
| 106-93-4                  | 1,2-Dibromoethane           | 1.00   | U         | 0.15     | 1.00       | ug/L    |
| 127-18-4                  | Tetrachloroethene           | 1.00   | U         | 0.23     | 1.00       | ug/L    |
| 108-90-7                  | Chlorobenzene               | 1.00   | U         | 0.12     | 1.00       | ug/L    |
| 100-41-4                  | Ethyl Benzene               | 1.00   | U         | 0.13     | 1.00       | ug/L    |
| 179601-23-1               | m/p-Xylenes                 | 2.00   | U         | 0.24     | 2.00       | ug/L    |
| 95-47-6                   | o-Xylene                    | 1.00   | U         | 0.12     | 1.00       | ug/L    |
| 100-42-5                  | Styrene                     | 1.00   | U         | 0.15     | 1.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 1.00   | U         | 0.19     | 1.00       | ug/L    |
| 98-82-8                   | Isopropylbenzene            | 1.00   | U         | 0.12     | 1.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 1.00   | U         | 0.26     | 1.00       | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene         | 1.00   | U         | 0.16     | 1.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 1.00   | U         | 0.19     | 1.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 1.00   | U         | 0.16     | 1.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 1.00   | U         | 0.53     | 1.00       | ug/L    |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 1.00   | U         | 0.20     | 1.00       | ug/L    |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 1.00   | U         | 0.20     | 1.00       | ug/L    |
| <b>SURROGATES</b>         |                             |        |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 48.2   |           | 74 - 125 | 96%        | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 43.8   |           | 75 - 124 | 88%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 43.4   |           | 86 - 113 | 87%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 42.3   |           | 77 - 121 | 85%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |        |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 503000 | 4.6       |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 837000 | 5.532     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 730000 | 8.776     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 318000 | 11.175    |          |            |         |

## Report of Analysis

|                    |                                    |           |                 |               |
|--------------------|------------------------------------|-----------|-----------------|---------------|
| Client:            | AECOM                              |           | Date Collected: |               |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  |               |
| Client Sample ID:  | VV0821WBL01                        |           | SDG No.:        | Q2924         |
| Lab Sample ID:     | VV0821WBL01                        |           | Matrix:         | Water         |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000 uL       |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |
| Prep Method :      |                                    |           |                 |               |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038984.D        | 1         | 08/21/25 16:34 | VV082125      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

U = Not Detected  
 LOQ = Limit of Quantitation  
 MDL = Method Detection Limit  
 LOD = Limit of Detection  
 E = Value Exceeds Calibration Range  
 Q = indicates LCS control criteria did not meet requirements  
 M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value  
 B = Analyte Found in Associated Method Blank  
 N = Presumptive Evidence of a Compound  
 \* = Values outside of QC limits  
 D = Dilution  
 () = Laboratory InHouse Limit  
 A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |            |                 |    |
|--------------------|------------------------------------|------------|-----------------|----|
| Client:            | AECOM                              |            | Date Collected: |    |
| Project:           | National Grid Equity - Brooklyn NY |            | Date Received:  |    |
| Client Sample ID:  | VV0822WBL01                        | SDG No.:   | Q2924           |    |
| Lab Sample ID:     | VV0822WBL01                        | Matrix:    | Water           |    |
| Analytical Method: | 8260D                              | % Solid:   | 0               |    |
| Sample Wt/Vol:     | 5                                  | Units:     | mL              |    |
| Soil Aliquot Vol:  |                                    |            | uL              |    |
| GC Column:         | DB-624UI                           | ID :       | 0.18            |    |
| Prep Method :      |                                    | Level :    | LOW             |    |
|                    |                                    | Final Vol: | 5000            | uL |
|                    |                                    | Test:      | VOC-TCLVOA-10   |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039010.D        | 1         | 08/22/25 10:55 | VV082225      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 1.00  | U         | 0.22  | 1.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 1.00  | U         | 0.32  | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 1.00  | U         | 0.26  | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 5.00  | U         | 1.40  | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 1.00  | U         | 0.47  | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 1.00  | U         | 0.33  | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 1.00  | U         | 0.25  | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 1.00  | U         | 0.23  | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 5.00  | U         | 1.50  | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 1.00  | U         | 0.21  | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 1.00  | U         | 0.16  | 1.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 1.00  | U         | 0.27  | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 1.00  | U         | 0.28  | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 1.00  | U         | 0.23  | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 1.00  | U         | 0.23  | 1.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 5.00  | U         | 1.50  | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 5.00  | U         | 0.98  | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 1.00  | U         | 0.25  | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 1.00  | U         | 0.19  | 1.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 1.00  | U         | 0.22  | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 1.00  | U         | 0.25  | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 1.00  | U         | 0.20  | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 1.00  | U         | 0.16  | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 1.00  | U         | 0.15  | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 1.00  | U         | 0.22  | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 1.00  | U         | 0.090 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 1.00  | U         | 0.20  | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 1.00  | U         | 0.22  | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 5.00  | U         | 0.68  | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 1.00  | U         | 0.14  | 1.00       | ug/L  |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: |               |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  |               |    |
| Client Sample ID:  | VV0822WBL01                        |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | VV0822WBL01                        |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039010.D        | 1         | 08/22/25 10:55 | VV082225      |

| CAS Number                | Parameter                   | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|---------|-----------|----------|------------|---------|
| 10061-02-6                | t-1,3-Dichloropropene       | 1.00    | U         | 0.17     | 1.00       | ug/L    |
| 10061-01-5                | cis-1,3-Dichloropropene     | 1.00    | U         | 0.16     | 1.00       | ug/L    |
| 79-00-5                   | 1,1,2-Trichloroethane       | 1.00    | U         | 0.21     | 1.00       | ug/L    |
| 591-78-6                  | 2-Hexanone                  | 5.00    | U         | 0.89     | 5.00       | ug/L    |
| 124-48-1                  | Dibromochloromethane        | 1.00    | U         | 0.18     | 1.00       | ug/L    |
| 106-93-4                  | 1,2-Dibromoethane           | 1.00    | U         | 0.15     | 1.00       | ug/L    |
| 127-18-4                  | Tetrachloroethene           | 1.00    | U         | 0.23     | 1.00       | ug/L    |
| 108-90-7                  | Chlorobenzene               | 1.00    | U         | 0.12     | 1.00       | ug/L    |
| 100-41-4                  | Ethyl Benzene               | 1.00    | U         | 0.13     | 1.00       | ug/L    |
| 179601-23-1               | m/p-Xylenes                 | 2.00    | U         | 0.24     | 2.00       | ug/L    |
| 95-47-6                   | o-Xylene                    | 1.00    | U         | 0.12     | 1.00       | ug/L    |
| 100-42-5                  | Styrene                     | 1.00    | U         | 0.15     | 1.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 1.00    | U         | 0.19     | 1.00       | ug/L    |
| 98-82-8                   | Isopropylbenzene            | 1.00    | U         | 0.12     | 1.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 1.00    | U         | 0.26     | 1.00       | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene         | 1.00    | U         | 0.16     | 1.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 1.00    | U         | 0.19     | 1.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 1.00    | U         | 0.16     | 1.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 1.00    | U         | 0.53     | 1.00       | ug/L    |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 1.00    | U         | 0.20     | 1.00       | ug/L    |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 1.00    | U         | 0.20     | 1.00       | ug/L    |
| <b>SURROGATES</b>         |                             |         |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 60.0    |           | 74 - 125 | 120%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 47.4    |           | 75 - 124 | 95%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 47.7    |           | 86 - 113 | 95%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 49.3    |           | 77 - 121 | 99%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |         |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 530000  | 4.603     |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 1080000 | 5.532     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 992000  | 8.776     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 417000  | 11.175    |          |            |         |

## Report of Analysis

|                    |                                    |           |                 |               |
|--------------------|------------------------------------|-----------|-----------------|---------------|
| Client:            | AECOM                              |           | Date Collected: |               |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  |               |
| Client Sample ID:  | VV0822WBL01                        |           | SDG No.:        | Q2924         |
| Lab Sample ID:     | VV0822WBL01                        |           | Matrix:         | Water         |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000 uL       |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |
| Prep Method :      |                                    |           |                 |               |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039010.D        | 1         | 08/22/25 10:55 | VV082225      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

U = Not Detected  
 LOQ = Limit of Quantitation  
 MDL = Method Detection Limit  
 LOD = Limit of Detection  
 E = Value Exceeds Calibration Range  
 Q = indicates LCS control criteria did not meet requirements  
 M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value  
 B = Analyte Found in Associated Method Blank  
 N = Presumptive Evidence of a Compound  
 \* = Values outside of QC limits  
 D = Dilution  
 () = Laboratory InHouse Limit  
 A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: |               |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  |               |    |
| Client Sample ID:  | VV0821WBS01                        |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | VV0821WBS01                        |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038985.D        | 1         | 08/21/25 17:09 | VV082125      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 17.3  |           | 0.22  | 1.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 15.8  |           | 0.32  | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 16.0  |           | 0.26  | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 19.0  |           | 1.40  | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 19.4  |           | 0.47  | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 19.0  |           | 0.33  | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 18.6  |           | 0.25  | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 19.2  |           | 0.23  | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 94.1  |           | 1.50  | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 19.2  |           | 0.21  | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 20.3  |           | 0.16  | 1.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 20.6  |           | 0.27  | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 18.7  |           | 0.28  | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 18.9  |           | 0.23  | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 17.8  |           | 0.23  | 1.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 18.2  |           | 1.50  | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 94.3  |           | 0.98  | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 16.5  |           | 0.25  | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 17.9  |           | 0.19  | 1.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 16.0  |           | 0.22  | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 18.1  |           | 0.25  | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 17.6  |           | 0.20  | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 18.7  |           | 0.16  | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 19.2  |           | 0.15  | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 18.4  |           | 0.22  | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 18.0  |           | 0.090 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 19.4  |           | 0.20  | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 18.1  |           | 0.22  | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 110   |           | 0.68  | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 19.4  |           | 0.14  | 1.00       | ug/L  |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: |               |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  |               |    |
| Client Sample ID:  | VV0821WBS01                        |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | VV0821WBS01                        |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038985.D        | 1         | 08/21/25 17:09 | VV082125      |

| CAS Number                | Parameter                   | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|--------|-----------|----------|------------|---------|
| 10061-02-6                | t-1,3-Dichloropropene       | 17.9   |           | 0.17     | 1.00       | ug/L    |
| 10061-01-5                | cis-1,3-Dichloropropene     | 19.5   |           | 0.16     | 1.00       | ug/L    |
| 79-00-5                   | 1,1,2-Trichloroethane       | 17.6   |           | 0.21     | 1.00       | ug/L    |
| 591-78-6                  | 2-Hexanone                  | 82.7   |           | 0.89     | 5.00       | ug/L    |
| 124-48-1                  | Dibromochloromethane        | 17.2   |           | 0.18     | 1.00       | ug/L    |
| 106-93-4                  | 1,2-Dibromoethane           | 18.3   |           | 0.15     | 1.00       | ug/L    |
| 127-18-4                  | Tetrachloroethene           | 18.6   |           | 0.23     | 1.00       | ug/L    |
| 108-90-7                  | Chlorobenzene               | 18.0   |           | 0.12     | 1.00       | ug/L    |
| 100-41-4                  | Ethyl Benzene               | 19.1   |           | 0.13     | 1.00       | ug/L    |
| 179601-23-1               | m/p-Xylenes                 | 36.4   |           | 0.24     | 2.00       | ug/L    |
| 95-47-6                   | o-Xylene                    | 19.6   |           | 0.12     | 1.00       | ug/L    |
| 100-42-5                  | Styrene                     | 20.4   |           | 0.15     | 1.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 18.6   |           | 0.19     | 1.00       | ug/L    |
| 98-82-8                   | Isopropylbenzene            | 19.7   |           | 0.12     | 1.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 17.8   |           | 0.26     | 1.00       | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene         | 18.6   |           | 0.16     | 1.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 18.8   |           | 0.19     | 1.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 18.6   |           | 0.16     | 1.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 19.1   |           | 0.53     | 1.00       | ug/L    |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 17.0   |           | 0.20     | 1.00       | ug/L    |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 18.4   |           | 0.20     | 1.00       | ug/L    |
| <b>SURROGATES</b>         |                             |        |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 55.0   |           | 74 - 125 | 110%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 44.8   |           | 75 - 124 | 90%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 49.6   |           | 86 - 113 | 99%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 48.8   |           | 77 - 121 | 98%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |        |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 490000 | 4.597     |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 864000 | 5.529     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 809000 | 8.773     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 440000 | 11.172    |          |            |         |

## Report of Analysis

|                    |                                    |           |                 |               |
|--------------------|------------------------------------|-----------|-----------------|---------------|
| Client:            | AECOM                              |           | Date Collected: |               |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  |               |
| Client Sample ID:  | VV0821WBS01                        |           | SDG No.:        | Q2924         |
| Lab Sample ID:     | VV0821WBS01                        |           | Matrix:         | Water         |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000 uL       |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |
| Prep Method :      |                                    |           |                 |               |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV038985.D        | 1         | 08/21/25 17:09 | VV082125      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

U = Not Detected  
 LOQ = Limit of Quantitation  
 MDL = Method Detection Limit  
 LOD = Limit of Detection  
 E = Value Exceeds Calibration Range  
 Q = indicates LCS control criteria did not meet requirements  
 M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value  
 B = Analyte Found in Associated Method Blank  
 N = Presumptive Evidence of a Compound  
 \* = Values outside of QC limits  
 D = Dilution  
 () = Laboratory InHouse Limit  
 A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: |               |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  |               |    |
| Client Sample ID:  | VV0822WBS01                        |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | VV0822WBS01                        |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039011.D        | 1         | 08/22/25 11:22 | VV082225      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 16.0  |           | 0.22  | 1.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 15.7  |           | 0.32  | 1.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 16.7  |           | 0.26  | 1.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 19.2  |           | 1.40  | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 21.3  |           | 0.47  | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 18.7  |           | 0.33  | 1.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 20.1  |           | 0.25  | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 21.7  |           | 0.23  | 1.00       | ug/L  |
| 67-64-1        | Acetone                        | 120   |           | 1.50  | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 22.2  |           | 0.21  | 1.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 23.6  |           | 0.16  | 1.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 26.9  |           | 0.27  | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 20.7  |           | 0.28  | 1.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 21.7  |           | 0.23  | 1.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 20.8  |           | 0.23  | 1.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 21.4  |           | 1.50  | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 120   |           | 0.98  | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 16.1  |           | 0.25  | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 20.9  |           | 0.19  | 1.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 25.1  |           | 0.22  | 1.00       | ug/L  |
| 67-66-3        | Chloroform                     | 19.1  |           | 0.25  | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 17.7  |           | 0.20  | 1.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 19.8  |           | 0.16  | 1.00       | ug/L  |
| 71-43-2        | Benzene                        | 19.4  |           | 0.15  | 1.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 17.5  |           | 0.22  | 1.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 17.6  |           | 0.090 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 19.5  |           | 0.20  | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 17.6  |           | 0.22  | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 110   |           | 0.68  | 5.00       | ug/L  |
| 108-88-3       | Toluene                        | 19.6  |           | 0.14  | 1.00       | ug/L  |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: |               |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  |               |    |
| Client Sample ID:  | VV0822WBS01                        |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | VV0822WBS01                        |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039011.D        | 1         | 08/22/25 11:22 | VV082225      |

| CAS Number                | Parameter                   | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|---------|-----------|----------|------------|---------|
| 10061-02-6                | t-1,3-Dichloropropene       | 18.8    |           | 0.17     | 1.00       | ug/L    |
| 10061-01-5                | cis-1,3-Dichloropropene     | 20.1    |           | 0.16     | 1.00       | ug/L    |
| 79-00-5                   | 1,1,2-Trichloroethane       | 17.6    |           | 0.21     | 1.00       | ug/L    |
| 591-78-6                  | 2-Hexanone                  | 85.6    |           | 0.89     | 5.00       | ug/L    |
| 124-48-1                  | Dibromochloromethane        | 16.8    |           | 0.18     | 1.00       | ug/L    |
| 106-93-4                  | 1,2-Dibromoethane           | 18.3    |           | 0.15     | 1.00       | ug/L    |
| 127-18-4                  | Tetrachloroethene           | 18.3    |           | 0.23     | 1.00       | ug/L    |
| 108-90-7                  | Chlorobenzene               | 18.8    |           | 0.12     | 1.00       | ug/L    |
| 100-41-4                  | Ethyl Benzene               | 20.7    |           | 0.13     | 1.00       | ug/L    |
| 179601-23-1               | m/p-Xylenes                 | 41.0    |           | 0.24     | 2.00       | ug/L    |
| 95-47-6                   | o-Xylene                    | 20.8    |           | 0.12     | 1.00       | ug/L    |
| 100-42-5                  | Styrene                     | 21.4    |           | 0.15     | 1.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 17.7    |           | 0.19     | 1.00       | ug/L    |
| 98-82-8                   | Isopropylbenzene            | 21.5    |           | 0.12     | 1.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 18.5    |           | 0.26     | 1.00       | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene         | 19.6    |           | 0.16     | 1.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 18.7    |           | 0.19     | 1.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 19.4    |           | 0.16     | 1.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 21.2    |           | 0.53     | 1.00       | ug/L    |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 19.9    |           | 0.20     | 1.00       | ug/L    |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 19.9    |           | 0.20     | 1.00       | ug/L    |
| <b>SURROGATES</b>         |                             |         |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 48.5    |           | 74 - 125 | 97%        | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 46.0    |           | 75 - 124 | 92%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 46.3    |           | 86 - 113 | 93%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 48.5    |           | 77 - 121 | 97%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |         |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 711000  | 4.6       |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 1220000 | 5.532     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 1110000 | 8.773     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 587000  | 11.172    |          |            |         |

## Report of Analysis

|                    |                                    |           |                 |               |
|--------------------|------------------------------------|-----------|-----------------|---------------|
| Client:            | AECOM                              |           | Date Collected: |               |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  |               |
| Client Sample ID:  | VV0822WBS01                        |           | SDG No.:        | Q2924         |
| Lab Sample ID:     | VV0822WBS01                        |           | Matrix:         | Water         |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000 uL       |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |
| Prep Method :      |                                    |           |                 |               |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039011.D        | 1         | 08/22/25 11:22 | VV082225      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

U = Not Detected  
 LOQ = Limit of Quantitation  
 MDL = Method Detection Limit  
 LOD = Limit of Detection  
 E = Value Exceeds Calibration Range  
 Q = indicates LCS control criteria did not meet requirements  
 M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value  
 B = Analyte Found in Associated Method Blank  
 N = Presumptive Evidence of a Compound  
 \* = Values outside of QC limits  
 D = Dilution  
 () = Laboratory InHouse Limit  
 A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-201-082025MS                    |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-03MS                         |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039005.D        | 1         | 08/22/25 00:20 | VV082125      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 38.4  |           | 0.22  | 5.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 32.4  |           | 0.32  | 5.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 39.7  |           | 0.26  | 5.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 24.9  |           | 1.40  | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 52.0  |           | 0.47  | 5.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 45.7  |           | 0.33  | 5.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 44.6  |           | 0.25  | 5.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 49.2  |           | 0.23  | 5.00       | ug/L  |
| 67-64-1        | Acetone                        | 210   |           | 1.50  | 25.0       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 47.5  |           | 0.21  | 5.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 55.4  |           | 0.16  | 5.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 47.7  |           | 0.27  | 5.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 43.3  |           | 0.28  | 5.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 220   | E         | 0.23  | 5.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 42.5  |           | 0.23  | 5.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 41.1  |           | 1.50  | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 230   |           | 0.98  | 25.0       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 42.2  |           | 0.25  | 5.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 170   | E         | 0.19  | 5.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 45.2  |           | 0.22  | 5.00       | ug/L  |
| 67-66-3        | Chloroform                     | 41.8  |           | 0.25  | 5.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 42.5  |           | 0.20  | 5.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 45.2  |           | 0.16  | 5.00       | ug/L  |
| 71-43-2        | Benzene                        | 1600  | E         | 0.15  | 5.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 180   | E         | 0.22  | 5.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 74.0  |           | 0.090 | 5.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 40.8  |           | 0.20  | 5.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 40.6  |           | 0.22  | 5.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 220   |           | 0.68  | 25.0       | ug/L  |
| 108-88-3       | Toluene                        | 3300  | E         | 0.14  | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-201-082025MS                    |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-03MS                         |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039005.D        | 1         | 08/22/25 00:20 | VV082125      |

| CAS Number                | Parameter                   | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|--------|-----------|----------|------------|---------|
| 10061-02-6                | t-1,3-Dichloropropene       | 41.0   |           | 0.17     | 5.00       | ug/L    |
| 10061-01-5                | cis-1,3-Dichloropropene     | 42.6   |           | 0.16     | 5.00       | ug/L    |
| 79-00-5                   | 1,1,2-Trichloroethane       | 37.2   |           | 0.21     | 5.00       | ug/L    |
| 591-78-6                  | 2-Hexanone                  | 160    |           | 0.89     | 25.0       | ug/L    |
| 124-48-1                  | Dibromochloromethane        | 38.7   |           | 0.18     | 5.00       | ug/L    |
| 106-93-4                  | 1,2-Dibromoethane           | 39.9   |           | 0.15     | 5.00       | ug/L    |
| 127-18-4                  | Tetrachloroethene           | 43.7   |           | 0.23     | 5.00       | ug/L    |
| 108-90-7                  | Chlorobenzene               | 44.2   |           | 0.12     | 5.00       | ug/L    |
| 100-41-4                  | Ethyl Benzene               | 650    | E         | 0.13     | 5.00       | ug/L    |
| 179601-23-1               | m/p-Xylenes                 | 2300   | E         | 0.24     | 10.0       | ug/L    |
| 95-47-6                   | o-Xylene                    | 2200   | E         | 0.12     | 5.00       | ug/L    |
| 100-42-5                  | Styrene                     | 1800   | E         | 0.15     | 5.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 45.4   |           | 0.19     | 5.00       | ug/L    |
| 98-82-8                   | Isopropylbenzene            | 43.6   |           | 0.12     | 5.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 28.9   |           | 0.26     | 5.00       | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene         | 45.0   |           | 0.16     | 5.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 43.8   |           | 0.19     | 5.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 46.1   |           | 0.16     | 5.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 46.0   |           | 0.53     | 5.00       | ug/L    |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 51.9   |           | 0.20     | 5.00       | ug/L    |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 51.6   |           | 0.20     | 5.00       | ug/L    |
| <b>SURROGATES</b>         |                             |        |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 50.4   |           | 74 - 125 | 101%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 44.4   |           | 75 - 124 | 89%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 43.9   |           | 86 - 113 | 88%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 45.0   |           | 77 - 121 | 90%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |        |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 528000 | 4.603     |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 848000 | 5.535     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 744000 | 8.78      |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 532000 | 11.178    |          |            |         |



## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-201-082025MSD                   |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-04MSD                        |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039006.D        | 1         | 08/22/25 00:42 | VV082125      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL   | LOQ / CRQL | Units |
|----------------|--------------------------------|-------|-----------|-------|------------|-------|
| <b>TARGETS</b> |                                |       |           |       |            |       |
| 75-71-8        | Dichlorodifluoromethane        | 44.7  |           | 0.22  | 5.00       | ug/L  |
| 74-87-3        | Chloromethane                  | 49.3  |           | 0.32  | 5.00       | ug/L  |
| 75-01-4        | Vinyl Chloride                 | 57.5  |           | 0.26  | 5.00       | ug/L  |
| 74-83-9        | Bromomethane                   | 31.1  |           | 1.40  | 5.00       | ug/L  |
| 75-00-3        | Chloroethane                   | 60.7  |           | 0.47  | 5.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane         | 47.5  |           | 0.33  | 5.00       | ug/L  |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 44.0  |           | 0.25  | 5.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene             | 54.6  |           | 0.23  | 5.00       | ug/L  |
| 67-64-1        | Acetone                        | 260   |           | 1.50  | 25.0       | ug/L  |
| 75-15-0        | Carbon Disulfide               | 55.4  |           | 0.21  | 5.00       | ug/L  |
| 1634-04-4      | Methyl tert-butyl Ether        | 67.5  |           | 0.16  | 5.00       | ug/L  |
| 79-20-9        | Methyl Acetate                 | 61.9  |           | 0.27  | 5.00       | ug/L  |
| 75-09-2        | Methylene Chloride             | 53.6  |           | 0.28  | 5.00       | ug/L  |
| 156-60-5       | trans-1,2-Dichloroethene       | 250   | E         | 0.23  | 5.00       | ug/L  |
| 75-34-3        | 1,1-Dichloroethane             | 49.1  |           | 0.23  | 5.00       | ug/L  |
| 110-82-7       | Cyclohexane                    | 38.3  |           | 1.50  | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                     | 260   |           | 0.98  | 25.0       | ug/L  |
| 56-23-5        | Carbon Tetrachloride           | 41.8  |           | 0.25  | 5.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene         | 180   | E         | 0.19  | 5.00       | ug/L  |
| 74-97-5        | Bromochloromethane             | 47.4  |           | 0.22  | 5.00       | ug/L  |
| 67-66-3        | Chloroform                     | 44.9  |           | 0.25  | 5.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane          | 43.5  |           | 0.20  | 5.00       | ug/L  |
| 108-87-2       | Methylcyclohexane              | 38.9  |           | 0.16  | 5.00       | ug/L  |
| 71-43-2        | Benzene                        | 1600  | E         | 0.15  | 5.00       | ug/L  |
| 107-06-2       | 1,2-Dichloroethane             | 180   | E         | 0.22  | 5.00       | ug/L  |
| 79-01-6        | Trichloroethene                | 72.9  |           | 0.090 | 5.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane            | 43.1  |           | 0.20  | 5.00       | ug/L  |
| 75-27-4        | Bromodichloromethane           | 43.1  |           | 0.22  | 5.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone           | 230   |           | 0.68  | 25.0       | ug/L  |
| 108-88-3       | Toluene                        | 3200  | E         | 0.14  | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |           |                 |               |    |
|--------------------|------------------------------------|-----------|-----------------|---------------|----|
| Client:            | AECOM                              |           | Date Collected: | 08/20/25      |    |
| Project:           | National Grid Equity - Brooklyn NY |           | Date Received:  | 08/20/25      |    |
| Client Sample ID:  | MW-201-082025MSD                   |           | SDG No.:        | Q2924         |    |
| Lab Sample ID:     | Q2924-04MSD                        |           | Matrix:         | Water         |    |
| Analytical Method: | 8260D                              |           | % Solid:        | 0             |    |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000          | uL |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW           |    |
| Prep Method :      |                                    |           |                 |               |    |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039006.D        | 1         | 08/22/25 00:42 | VV082125      |

| CAS Number                | Parameter                   | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|--------|-----------|----------|------------|---------|
| 10061-02-6                | t-1,3-Dichloropropene       | 43.6   |           | 0.17     | 5.00       | ug/L    |
| 10061-01-5                | cis-1,3-Dichloropropene     | 45.0   |           | 0.16     | 5.00       | ug/L    |
| 79-00-5                   | 1,1,2-Trichloroethane       | 39.4   |           | 0.21     | 5.00       | ug/L    |
| 591-78-6                  | 2-Hexanone                  | 170    |           | 0.89     | 25.0       | ug/L    |
| 124-48-1                  | Dibromochloromethane        | 41.3   |           | 0.18     | 5.00       | ug/L    |
| 106-93-4                  | 1,2-Dibromoethane           | 42.9   |           | 0.15     | 5.00       | ug/L    |
| 127-18-4                  | Tetrachloroethene           | 44.9   |           | 0.23     | 5.00       | ug/L    |
| 108-90-7                  | Chlorobenzene               | 45.7   |           | 0.12     | 5.00       | ug/L    |
| 100-41-4                  | Ethyl Benzene               | 650    | E         | 0.13     | 5.00       | ug/L    |
| 179601-23-1               | m/p-Xylenes                 | 2300   | E         | 0.24     | 10.0       | ug/L    |
| 95-47-6                   | o-Xylene                    | 2500   | E         | 0.12     | 5.00       | ug/L    |
| 100-42-5                  | Styrene                     | 1900   | E         | 0.15     | 5.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 62.5   |           | 0.19     | 5.00       | ug/L    |
| 98-82-8                   | Isopropylbenzene            | 61.7   |           | 0.12     | 5.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 48.4   |           | 0.26     | 5.00       | ug/L    |
| 541-73-1                  | 1,3-Dichlorobenzene         | 49.8   |           | 0.16     | 5.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 46.8   |           | 0.19     | 5.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 51.5   |           | 0.16     | 5.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 61.3   |           | 0.53     | 5.00       | ug/L    |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 55.9   |           | 0.20     | 5.00       | ug/L    |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 55.4   |           | 0.20     | 5.00       | ug/L    |
| <b>SURROGATES</b>         |                             |        |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 52.6   |           | 74 - 125 | 105%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 47.2   |           | 75 - 124 | 94%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 44.3   |           | 86 - 113 | 89%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 60.7   |           | 77 - 121 | 121%       | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |        |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 517000 | 4.603     |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 846000 | 5.532     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 728000 | 8.78      |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 478000 | 11.178    |          |            |         |





# CALIBRATION SUMMARY

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

|                                                 |                                                          |
|-------------------------------------------------|----------------------------------------------------------|
| Lab Name: <u>Alliance</u>                       | Contract: <u>AECO02</u>                                  |
| Lab Code: <u>ACE</u>                            | SDG No.: <u>Q2924</u>                                    |
| Instrument ID: <u>MSVOA_V</u>                   | Calibration Date(s): <u>08/21/2025</u> <u>08/21/2025</u> |
| Heated Purge: (Y/N) <u>N</u>                    | Calibration Time(s): <u>09:05</u> <u>11:45</u>           |
| GC Column: <u>DB-624UI</u> ID: <u>0.18</u> (mm) |                                                          |

|                                |        |                     |        |                     |        |                     |       |       |
|--------------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|
| LAB FILE ID:                   |        | RRF001 = VV038973.D |        | RRF005 = VV038974.D |        | RRF020 = VV038975.D |       |       |
|                                |        | RRF050 = VV038976.D |        | RRF100 = VV038977.D |        | RRF010 = VV038979.D |       |       |
| COMPOUND                       | RRF001 | RRF005              | RRF020 | RRF050              | RRF100 | RRF010              | RRF   | % RSD |
| Dichlorodifluoromethane        | 0.952  | 0.756               | 0.675  | 0.722               | 0.633  | 0.747               | 0.747 | 14.7  |
| Chloromethane                  | 0.878  | 0.853               | 0.712  | 0.768               | 0.654  | 0.769               | 0.772 | 10.9  |
| Vinyl Chloride                 | 0.985  | 0.849               | 0.688  | 0.823               | 0.714  | 0.829               | 0.815 | 13.1  |
| Bromomethane                   |        | 0.459               | 0.389  | 0.448               | 0.423  | 0.482               | 0.440 | 8.1   |
| Chloroethane                   | 0.346  | 0.559               | 0.365  | 0.414               | 0.359  | 0.448               | 0.415 | 19.3  |
| Trichlorofluoromethane         | 1.342  | 1.178               | 1.026  | 1.088               | 0.961  | 1.102               | 1.116 | 11.9  |
| 1,1,2-Trichlorotrifluoroethane | 0.823  | 0.628               | 0.558  | 0.554               | 0.513  | 0.619               | 0.616 | 17.9  |
| 1,1-Dichloroethene             | 0.675  | 0.593               | 0.519  | 0.536               | 0.495  | 0.586               | 0.567 | 11.4  |
| Acetone                        | 0.390  | 0.264               | 0.261  | 0.249               | 0.242  | 0.284               | 0.282 | 19.5  |
| Carbon Disulfide               | 1.825  | 1.631               | 1.475  | 1.476               | 1.354  | 1.709               | 1.578 | 11.1  |
| Methyl tert-butyl Ether        | 1.629  | 1.485               | 1.792  | 1.715               | 1.614  | 1.832               | 1.678 | 7.6   |
| Methyl Acetate                 | 0.610  | 0.589               | 0.638  | 0.560               | 0.554  | 0.630               | 0.597 | 5.9   |
| Methylene Chloride             | 0.805  | 0.659               | 0.694  | 0.581               | 0.519  | 0.668               | 0.654 | 15    |
| trans-1,2-Dichloroethene       | 0.577  | 0.577               | 0.640  | 0.573               | 0.529  | 0.651               | 0.591 | 7.8   |
| 1,1-Dichloroethane             | 1.293  | 1.028               | 1.122  | 0.940               | 0.941  | 1.288               | 1.102 | 14.6  |
| Cyclohexane                    |        | 1.084               | 0.955  | 0.808               | 0.961  | 1.109               | 0.983 | 12.2  |
| 2-Butanone                     | 0.285  | 0.336               | 0.397  | 0.332               | 0.406  | 0.442               | 0.366 | 15.9  |
| Carbon Tetrachloride           | 0.725  | 0.726               | 0.608  | 0.616               | 0.577  | 0.660               | 0.652 | 9.7   |
| cis-1,2-Dichloroethene         | 0.725  | 0.710               | 0.686  | 0.640               | 0.712  | 0.806               | 0.713 | 7.6   |
| Bromochloromethane             | 0.463  | 0.419               | 0.283  | 0.221               | 0.266  | 0.271               | 0.320 | 30.2  |
| Chloroform                     | 1.373  | 1.345               | 1.244  | 1.096               | 1.150  | 1.347               | 1.259 | 9.2   |
| 1,1,1-Trichloroethane          | 1.408  | 1.248               | 1.107  | 1.038               | 1.051  | 1.246               | 1.183 | 12.1  |
| Methylcyclohexane              | 0.574  | 0.590               | 0.612  | 0.632               | 0.677  | 0.649               | 0.622 | 6.1   |
| Benzene                        | 1.384  | 1.558               | 1.602  | 1.430               | 1.510  | 1.667               | 1.525 | 7     |
| 1,2-Dichloroethane             | 0.602  | 0.597               | 0.546  | 0.514               | 0.506  | 0.580               | 0.558 | 7.5   |
| Trichloroethene                | 0.444  | 0.433               | 0.392  | 0.391               | 0.389  | 0.433               | 0.414 | 6.2   |
| 1,2-Dichloropropane            | 0.341  | 0.335               | 0.397  | 0.330               | 0.366  | 0.425               | 0.366 | 10.5  |
| Bromodichloromethane           | 0.686  | 0.635               | 0.616  | 0.584               | 0.571  | 0.657               | 0.625 | 7     |
| 4-Methyl-2-Pentanone           | 0.327  | 0.440               | 0.525  | 0.505               | 0.514  | 0.549               | 0.477 | 17.1  |
| Toluene                        | 0.791  | 0.916               | 1.009  | 1.070               | 0.984  | 1.026               | 0.966 | 10.3  |

\* Compounds with required minimum RRF and maximum %RSD values.  
All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

|                                                 |                                                          |
|-------------------------------------------------|----------------------------------------------------------|
| Lab Name: <u>Alliance</u>                       | Contract: <u>AECO02</u>                                  |
| Lab Code: <u>ACE</u>                            | SDG No.: <u>Q2924</u>                                    |
| Instrument ID: <u>MSVOA_V</u>                   | Calibration Date(s): <u>08/21/2025</u> <u>08/21/2025</u> |
| Heated Purge: (Y/N) <u>N</u>                    | Calibration Time(s): <u>09:05</u> <u>11:45</u>           |
| GC Column: <u>DB-624UI</u> ID: <u>0.18</u> (mm) |                                                          |

| LAB FILE ID:                |        |        |                     |        |        |                     |       |       |
|-----------------------------|--------|--------|---------------------|--------|--------|---------------------|-------|-------|
| RRF001 = VV038973.D         |        |        | RRF005 = VV038974.D |        |        | RRF020 = VV038975.D |       |       |
| RRF050 = VV038976.D         |        |        | RRF100 = VV038977.D |        |        | RRF010 = VV038979.D |       |       |
| COMPOUND                    | RRF001 | RRF005 | RRF020              | RRF050 | RRF100 | RRF010              | RRF   | % RSD |
| t-1,3-Dichloropropene       | 0.539  | 0.509  | 0.567               | 0.696  | 0.602  | 0.591               | 0.584 | 11.1  |
| cis-1,3-Dichloropropene     | 0.502  | 0.572  | 0.634               | 0.628  | 0.635  | 0.660               | 0.605 | 9.6   |
| 1,1,2-Trichloroethane       | 0.404  | 0.432  | 0.417               | 0.466  | 0.381  | 0.427               | 0.421 | 6.8   |
| 2-Hexanone                  | 0.208  | 0.302  | 0.385               | 0.465  | 0.384  | 0.388               | 0.355 | 24.9  |
| Dibromochloromethane        | 0.498  | 0.522  | 0.495               | 0.559  | 0.462  | 0.516               | 0.509 | 6.4   |
| 1,2-Dibromoethane           | 0.367  | 0.437  | 0.442               | 0.498  | 0.417  | 0.459               | 0.437 | 10    |
| Tetrachloroethene           | 0.359  | 0.382  | 0.368               | 0.364  | 0.362  | 0.397               | 0.372 | 4     |
| Chlorobenzene               | 1.265  | 1.206  | 1.166               | 1.163  | 1.142  | 1.291               | 1.206 | 5     |
| Ethyl Benzene               | 1.654  | 1.659  | 1.848               | 1.952  | 1.998  | 1.931               | 1.840 | 8.2   |
| m/p-Xylenes                 | 0.574  | 0.615  | 0.762               | 0.784  | 0.785  | 0.769               | 0.715 | 13.2  |
| o-Xylene                    | 0.529  | 0.572  | 0.669               | 0.712  | 0.740  | 0.686               | 0.651 | 12.7  |
| Styrene                     | 0.853  | 0.911  | 1.210               | 1.291  | 1.298  | 1.143               | 1.118 | 17.2  |
| Bromoform                   | 0.369  | 0.360  | 0.351               | 0.350  | 0.344  | 0.365               | 0.357 | 2.7   |
| Isopropylbenzene            | 3.332  | 2.780  | 3.605               | 3.326  | 3.515  | 3.573               | 3.355 | 9.1   |
| 1,1,2,2-Tetrachloroethane   | 1.745  | 1.323  | 1.324               | 1.111  | 1.122  | 1.368               | 1.332 | 17.3  |
| 1,3-Dichlorobenzene         | 1.735  | 1.691  | 1.660               | 1.623  | 1.663  | 1.841               | 1.702 | 4.6   |
| 1,4-Dichlorobenzene         | 2.000  | 1.803  | 1.733               | 1.666  | 1.683  | 2.011               | 1.816 | 8.5   |
| 1,2-Dichlorobenzene         | 1.685  | 1.559  | 1.707               | 1.521  | 1.564  | 1.702               | 1.623 | 5.2   |
| 1,2-Dibromo-3-Chloropropane | 0.287  | 0.261  | 0.274               | 0.236  | 0.255  | 0.242               | 0.259 | 7.4   |
| 1,2,4-Trichlorobenzene      | 0.938  | 0.865  | 1.004               | 0.973  | 1.065  | 0.965               | 0.969 | 6.9   |
| 1,2,3-Trichlorobenzene      | 0.878  | 0.847  | 1.059               | 0.983  | 1.064  | 1.028               | 0.977 | 9.6   |
| 1,2-Dichloroethane-d4       |        | 0.815  | 0.696               | 0.643  | 0.647  | 0.707               | 0.702 | 9.9   |
| Dibromofluoromethane        |        | 0.420  | 0.386               | 0.346  | 0.347  | 0.377               | 0.375 | 8.2   |
| Toluene-d8                  |        | 1.355  | 1.346               | 1.338  | 1.246  | 1.253               | 1.308 | 4.1   |
| 4-Bromofluorobenzene        |        | 0.425  | 0.472               | 0.559  | 0.477  | 0.459               | 0.478 | 10.3  |

\* Compounds with required minimum RRF and maximum %RSD values.  
All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.

VOLATILE CONTINUING CALIBRATION CHECK

|                     |                   |                        |                              |
|---------------------|-------------------|------------------------|------------------------------|
| Lab Name:           | <u>Alliance</u>   | Contract:              | <u>AECO02</u>                |
| Lab Code:           | <u>ACE</u>        | SDG No.:               | <u>Q2924</u>                 |
| Instrument ID:      | <u>MSVOA_V</u>    | Calibration Date/Time: | <u>08/21/2025 15:42</u>      |
| Lab File ID:        | <u>VV038982.D</u> | Init. Calib. Date(s):  | <u>08/21/2025 08/21/2025</u> |
| Heated Purge: (Y/N) | <u>N</u>          | Init. Calib. Time(s):  | <u>09:05 11:45</u>           |
| GC Column:          | <u>DB-624UI</u>   | ID:                    | <u>0.18</u> (mm)             |

| COMPOUND                       | RRF   | RRF050 | MIN RRF | %D     | MAX%D |
|--------------------------------|-------|--------|---------|--------|-------|
| Dichlorodifluoromethane        | 0.747 | 0.644  |         | -13.79 | 20    |
| Chloromethane                  | 0.772 | 0.671  | 0.1     | -13.08 | 20    |
| Vinyl Chloride                 | 0.815 | 0.716  |         | -12.15 | 20    |
| Bromomethane                   | 0.440 | 0.434  |         | -1.36  | 20    |
| Chloroethane                   | 0.415 | 0.391  |         | -5.78  | 20    |
| Trichlorofluoromethane         | 1.116 | 1.035  |         | -7.26  | 20    |
| 1,1,2-Trichlorotrifluoroethane | 0.616 | 0.554  |         | -10.06 | 20    |
| 1,1-Dichloroethene             | 0.567 | 0.536  |         | -5.47  | 20    |
| Acetone                        | 0.282 | 0.277  |         | -1.77  | 20    |
| Carbon Disulfide               | 1.578 | 1.548  |         | -1.9   | 20    |
| Methyl tert-butyl Ether        | 1.678 | 1.921  |         | 14.48  | 20    |
| Methyl Acetate                 | 0.597 | 0.750  |         | 25.63  | 20    |
| Methylene Chloride             | 0.654 | 0.690  |         | 5.51   | 20    |
| trans-1,2-Dichloroethene       | 0.591 | 0.641  |         | 8.46   | 20    |
| 1,1-Dichloroethane             | 1.102 | 1.114  | 0.1     | 1.09   | 20    |
| Cyclohexane                    | 0.983 | 0.959  |         | -2.44  | 20    |
| 2-Butanone                     | 0.366 | 0.416  |         | 13.66  | 20    |
| Carbon Tetrachloride           | 0.652 | 0.600  |         | -7.97  | 20    |
| cis-1,2-Dichloroethene         | 0.713 | 0.718  |         | 0.7    | 20    |
| Bromochloromethane             | 0.320 | 0.269  |         | -15.94 | 20    |
| Chloroform                     | 1.259 | 1.201  |         | -4.61  | 20    |
| 1,1,1-Trichloroethane          | 1.183 | 1.077  |         | -8.96  | 20    |
| Methylcyclohexane              | 0.622 | 0.666  |         | 7.07   | 20    |
| Benzene                        | 1.525 | 1.571  |         | 3.02   | 20    |
| 1,2-Dichloroethane             | 0.558 | 0.537  |         | -3.76  | 20    |
| Trichloroethene                | 0.414 | 0.401  |         | -3.14  | 20    |
| 1,2-Dichloropropane            | 0.366 | 0.389  |         | 6.28   | 20    |
| Bromodichloromethane           | 0.625 | 0.597  |         | -4.48  | 20    |
| 4-Methyl-2-Pentanone           | 0.477 | 0.544  |         | 14.05  | 20    |
| Toluene                        | 0.966 | 1.015  |         | 5.07   | 20    |
| t-1,3-Dichloropropene          | 0.584 | 0.604  |         | 3.42   | 20    |
| cis-1,3-Dichloropropene        | 0.605 | 0.643  |         | 6.28   | 20    |
| 1,1,2-Trichloroethane          | 0.421 | 0.414  |         | -1.66  | 20    |
| 2-Hexanone                     | 0.355 | 0.405  |         | 14.09  | 20    |
| Dibromochloromethane           | 0.509 | 0.477  |         | -6.29  | 20    |
| 1,2-Dibromoethane              | 0.437 | 0.436  |         | -0.23  | 20    |
| Tetrachloroethene              | 0.372 | 0.377  |         | 1.34   | 20    |
| Chlorobenzene                  | 1.206 | 1.194  | 0.3     | -1     | 20    |

All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.

VOLATILE CONTINUING CALIBRATION CHECK

|                     |                   |                        |                              |
|---------------------|-------------------|------------------------|------------------------------|
| Lab Name:           | <u>Alliance</u>   | Contract:              | <u>AECO02</u>                |
| Lab Code:           | <u>ACE</u>        | SDG No.:               | <u>Q2924</u>                 |
| Instrument ID:      | <u>MSVOA_V</u>    | Calibration Date/Time: | <u>08/21/2025 15:42</u>      |
| Lab File ID:        | <u>VV038982.D</u> | Init. Calib. Date(s):  | <u>08/21/2025 08/21/2025</u> |
| Heated Purge: (Y/N) | <u>N</u>          | Init. Calib. Time(s):  | <u>09:05 11:45</u>           |
| GC Column:          | <u>DB-624UI</u>   | ID:                    | <u>0.18</u> (mm)             |

| COMPOUND                    | RRF   | RRF050 | MIN<br>RRF | %D    | MAX%D |
|-----------------------------|-------|--------|------------|-------|-------|
| Ethyl Benzene               | 1.840 | 2.027  |            | 10.16 | 20    |
| m/p-Xylenes                 | 0.715 | 0.813  |            | 13.71 | 20    |
| o-Xylene                    | 0.651 | 0.749  |            | 15.05 | 20    |
| Styrene                     | 1.118 | 1.344  |            | 20.22 | 20    |
| Bromoform                   | 0.357 | 0.356  | 0.1        | -0.28 | 20    |
| Isopropylbenzene            | 3.355 | 3.588  |            | 6.95  | 20    |
| 1,1,2,2-Tetrachloroethane   | 1.332 | 1.201  | 0.3        | -9.84 | 20    |
| 1,3-Dichlorobenzene         | 1.702 | 1.726  |            | 1.41  | 20    |
| 1,4-Dichlorobenzene         | 1.816 | 1.752  |            | -3.52 | 20    |
| 1,2-Dichlorobenzene         | 1.623 | 1.634  |            | 0.68  | 20    |
| 1,2-Dibromo-3-Chloropropane | 0.259 | 0.256  |            | -1.16 | 20    |
| 1,2,4-Trichlorobenzene      | 0.969 | 1.048  |            | 8.15  | 20    |
| 1,2,3-Trichlorobenzene      | 0.977 | 1.075  |            | 10.03 | 20    |
| 1,2-Dichloroethane-d4       | 0.702 | 0.672  |            | -4.27 | 20    |
| Dibromofluoromethane        | 0.375 | 0.368  |            | -1.87 | 20    |
| Toluene-d8                  | 1.308 | 1.282  |            | -1.99 | 20    |
| 4-Bromofluorobenzene        | 0.478 | 0.486  |            | 1.67  | 20    |

All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.

VOLATILE CONTINUING CALIBRATION CHECK

|                     |                   |                        |                                     |
|---------------------|-------------------|------------------------|-------------------------------------|
| Lab Name:           | <u>Alliance</u>   | Contract:              | <u>AECO02</u>                       |
| Lab Code:           | <u>ACE</u>        | SDG No.:               | <u>Q2924</u>                        |
| Instrument ID:      | <u>MSVOA_V</u>    | Calibration Date/Time: | <u>08/22/2025</u> <u>09:03</u>      |
| Lab File ID:        | <u>VV039009.D</u> | Init. Calib. Date(s):  | <u>08/21/2025</u> <u>08/21/2025</u> |
| Heated Purge: (Y/N) | <u>N</u>          | Init. Calib. Time(s):  | <u>09:05</u> <u>11:45</u>           |
| GC Column:          | <u>DB-624UI</u>   | ID:                    | <u>0.18</u> (mm)                    |

| COMPOUND                       | RRF   | RRF050 | MIN<br>RRF | %D     | MAX%D |
|--------------------------------|-------|--------|------------|--------|-------|
| Dichlorodifluoromethane        | 0.747 | 0.601  |            | -19.55 | 20    |
| Chloromethane                  | 0.772 | 0.585  | 0.1        | -24.22 | 20    |
| Vinyl Chloride                 | 0.815 | 0.691  |            | -15.22 | 20    |
| Bromomethane                   | 0.440 | 0.347  |            | -21.14 | 20    |
| Chloroethane                   | 0.415 | 0.469  |            | 13.01  | 20    |
| Trichlorofluoromethane         | 1.116 | 1.051  |            | -5.82  | 20    |
| 1,1,2-Trichlorotrifluoroethane | 0.616 | 0.640  |            | 3.9    | 20    |
| 1,1-Dichloroethene             | 0.567 | 0.607  |            | 7.05   | 20    |
| Acetone                        | 0.282 | 0.366  |            | 29.79  | 20    |
| Carbon Disulfide               | 1.578 | 1.789  |            | 13.31  | 20    |
| Methyl tert-butyl Ether        | 1.678 | 2.087  |            | 24.37  | 20    |
| Methyl Acetate                 | 0.597 | 0.801  |            | 34.17  | 20    |
| Methylene Chloride             | 0.654 | 0.668  |            | 2.14   | 20    |
| trans-1,2-Dichloroethene       | 0.591 | 0.630  |            | 6.6    | 20    |
| 1,1-Dichloroethane             | 1.102 | 1.141  | 0.1        | 3.54   | 20    |
| Cyclohexane                    | 0.983 | 1.041  |            | 5.9    | 20    |
| 2-Butanone                     | 0.366 | 0.460  |            | 25.68  | 20    |
| Carbon Tetrachloride           | 0.652 | 0.526  |            | -19.33 | 20    |
| cis-1,2-Dichloroethene         | 0.713 | 0.751  |            | 5.33   | 20    |
| Bromochloromethane             | 0.320 | 0.277  |            | -13.44 | 20    |
| Chloroform                     | 1.259 | 1.179  |            | -6.35  | 20    |
| 1,1,1-Trichloroethane          | 1.183 | 1.068  |            | -9.72  | 20    |
| Methylcyclohexane              | 0.622 | 0.651  |            | 4.66   | 20    |
| Benzene                        | 1.525 | 1.478  |            | -3.08  | 20    |
| 1,2-Dichloroethane             | 0.558 | 0.490  |            | -12.19 | 20    |
| Trichloroethene                | 0.414 | 0.376  |            | -9.18  | 20    |
| 1,2-Dichloropropane            | 0.366 | 0.371  |            | 1.37   | 20    |
| Bromodichloromethane           | 0.625 | 0.561  |            | -10.24 | 20    |
| 4-Methyl-2-Pentanone           | 0.477 | 0.534  |            | 11.95  | 20    |
| Toluene                        | 0.966 | 0.942  |            | -2.48  | 20    |
| t-1,3-Dichloropropene          | 0.584 | 0.588  |            | 0.69   | 20    |
| cis-1,3-Dichloropropene        | 0.605 | 0.623  |            | 2.97   | 20    |
| 1,1,2-Trichloroethane          | 0.421 | 0.375  |            | -10.93 | 20    |
| 2-Hexanone                     | 0.355 | 0.396  |            | 11.55  | 20    |
| Dibromochloromethane           | 0.509 | 0.441  |            | -13.36 | 20    |
| 1,2-Dibromoethane              | 0.437 | 0.409  |            | -6.41  | 20    |
| Tetrachloroethene              | 0.372 | 0.348  |            | -6.45  | 20    |
| Chlorobenzene                  | 1.206 | 1.155  | 0.3        | -4.23  | 20    |

All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.

VOLATILE CONTINUING CALIBRATION CHECK

|                     |                   |                        |                                     |
|---------------------|-------------------|------------------------|-------------------------------------|
| Lab Name:           | <u>Alliance</u>   | Contract:              | <u>AECO02</u>                       |
| Lab Code:           | <u>ACE</u>        | SDG No.:               | <u>Q2924</u>                        |
| Instrument ID:      | <u>MSVOA_V</u>    | Calibration Date/Time: | <u>08/22/2025</u> <u>09:03</u>      |
| Lab File ID:        | <u>VV039009.D</u> | Init. Calib. Date(s):  | <u>08/21/2025</u> <u>08/21/2025</u> |
| Heated Purge: (Y/N) | <u>N</u>          | Init. Calib. Time(s):  | <u>09:05</u> <u>11:45</u>           |
| GC Column:          | <u>DB-624UI</u>   | ID:                    | <u>0.18</u> (mm)                    |

| COMPOUND                    | RRF   | RRF050 | MIN<br>RRF | %D    | MAX%D |
|-----------------------------|-------|--------|------------|-------|-------|
| Ethyl Benzene               | 1.840 | 2.002  |            | 8.8   | 20    |
| m/p-Xylenes                 | 0.715 | 0.770  |            | 7.69  | 20    |
| o-Xylene                    | 0.651 | 0.736  |            | 13.06 | 20    |
| Styrene                     | 1.118 | 1.245  |            | 11.36 | 20    |
| Bromoform                   | 0.357 | 0.338  | 0.1        | -5.32 | 20    |
| Isopropylbenzene            | 3.355 | 3.778  |            | 12.61 | 20    |
| 1,1,2,2-Tetrachloroethane   | 1.332 | 1.280  | 0.3        | -3.9  | 20    |
| 1,3-Dichlorobenzene         | 1.702 | 1.744  |            | 2.47  | 20    |
| 1,4-Dichlorobenzene         | 1.816 | 1.769  |            | -2.59 | 20    |
| 1,2-Dichlorobenzene         | 1.623 | 1.615  |            | -0.49 | 20    |
| 1,2-Dibromo-3-Chloropropane | 0.259 | 0.274  |            | 5.79  | 20    |
| 1,2,4-Trichlorobenzene      | 0.969 | 1.043  |            | 7.64  | 20    |
| 1,2,3-Trichlorobenzene      | 0.977 | 1.083  |            | 10.85 | 20    |
| 1,2-Dichloroethane-d4       | 0.702 | 0.659  |            | -6.13 | 20    |
| Dibromofluoromethane        | 0.375 | 0.333  |            | -11.2 | 20    |
| Toluene-d8                  | 1.308 | 1.155  |            | -11.7 | 20    |
| 4-Bromofluorobenzene        | 0.478 | 0.443  |            | -7.32 | 20    |

All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.

## LAB CHRONICLE

|                 |         |                   |                                    |
|-----------------|---------|-------------------|------------------------------------|
| <b>OrderID:</b> | Q2924   | <b>OrderDate:</b> | 8/20/2025 3:43:00 PM               |
| <b>Client:</b>  | AECOM   | <b>Project:</b>   | National Grid Equity - Brooklyn NY |
| <b>Contact:</b> | Peter S | <b>Location:</b>  | J23,VOA Lab                        |

| LabID                   | ClientID                | Matrix       | Test             | Method | Sample Date     | Prep Date | Anal Date | Received        |
|-------------------------|-------------------------|--------------|------------------|--------|-----------------|-----------|-----------|-----------------|
| <b>Q2924-01</b>         | <b>MW-10C-082025</b>    | <b>Water</b> | SVOC-TCL BNA -20 | 8270E  | <b>08/20/25</b> | 08/22/25  | 08/22/25  | <b>08/20/25</b> |
| <b>Q2924-02</b>         | <b>MW-201-082025</b>    | <b>Water</b> | SVOC-TCL BNA -20 | 8270E  | <b>08/20/25</b> | 08/22/25  | 08/27/25  | <b>08/20/25</b> |
| <b>Q2924-02DL</b>       | <b>MW-201-082025DL</b>  | <b>Water</b> | SVOC-TCL BNA -20 | 8270E  | <b>08/20/25</b> | 08/22/25  | 08/27/25  | <b>08/20/25</b> |
| <b>Q2924-02DL<br/>2</b> | <b>MW-201-082025DL2</b> | <b>Water</b> | SVOC-TCL BNA -20 | 8270E  | <b>08/20/25</b> | 08/22/25  | 08/27/25  | <b>08/20/25</b> |
| <b>Q2924-05</b>         | <b>MW-2B-082025</b>     | <b>Water</b> | SVOC-TCL BNA -20 | 8270E  | <b>08/20/25</b> | 08/22/25  | 08/22/25  | <b>08/20/25</b> |
| <b>Q2924-05DL</b>       | <b>MW-2B-082025DL</b>   | <b>Water</b> | SVOC-TCL BNA -20 | 8270E  | <b>08/20/25</b> | 08/22/25  | 08/26/25  | <b>08/20/25</b> |
| <b>Q2924-05DL<br/>2</b> | <b>MW-2B-082025DL2</b>  | <b>Water</b> | SVOC-TCL BNA -20 | 8270E  | <b>08/20/25</b> | 08/22/25  | 08/26/25  | <b>08/20/25</b> |
| <b>Q2924-06</b>         | <b>MW-2A-082025</b>     | <b>Water</b> | SVOC-TCL BNA -20 | 8270E  | <b>08/20/25</b> | 08/22/25  | 08/26/25  | <b>08/20/25</b> |
| <b>Q2924-07</b>         | <b>MW-18C-082025</b>    | <b>Water</b> | SVOC-TCL BNA -20 | 8270E  | <b>08/20/25</b> | 08/22/25  | 08/27/25  | <b>08/20/25</b> |
| <b>Q2924-08</b>         | <b>MW-16A-082025</b>    | <b>Water</b> | SVOC-TCL BNA -20 | 8270E  | <b>08/20/25</b> | 08/22/25  | 08/26/25  | <b>08/20/25</b> |
| <b>Q2924-09</b>         | <b>MW-16C-082025</b>    | <b>Water</b> | SVOC-TCL BNA -20 | 8270E  | <b>08/20/25</b> | 08/22/25  | 08/26/25  | <b>08/20/25</b> |

## LAB CHRONICLE

|                 |                     |              |                  |       |                 |          |          |                 |
|-----------------|---------------------|--------------|------------------|-------|-----------------|----------|----------|-----------------|
| <b>Q2924-10</b> | <b>MW-8A-082025</b> | <b>Water</b> |                  |       | <b>08/20/25</b> |          |          | <b>08/20/25</b> |
|                 |                     |              | SVOC-TCL BNA -20 | 8270E |                 | 08/22/25 | 08/22/25 |                 |
| <b>Q2924-11</b> | <b>FB-01-082025</b> | <b>Water</b> |                  |       | <b>08/20/25</b> |          |          | <b>08/20/25</b> |
|                 |                     |              | SVOC-TCL BNA -20 | 8270E |                 | 08/22/25 | 08/27/25 |                 |

### Hit Summary Sheet SW-846

SDG No.: Q2924  
Client: AECOM

| Sample ID   | Client ID     |       | Parameter                           | Concentration | C  | MDL      | RDL  | Units |
|-------------|---------------|-------|-------------------------------------|---------------|----|----------|------|-------|
| Client ID : | MW-10C-082025 |       |                                     |               |    |          |      |       |
| Q2924-01    | MW-10C-082025 | WATER | 1,4-Benzenedicarboxylic acid, bis * | 2.500         | J  | 0        | 0    | ug/L  |
| Q2924-01    | MW-10C-082025 | WATER | 2-Pentanone, 4-hydroxy-4-methyl *   | 5.500         | AB | 0        | 0    | ug/L  |
| Q2924-01    | MW-10C-082025 | WATER | Butane, 2-methoxy-2-methyl- *       | 87.300        | J  | 0        | 0    | ug/L  |
| Q2924-01    | MW-10C-082025 | WATER | Hexanoic acid, 2-ethyl- *           | 5.300         | J  | 0        | 0    | ug/L  |
|             |               |       | Total Tics :                        |               |    | 100.60   |      |       |
|             |               |       | Total Concentration:                |               |    | 100.60   |      |       |
| Client ID : | MW-201-082025 |       |                                     |               |    |          |      |       |
| Q2924-02    | MW-201-082025 | WATER | 3+4-Methylphenols                   | 10.000        | J  | 1.2      | 10.5 | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Naphthalene                         | 3,100.000     | E  | 0.53     | 5.3  | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | 2-Methylnaphthalene                 | 920.000       | E  | 0.59     | 5.3  | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | 1,1-Biphenyl                        | 26.000        |    | 0.56     | 5.3  | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Acenaphthylene                      | 160.000       | E  | 0.79     | 5.3  | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Acenaphthene                        | 8.700         |    | 0.58     | 5.3  | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Dibenzofuran                        | 4.000         | J  | 0.64     | 5.3  | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Diethylphthalate                    | 2.100         | J  | 0.73     | 5.3  | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Fluorene                            | 27.400        |    | 0.66     | 5.3  | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Phenanthrene                        | 29.600        |    | 0.53     | 5.3  | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Anthracene                          | 6.100         |    | 0.64     | 5.3  | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Carbazole                           | 13.800        |    | 0.76     | 5.3  | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Di-n-butylphthalate                 | 2.700         | J  | 1.3      | 5.3  | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Fluoranthene                        | 2.300         | J  | 0.86     | 5.3  | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Pyrene                              | 3.000         | J  | 0.53     | 5.3  | ug/L  |
|             |               |       | Total Svoc :                        |               |    | 4,315.70 |      |       |
| Q2924-02    | MW-201-082025 | WATER | .alpha.-Methylstyrene *             | 11.100        | J  | 0        | 0    | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | 4-Methylcarbazole *                 | 16.200        | J  | 0        | 0    | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Azulene *                           | 21.100        | J  | 0        | 0    | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Benzene, (2-methyl-1-propenyl)- *   | 7.900         | J  | 0        | 0    | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Benzene, 1-ethyl-3-methyl- *        | 12.800        | J  | 0        | 0    | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Benzene, 1-ethyl-4-methyl- *        | 5.300         | J  | 0        | 0    | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Benzene, 1-propenyl- *              | 9.500         | J  | 0        | 0    | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Butane, 2-methoxy-2-methyl- *       | 11.500        | J  | 0        | 0    | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Indane *                            | 12.700        | J  | 0        | 0    | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Naphthalene, 1,3-dimethyl- *        | 5.900         | J  | 0        | 0    | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Naphthalene, 1,7-dimethyl- *        | 4.600         | J  | 0        | 0    | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | p-Xylene *                          | 160.000       | J  | 0        | 0    | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | Thiophene, 3-methyl- *              | 4.600         | J  | 0        | 0    | ug/L  |
| Q2924-02    | MW-201-082025 | WATER | unknown7.449 *                      | 99.800        | J  | 0        | 0    | ug/L  |

### Hit Summary Sheet SW-846

**SDG No.:** Q2924  
**Client:** AECOM

| Sample ID                    | Client ID        |       | Parameter                         | Concentration | C       | MDL    | RDL  | Units |
|------------------------------|------------------|-------|-----------------------------------|---------------|---------|--------|------|-------|
| Q2924-02                     | MW-201-082025    | WATER | unknown7.525                      | *             | 31.200  | J 0    | 0    | ug/L  |
| Q2924-02                     | MW-201-082025    | WATER | unknown8.372                      | *             | 310.000 | J 0    | 0    | ug/L  |
| Q2924-02                     | MW-201-082025    | WATER | 1-Methylnaphthalene               | *             | 600.000 | J 0.69 | 5.3  | ug/L  |
| Total Tics :                 |                  |       |                                   | 1,324.20      |         |        |      |       |
| Total Concentration:         |                  |       |                                   | 5,639.90      |         |        |      |       |
| Client ID : MW-201-082025DL  |                  |       |                                   |               |         |        |      |       |
| Q2924-02DL                   | MW-201-082025DL  | WATER | Naphthalene                       | 6,000.000     | ED      | 10.5   | 110  | ug/L  |
| Q2924-02DL                   | MW-201-082025DL  | WATER | 2-Methylnaphthalene               | 610.000       | D       | 11.8   | 110  | ug/L  |
| Q2924-02DL                   | MW-201-082025DL  | WATER | Acenaphthylene                    | 180.000       | D       | 15.8   | 110  | ug/L  |
| Total Svoc :                 |                  |       |                                   | 6,790.00      |         |        |      |       |
| Total Concentration:         |                  |       |                                   | 6,790.00      |         |        |      |       |
| Client ID : MW-201-082025DL2 |                  |       |                                   |               |         |        |      |       |
| Q2924-02DL2                  | MW-201-082025DL2 | WATER | Naphthalene                       | 6,800.000     | D       | 110    | 1100 | ug/L  |
| Q2924-02DL2                  | MW-201-082025DL2 | WATER | 2-Methylnaphthalene               | 570.000       | JD      | 120    | 1100 | ug/L  |
| Total Svoc :                 |                  |       |                                   | 7,370.00      |         |        |      |       |
| Total Concentration:         |                  |       |                                   | 7,370.00      |         |        |      |       |
| Client ID : MW-2B-082025     |                  |       |                                   |               |         |        |      |       |
| Q2924-05                     | MW-2B-082025     | WATER | Phenol                            | 5.100         | J       | 0.95   | 5.2  | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | Naphthalene                       | 930.000       | E       | 0.52   | 5.2  | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | 2-Methylnaphthalene               | 160.000       | E       | 0.58   | 5.2  | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | 1,1-Biphenyl                      | 12.700        |         | 0.55   | 5.2  | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | Acenaphthene                      | 61.200        |         | 0.57   | 5.2  | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | Dibenzofuran                      | 2.600         | J       | 0.64   | 5.2  | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | Fluorene                          | 13.000        |         | 0.66   | 5.2  | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | Phenanthrene                      | 17.200        |         | 0.52   | 5.2  | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | Anthracene                        | 2.600         | J       | 0.64   | 5.2  | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | Carbazole                         | 5.800         |         | 0.75   | 5.2  | ug/L  |
| Total Svoc :                 |                  |       |                                   | 1,210.20      |         |        |      |       |
| Q2924-05                     | MW-2B-082025     | WATER | Naphthalene, 1,5-dimethyl-        | *             | 8.900   | J 0    | 0    | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | Naphthalene, 1,6-dimethyl-        | *             | 15.600  | J 0    | 0    | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | Naphthalene, 1-ethyl-             | *             | 8.400   | J 0    | 0    | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | Naphthalene, 2,3-dimethyl-        | *             | 18.700  | J 0    | 0    | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | Naphthalene, 2,6-dimethyl-        | *             | 5.900   | J 0    | 0    | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | 1,3-Cyclopentadiene, 5-(1-methyl- | *             | 290.000 | J 0    | 0    | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | 1-Naphthalenol                    | *             | 7.100   | J 0    | 0    | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | 1-Naphthalenol, 2-methyl-         | *             | 8.800   | J 0    | 0    | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | 9H-Carbazole, 9-methyl-           | *             | 7.700   | J 0    | 0    | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | Benzene, 1,3-dimethyl-            | *             | 28.600  | J 0    | 0    | ug/L  |
| Q2924-05                     | MW-2B-082025     | WATER | Butane, 2-methoxy-2-methyl-       | *             | 110.000 | J 0    | 0    | ug/L  |

### Hit Summary Sheet SW-846

SDG No.: Q2924  
Client: AECOM

| Sample ID                   | Client ID       |       | Parameter                         | Concentration | C         | MDL |      | RDL | Units |
|-----------------------------|-----------------|-------|-----------------------------------|---------------|-----------|-----|------|-----|-------|
| Q2924-05                    | MW-2B-082025    | WATER | unknown6.387                      | *             | 14.000    | J   | 0    | 0   | ug/L  |
| Q2924-05                    | MW-2B-082025    | WATER | unknown6.751                      | *             | 40.500    | J   | 0    | 0   | ug/L  |
| Q2924-05                    | MW-2B-082025    | WATER | unknown6.987                      | *             | 76.500    | J   | 0    | 0   | ug/L  |
| Q2924-05                    | MW-2B-082025    | WATER | unknown7.034                      | *             | 14.100    | J   | 0    | 0   | ug/L  |
| Q2924-05                    | MW-2B-082025    | WATER | unknown7.134                      | *             | 770.000   | J   | 0    | 0   | ug/L  |
| Q2924-05                    | MW-2B-082025    | WATER | unknown7.193                      | *             | 35.400    | J   | 0    | 0   | ug/L  |
| Q2924-05                    | MW-2B-082025    | WATER | unknown7.951                      | *             | 2.300     | J   | 0    | 0   | ug/L  |
| Q2924-05                    | MW-2B-082025    | WATER | 1-Methylnaphthalene               | *             | 240.000   | J   | 0.69 | 5.2 | ug/L  |
| Total Tics :                |                 |       |                                   | 1,702.50      |           |     |      |     |       |
| Total Concentration:        |                 |       |                                   | 2,912.70      |           |     |      |     |       |
| Client ID : MW-2B-082025DL  |                 |       |                                   |               |           |     |      |     |       |
| Q2924-05DL                  | MW-2B-082025DL  | WATER | Naphthalene                       |               | 1,800.000 | ED  | 10.4 | 100 | ug/L  |
| Q2924-05DL                  | MW-2B-082025DL  | WATER | 2-Methylnaphthalene               |               | 180.000   | D   | 11.7 | 100 | ug/L  |
| Q2924-05DL                  | MW-2B-082025DL  | WATER | Acenaphthene                      |               | 72.500    | JD  | 11.5 | 100 | ug/L  |
| Total Svoc :                |                 |       |                                   | 2,052.50      |           |     |      |     |       |
| Total Concentration:        |                 |       |                                   | 2,052.50      |           |     |      |     |       |
| Client ID : MW-2B-082025DL2 |                 |       |                                   |               |           |     |      |     |       |
| Q2924-05DL2                 | MW-2B-082025DL2 | WATER | Naphthalene                       |               | 1,800.000 | D   | 20.8 | 210 | ug/L  |
| Q2924-05DL2                 | MW-2B-082025DL2 | WATER | 2-Methylnaphthalene               |               | 180.000   | JD  | 23.3 | 210 | ug/L  |
| Total Svoc :                |                 |       |                                   | 1,980.00      |           |     |      |     |       |
| Total Concentration:        |                 |       |                                   | 1,980.00      |           |     |      |     |       |
| Client ID : MW-2A-082025    |                 |       |                                   |               |           |     |      |     |       |
| Q2924-06                    | MW-2A-082025    | WATER | 1-Octanamine, N-methyl-N-octyl-   | *             | 2.900     | J   | 0    | 0   | ug/L  |
| Q2924-06                    | MW-2A-082025    | WATER | 1-Propanol, 2-(2-hydroxypropoxy-  | *             | 3.700     | J   | 0    | 0   | ug/L  |
| Q2924-06                    | MW-2A-082025    | WATER | 2-Pentanone, 4-hydroxy-4-methyl   | *             | 4.900     | AB  | 0    | 0   | ug/L  |
| Q2924-06                    | MW-2A-082025    | WATER | 2-Propanol, 1,1-[(1-methyl-1,2-et | *             | 4.200     | J   | 0    | 0   | ug/L  |
| Q2924-06                    | MW-2A-082025    | WATER | Butane, 2-methoxy-2-methyl-       | *             | 96.500    | J   | 0    | 0   | ug/L  |
| Q2924-06                    | MW-2A-082025    | WATER | Butanoic acid, 4-(2-methoxy-1-m   | *             | 2.700     | J   | 0    | 0   | ug/L  |
| Q2924-06                    | MW-2A-082025    | WATER | di-2-Pyridyl ketone               | *             | 5.000     | J   | 0    | 0   | ug/L  |
| Q2924-06                    | MW-2A-082025    | WATER | Ethyl cyclopropanecarboxylate     | *             | 17.200    | J   | 0    | 0   | ug/L  |
| Q2924-06                    | MW-2A-082025    | WATER | n-Hexadecanoic acid               | *             | 4.300     | J   | 0    | 0   | ug/L  |
| Q2924-06                    | MW-2A-082025    | WATER | unknown12.196                     | *             | 17.600    | J   | 0    | 0   | ug/L  |
| Q2924-06                    | MW-2A-082025    | WATER | unknown19.683                     | *             | 4.800     | J   | 0    | 0   | ug/L  |
| Total Tics :                |                 |       |                                   | 163.80        |           |     |      |     |       |
| Total Concentration:        |                 |       |                                   | 163.80        |           |     |      |     |       |
| Client ID : MW-18C-082025   |                 |       |                                   |               |           |     |      |     |       |
| Q2924-07                    | MW-18C-082025   | WATER | 2-Pentanone, 4-hydroxy-4-methyl   | *             | 5.800     | AB  | 0    | 0   | ug/L  |
| Q2924-07                    | MW-18C-082025   | WATER | Benzophenone                      | *             | 2.300     | J   | 0    | 0   | ug/L  |
| Q2924-07                    | MW-18C-082025   | WATER | n-Hexadecanoic acid               | *             | 2.700     | J   | 0    | 0   | ug/L  |

### Hit Summary Sheet SW-846

SDG No.: Q2924  
Client: AECOM

| Sample ID                 | Client ID     |       | Parameter                           | Concentration | C       | MDL    | RDL | Units |
|---------------------------|---------------|-------|-------------------------------------|---------------|---------|--------|-----|-------|
| Q2924-07                  | MW-18C-082025 | WATER | Palmitoleic acid                    | *             | 2.300   | J 0    | 0   | ug/L  |
| Total Tics :              |               |       |                                     |               |         | 13.10  |     |       |
| Total Concentration:      |               |       |                                     |               |         | 13.10  |     |       |
| Client ID : MW-16A-082025 |               |       |                                     |               |         |        |     |       |
| Q2924-08                  | MW-16A-082025 | WATER | 2-Pentanone, 4-hydroxy-4-methyl *   | 5.600         | AB      | 0      | 0   | ug/L  |
| Q2924-08                  | MW-16A-082025 | WATER | Butane, 2-methoxy-2-methyl-         | *             | 110.000 | J 0    | 0   | ug/L  |
| Q2924-08                  | MW-16A-082025 | WATER | cis-13-Octadecenoic acid            | *             | 7.300   | J 0    | 0   | ug/L  |
| Q2924-08                  | MW-16A-082025 | WATER | n-Hexadecanoic acid                 | *             | 5.200   | J 0    | 0   | ug/L  |
| Total Tics :              |               |       |                                     |               |         | 128.10 |     |       |
| Total Concentration:      |               |       |                                     |               |         | 128.10 |     |       |
| Client ID : MW-16C-082025 |               |       |                                     |               |         |        |     |       |
| Q2924-09                  | MW-16C-082025 | WATER | 1H-Indene, 2,3-dihydro-4,7-dimet *  | 2.900         | J       | 0      | 0   | ug/L  |
| Q2924-09                  | MW-16C-082025 | WATER | 2,3,4,5,6,7-Hexahydro-1H-cyclop *   | 2.600         | J       | 0      | 0   | ug/L  |
| Q2924-09                  | MW-16C-082025 | WATER | 2-Pentanone, 4-hydroxy-4-methyl *   | 5.300         | AB      | 0      | 0   | ug/L  |
| Q2924-09                  | MW-16C-082025 | WATER | 3-(2,3-Dihydro-1-benzofuran-5-yl *  | 3.900         | J       | 0      | 0   | ug/L  |
| Q2924-09                  | MW-16C-082025 | WATER | Benzenemethanol, .alpha.,.alpha.- * | 2.900         | J       | 0      | 0   | ug/L  |
| Q2924-09                  | MW-16C-082025 | WATER | Butane, 2-methoxy-2-methyl-         | *             | 110.000 | J 0    | 0   | ug/L  |
| Q2924-09                  | MW-16C-082025 | WATER | Erucic acid                         | *             | 2.700   | J 0    | 0   | ug/L  |
| Q2924-09                  | MW-16C-082025 | WATER | n-Hexadecanoic acid                 | *             | 4.000   | J 0    | 0   | ug/L  |
| Total Tics :              |               |       |                                     |               |         | 134.30 |     |       |
| Total Concentration:      |               |       |                                     |               |         | 134.30 |     |       |
| Client ID : MW-8A-082025  |               |       |                                     |               |         |        |     |       |
| Q2924-10                  | MW-8A-082025  | WATER | 2-Pentanone, 4-hydroxy-4-methyl *   | 4.400         | AB      | 0      | 0   | ug/L  |
| Q2924-10                  | MW-8A-082025  | WATER | Benzophenone                        | *             | 2.800   | J 0    | 0   | ug/L  |
| Q2924-10                  | MW-8A-082025  | WATER | Butane, 2-methoxy-2-methyl-         | *             | 110.000 | J 0    | 0   | ug/L  |
| Q2924-10                  | MW-8A-082025  | WATER | n-Hexadecanoic acid                 | *             | 5.300   | J 0    | 0   | ug/L  |
| Total Tics :              |               |       |                                     |               |         | 122.50 |     |       |
| Total Concentration:      |               |       |                                     |               |         | 122.50 |     |       |
| Client ID : FB-01-082025  |               |       |                                     |               |         |        |     |       |
| Q2924-11                  | FB-01-082025  | WATER | 2-Pentanone, 4-hydroxy-4-methyl *   | 6.900         | AB      | 0      | 0   | ug/L  |
| Total Tics :              |               |       |                                     |               |         | 6.90   |     |       |
| Total Concentration:      |               |       |                                     |               |         | 6.90   |     |       |



# SAMPLE DATA

## Report of Analysis

|                    |                                    |                 |                  |
|--------------------|------------------------------------|-----------------|------------------|
| Client:            | AECOM                              | Date Collected: | 08/20/25         |
| Project:           | National Grid Equity - Brooklyn NY | Date Received:  | 08/20/25         |
| Client Sample ID:  | MW-10C-082025                      | SDG No.:        | Q2924            |
| Lab Sample ID:     | Q2924-01                           | Matrix:         | Water            |
| Analytical Method: | 8270E                              | % Solid:        | 0                |
| Sample Wt/Vol:     | 1000 Units: mL                     | Final Vol:      | 1000 uL          |
| Soil Aliquot Vol:  | uL                                 | Test:           | SVOC-TCL BNA -20 |
| Extraction Type :  | Decanted : N                       | Level :         | LOW              |
| Injection Volume : | GPC Factor : 1.0                   | GPC Cleanup :   | N PH :           |
| Prep Method :      |                                    |                 |                  |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BF143581.D        | 1         | 08/22/25 10:34 | 08/22/25 19:21 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 100-52-7       | Benzaldehyde                | 10.0  | U         | 3.90 | 10.0       | ug/L  |
| 108-95-2       | Phenol                      | 5.00  | U         | 0.91 | 5.00       | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 5.00  | U         | 0.81 | 5.00       | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 5.00  | U         | 0.58 | 5.00       | ug/L  |
| 95-48-7        | 2-Methylphenol              | 5.00  | U         | 1.10 | 5.00       | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 5.00  | U         | 1.30 | 5.00       | ug/L  |
| 98-86-2        | Acetophenone                | 5.00  | U         | 0.74 | 5.00       | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 10.0  | U         | 1.10 | 10.0       | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 2.50  | U         | 1.40 | 2.50       | ug/L  |
| 67-72-1        | Hexachloroethane            | 5.00  | U         | 0.65 | 5.00       | ug/L  |
| 98-95-3        | Nitrobenzene                | 5.00  | U         | 0.76 | 5.00       | ug/L  |
| 78-59-1        | Isophorone                  | 5.00  | U         | 0.75 | 5.00       | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 5.00  | U         | 1.80 | 5.00       | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 5.00  | U         | 1.90 | 5.00       | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 5.00  | U         | 0.68 | 5.00       | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 5.00  | U         | 0.52 | 5.00       | ug/L  |
| 91-20-3        | Naphthalene                 | 5.00  | U         | 0.50 | 5.00       | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 5.00  | U         | 0.84 | 5.00       | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 5.00  | U         | 0.54 | 5.00       | ug/L  |
| 105-60-2       | Caprolactam                 | 10.0  | U         | 1.10 | 10.0       | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 5.00  | U         | 0.59 | 5.00       | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 5.00  | U         | 0.56 | 5.00       | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 10.0  | U         | 3.60 | 10.0       | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 5.00  | U         | 0.51 | 5.00       | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 5.00  | U         | 0.62 | 5.00       | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 5.00  | U         | 0.53 | 5.00       | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 5.00  | U         | 0.61 | 5.00       | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 5.00  | U         | 1.30 | 5.00       | ug/L  |
| 131-11-3       | Dimethylphthalate           | 5.00  | U         | 0.61 | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-10C-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-01                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 1000                               | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BF143581.D        | 1         | 08/22/25 10:34 | 08/22/25 19:21 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 5.00  | U         | 0.75 | 5.00       | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 5.00  | U         | 0.92 | 5.00       | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 5.00  | U         | 1.10 | 5.00       | ug/L  |
| 83-32-9    | Acenaphthene               | 5.00  | U         | 0.55 | 5.00       | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 10.0  | U         | 6.00 | 10.0       | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 10.0  | U         | 2.40 | 10.0       | ug/L  |
| 132-64-9   | Dibenzofuran               | 5.00  | U         | 0.61 | 5.00       | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 5.00  | U         | 1.20 | 5.00       | ug/L  |
| 84-66-2    | Diethylphthalate           | 5.00  | U         | 0.69 | 5.00       | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 5.00  | U         | 0.68 | 5.00       | ug/L  |
| 86-73-7    | Fluorene                   | 5.00  | U         | 0.63 | 5.00       | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 5.00  | U         | 1.50 | 5.00       | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 10.0  | U         | 2.90 | 10.0       | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 5.00  | U         | 0.58 | 5.00       | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 5.00  | U         | 0.40 | 5.00       | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 5.00  | U         | 0.52 | 5.00       | ug/L  |
| 1912-24-9  | Atrazine                   | 5.00  | U         | 1.00 | 5.00       | ug/L  |
| 87-86-5    | Pentachlorophenol          | 10.0  | U         | 1.60 | 10.0       | ug/L  |
| 85-01-8    | Phenanthrene               | 5.00  | U         | 0.50 | 5.00       | ug/L  |
| 120-12-7   | Anthracene                 | 5.00  | U         | 0.61 | 5.00       | ug/L  |
| 86-74-8    | Carbazole                  | 5.00  | U         | 0.72 | 5.00       | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 5.00  | U         | 1.20 | 5.00       | ug/L  |
| 206-44-0   | Fluoranthene               | 5.00  | U         | 0.82 | 5.00       | ug/L  |
| 129-00-0   | Pyrene                     | 5.00  | U         | 0.50 | 5.00       | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 5.00  | U         | 1.90 | 5.00       | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 10.0  | U         | 0.93 | 10.0       | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 5.00  | U         | 0.45 | 5.00       | ug/L  |
| 218-01-9   | Chrysene                   | 5.00  | U         | 0.44 | 5.00       | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 5.00  | U         | 1.60 | 5.00       | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 10.0  | U         | 2.30 | 10.0       | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 5.00  | U         | 0.49 | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-10C-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-01                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 1000                               | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BF143581.D        | 1         | 08/22/25 10:34 | 08/22/25 19:21 | PB169362      |

| CAS Number                            | Parameter                          | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------------------|------------------------------------|--------|-----------|----------|------------|----------|
| 207-08-9                              | Benzo(k)fluoranthene               | 5.00   | U         | 0.48     | 5.00       | ug/L     |
| 50-32-8                               | Benzo(a)pyrene                     | 5.00   | U         | 0.55     | 5.00       | ug/L     |
| 193-39-5                              | Indeno(1,2,3-cd)pyrene             | 5.00   | U         | 0.59     | 5.00       | ug/L     |
| 53-70-3                               | Dibenzo(a,h)anthracene             | 5.00   | U         | 0.67     | 5.00       | ug/L     |
| 191-24-2                              | Benzo(g,h,i)perylene               | 5.00   | U         | 0.69     | 5.00       | ug/L     |
| 95-94-3                               | 1,2,4,5-Tetrachlorobenzene         | 5.00   | U         | 0.52     | 5.00       | ug/L     |
| 123-91-1                              | 1,4-Dioxane                        | 5.00   | U         | 1.00     | 5.00       | ug/L     |
| 58-90-2                               | 2,3,4,6-Tetrachlorophenol          | 5.00   | U         | 0.72     | 5.00       | ug/L     |
| <b>SURROGATES</b>                     |                                    |        |           |          |            |          |
| 367-12-4                              | 2-Fluorophenol                     | 59.1   |           | 23 - 138 | 39%        | SPK: 150 |
| 13127-88-3                            | Phenol-d6                          | 38.3   |           | 10 - 134 | 26%        | SPK: 150 |
| 4165-60-0                             | Nitrobenzene-d5                    | 99.8   |           | 67 - 132 | 100%       | SPK: 100 |
| 321-60-8                              | 2-Fluorobiphenyl                   | 88.3   |           | 52 - 132 | 88%        | SPK: 100 |
| 118-79-6                              | 2,4,6-Tribromophenol               | 151    |           | 44 - 137 | 101%       | SPK: 150 |
| 1718-51-0                             | Terphenyl-d14                      | 102    |           | 42 - 152 | 102%       | SPK: 100 |
| <b>INTERNAL STANDARDS</b>             |                                    |        |           |          |            |          |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4             | 105000 | 6.922     |          |            |          |
| 1146-65-2                             | Naphthalene-d8                     | 416000 | 8.204     |          |            |          |
| 15067-26-2                            | Acenaphthene-d10                   | 252000 | 9.963     |          |            |          |
| 1517-22-2                             | Phenanthrene-d10                   | 423000 | 11.451    |          |            |          |
| 1719-03-5                             | Chrysene-d12                       | 233000 | 14.092    |          |            |          |
| 1520-96-3                             | Perylene-d12                       | 204000 | 15.592    |          |            |          |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                                    |        |           |          |            |          |
| 000994-05-8                           | Butane, 2-methoxy-2-methyl-        | 87.3   | J         |          | 2.25       | ug/L     |
| 000123-42-2                           | 2-Pentanone, 4-hydroxy-4-methyl-   | 5.50   | AB        |          | 5.15       | ug/L     |
| 000149-57-5                           | Hexanoic acid, 2-ethyl-            | 5.30   | J         |          | 7.61       | ug/L     |
| 006422-86-2                           | 1,4-Benzenedicarboxylic acid, bis( | 2.50   | J         |          | 14.7       | ug/L     |

## Report of Analysis

|                    |                                    |              |                 |                  |        |
|--------------------|------------------------------------|--------------|-----------------|------------------|--------|
| Client:            | AECOM                              |              | Date Collected: | 08/20/25         |        |
| Project:           | National Grid Equity - Brooklyn NY |              | Date Received:  | 08/20/25         |        |
| Client Sample ID:  | MW-10C-082025                      |              | SDG No.:        | Q2924            |        |
| Lab Sample ID:     | Q2924-01                           |              | Matrix:         | Water            |        |
| Analytical Method: | 8270E                              |              | % Solid:        | 0                |        |
| Sample Wt/Vol:     | 1000                               | Units: mL    | Final Vol:      | 1000             | uL     |
| Soil Aliquot Vol:  |                                    | uL           | Test:           | SVOC-TCL BNA -20 |        |
| Extraction Type :  |                                    | Decanted :   | N               | Level :          | LOW    |
| Injection Volume : |                                    | GPC Factor : | 1.0             | GPC Cleanup :    | N PH : |
| Prep Method :      |                                    |              |                 |                  |        |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BF143581.D        | 1         | 08/22/25 10:34 | 08/22/25 19:21 | PB169362      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-201-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-02                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 950                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025601.D        | 1         | 08/22/25 10:34 | 08/27/25 13:37 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 100-52-7       | Benzaldehyde                | 10.5  | U         | 4.10 | 10.5       | ug/L  |
| 108-95-2       | Phenol                      | 5.30  | U         | 0.96 | 5.30       | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 5.30  | U         | 0.85 | 5.30       | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 5.30  | U         | 0.61 | 5.30       | ug/L  |
| 95-48-7        | 2-Methylphenol              | 5.30  | U         | 1.20 | 5.30       | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 5.30  | U         | 1.30 | 5.30       | ug/L  |
| 98-86-2        | Acetophenone                | 5.30  | U         | 0.78 | 5.30       | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 10.0  | J         | 1.20 | 10.5       | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 2.60  | U         | 1.50 | 2.60       | ug/L  |
| 67-72-1        | Hexachloroethane            | 5.30  | U         | 0.68 | 5.30       | ug/L  |
| 98-95-3        | Nitrobenzene                | 5.30  | U         | 0.80 | 5.30       | ug/L  |
| 78-59-1        | Isophorone                  | 5.30  | U         | 0.79 | 5.30       | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 5.30  | U         | 1.90 | 5.30       | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 5.30  | U         | 1.90 | 5.30       | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 5.30  | U         | 0.72 | 5.30       | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 5.30  | U         | 0.55 | 5.30       | ug/L  |
| 91-20-3        | Naphthalene                 | 3100  | E         | 0.53 | 5.30       | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 5.30  | U         | 0.88 | 5.30       | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 5.30  | U         | 0.57 | 5.30       | ug/L  |
| 105-60-2       | Caprolactam                 | 10.5  | U         | 1.20 | 10.5       | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 5.30  | U         | 0.62 | 5.30       | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 920   | E         | 0.59 | 5.30       | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 10.5  | U         | 3.80 | 10.5       | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 5.30  | U         | 0.54 | 5.30       | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 5.30  | U         | 0.65 | 5.30       | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 26.0  |           | 0.56 | 5.30       | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 5.30  | U         | 0.64 | 5.30       | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 5.30  | U         | 1.30 | 5.30       | ug/L  |
| 131-11-3       | Dimethylphthalate           | 5.30  | U         | 0.64 | 5.30       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-201-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-02                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 950                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025601.D        | 1         | 08/22/25 10:34 | 08/27/25 13:37 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 160   | E         | 0.79 | 5.30       | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 5.30  | U         | 0.97 | 5.30       | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 5.30  | U         | 1.10 | 5.30       | ug/L  |
| 83-32-9    | Acenaphthene               | 8.70  |           | 0.58 | 5.30       | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 10.5  | U         | 6.30 | 10.5       | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 10.5  | U         | 2.50 | 10.5       | ug/L  |
| 132-64-9   | Dibenzofuran               | 4.00  | J         | 0.64 | 5.30       | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 5.30  | U         | 1.30 | 5.30       | ug/L  |
| 84-66-2    | Diethylphthalate           | 2.10  | J         | 0.73 | 5.30       | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 5.30  | U         | 0.72 | 5.30       | ug/L  |
| 86-73-7    | Fluorene                   | 27.4  |           | 0.66 | 5.30       | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 5.30  | U         | 1.60 | 5.30       | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 10.5  | U         | 3.00 | 10.5       | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 5.30  | U         | 0.61 | 5.30       | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 5.30  | U         | 0.42 | 5.30       | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 5.30  | U         | 0.55 | 5.30       | ug/L  |
| 1912-24-9  | Atrazine                   | 5.30  | U         | 1.10 | 5.30       | ug/L  |
| 87-86-5    | Pentachlorophenol          | 10.5  | U         | 1.70 | 10.5       | ug/L  |
| 85-01-8    | Phenanthrene               | 29.6  |           | 0.53 | 5.30       | ug/L  |
| 120-12-7   | Anthracene                 | 6.10  |           | 0.64 | 5.30       | ug/L  |
| 86-74-8    | Carbazole                  | 13.8  |           | 0.76 | 5.30       | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 2.70  | J         | 1.30 | 5.30       | ug/L  |
| 206-44-0   | Fluoranthene               | 2.30  | J         | 0.86 | 5.30       | ug/L  |
| 129-00-0   | Pyrene                     | 3.00  | J         | 0.53 | 5.30       | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 5.30  | U         | 2.00 | 5.30       | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 10.5  | U         | 0.98 | 10.5       | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 5.30  | U         | 0.47 | 5.30       | ug/L  |
| 218-01-9   | Chrysene                   | 5.30  | U         | 0.46 | 5.30       | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 5.30  | U         | 1.70 | 5.30       | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 10.5  | U         | 2.50 | 10.5       | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 5.30  | U         | 0.52 | 5.30       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-201-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-02                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 950                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025601.D        | 1         | 08/22/25 10:34 | 08/27/25 13:37 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 207-08-9   | Benzo(k)fluoranthene       | 5.30  | U         | 0.51 | 5.30       | ug/L  |
| 50-32-8    | Benzo(a)pyrene             | 5.30  | U         | 0.58 | 5.30       | ug/L  |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 5.30  | U         | 0.62 | 5.30       | ug/L  |
| 53-70-3    | Dibenzo(a,h)anthracene     | 5.30  | U         | 0.71 | 5.30       | ug/L  |
| 191-24-2   | Benzo(g,h,i)perylene       | 5.30  | U         | 0.73 | 5.30       | ug/L  |
| 95-94-3    | 1,2,4,5-Tetrachlorobenzene | 5.30  | U         | 0.55 | 5.30       | ug/L  |
| 123-91-1   | 1,4-Dioxane                | 5.30  | U         | 1.10 | 5.30       | ug/L  |
| 58-90-2    | 2,3,4,6-Tetrachlorophenol  | 5.30  | U         | 0.76 | 5.30       | ug/L  |

### SURROGATES

|            |                      |      |   |          |      |          |
|------------|----------------------|------|---|----------|------|----------|
| 367-12-4   | 2-Fluorophenol       | 65.2 |   | 23 - 138 | 43%  | SPK: 150 |
| 13127-88-3 | Phenol-d6            | 38.7 |   | 10 - 134 | 26%  | SPK: 150 |
| 4165-60-0  | Nitrobenzene-d5      | 171  | * | 67 - 132 | 171% | SPK: 100 |
| 321-60-8   | 2-Fluorobiphenyl     | 78.8 |   | 52 - 132 | 79%  | SPK: 100 |
| 118-79-6   | 2,4,6-Tribromophenol | 136  |   | 44 - 137 | 91%  | SPK: 150 |
| 1718-51-0  | Terphenyl-d14        | 74.0 |   | 42 - 152 | 74%  | SPK: 100 |

### INTERNAL STANDARDS

|            |                        |         |        |  |
|------------|------------------------|---------|--------|--|
| 3855-82-1  | 1,4-Dichlorobenzene-d4 | 185000  | 7.778  |  |
| 1146-65-2  | Naphthalene-d8         | 347000  | 10.566 |  |
| 15067-26-2 | Acenaphthene-d10       | 427000  | 14.389 |  |
| 1517-22-2  | Phenanthrene-d10       | 843000  | 17.195 |  |
| 1719-03-5  | Chrysene-d12           | 909000  | 21.636 |  |
| 1520-96-3  | Perylene-d12           | 1060000 | 25.001 |  |

### TENTATIVE IDENTIFIED COMPOUNDS

|             |                             |      |   |      |      |
|-------------|-----------------------------|------|---|------|------|
| 000994-05-8 | Butane, 2-methoxy-2-methyl- | 11.5 | J | 3.03 | ug/L |
| 000616-44-4 | Thiophene, 3-methyl-        | 4.60 | J | 4.19 | ug/L |
| 000106-42-3 | p-Xylene                    | 160  | J | 5.47 | ug/L |
| 000620-14-4 | Benzene, 1-ethyl-3-methyl-  | 12.8 | J | 6.88 | ug/L |
| 000622-96-8 | Benzene, 1-ethyl-4-methyl-  | 5.30 | J | 7.18 | ug/L |
| 000098-83-9 | .alpha.-Methylstyrene       | 11.1 | J | 7.21 | ug/L |
|             | unknown7.449                | 99.8 | J | 7.45 | ug/L |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-201-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-02                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 950                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025601.D        | 1         | 08/22/25 10:34 | 08/27/25 13:37 | PB169362      |

| CAS Number  | Parameter                       | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|-------------|---------------------------------|-------|-----------|-----|------------|-------|
|             | unknown7.525                    | 31.2  | J         |     | 7.53       | ug/L  |
| 000637-50-3 | Benzene, 1-propenyl-            | 9.50  | J         |     | 7.96       | ug/L  |
| 000496-11-7 | Indane                          | 12.7  | J         |     | 8.15       | ug/L  |
|             | unknown8.372                    | 310   | J         |     | 8.37       | ug/L  |
| 000768-49-0 | Benzene, (2-methyl-1-propenyl)- | 7.90  | J         |     | 9.03       | ug/L  |
| 000275-51-4 | Azulene                         | 21.1  | J         |     | 10.7       | ug/L  |
| 90-12-0     | 1-Methylnaphthalene             | 600   | J         |     | 12.4       | ug/L  |
| 000575-37-1 | Naphthalene, 1,7-dimethyl-      | 4.60  | J         |     | 13.6       | ug/L  |
| 000575-41-7 | Naphthalene, 1,3-dimethyl-      | 5.90  | J         |     | 13.7       | ug/L  |
| 003770-48-7 | 4-Methylcarbazole               | 16.2  | J         |     | 18.1       | ug/L  |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-201-082025DL                    |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-02DL                         |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 950                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025607.D        | 20        | 08/22/25 10:34 | 08/27/25 17:44 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 100-52-7       | Benzaldehyde                | 210   | UD        | 82.3 | 210        | ug/L  |
| 108-95-2       | Phenol                      | 110   | UD        | 19.2 | 110        | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 110   | UD        | 17.1 | 110        | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 110   | UD        | 12.2 | 110        | ug/L  |
| 95-48-7        | 2-Methylphenol              | 110   | UD        | 23.6 | 110        | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 110   | UD        | 26.9 | 110        | ug/L  |
| 98-86-2        | Acetophenone                | 110   | UD        | 15.6 | 110        | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 210   | UD        | 23.2 | 210        | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 52.6  | UD        | 29.7 | 52.6       | ug/L  |
| 67-72-1        | Hexachloroethane            | 110   | UD        | 13.7 | 110        | ug/L  |
| 98-95-3        | Nitrobenzene                | 110   | UD        | 16.0 | 110        | ug/L  |
| 78-59-1        | Isophorone                  | 110   | UD        | 15.8 | 110        | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 110   | UD        | 37.1 | 110        | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 110   | UD        | 38.9 | 110        | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 110   | UD        | 14.3 | 110        | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 110   | UD        | 10.9 | 110        | ug/L  |
| 91-20-3        | Naphthalene                 | 6000  | ED        | 10.5 | 110        | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 110   | UD        | 17.7 | 110        | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 110   | UD        | 11.4 | 110        | ug/L  |
| 105-60-2       | Caprolactam                 | 210   | UD        | 23.8 | 210        | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 110   | UD        | 12.4 | 110        | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 610   | D         | 11.8 | 110        | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 210   | UD        | 76.4 | 210        | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 110   | UD        | 10.7 | 110        | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 110   | UD        | 13.1 | 110        | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 110   | UD        | 11.2 | 110        | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 110   | UD        | 12.8 | 110        | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 110   | UD        | 26.5 | 110        | ug/L  |
| 131-11-3       | Dimethylphthalate           | 110   | UD        | 12.8 | 110        | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-201-082025DL                    |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-02DL                         |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 950                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025607.D        | 20        | 08/22/25 10:34 | 08/27/25 17:44 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 180   | D         | 15.8 | 110        | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 110   | UD        | 19.4 | 110        | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 110   | UD        | 22.1 | 110        | ug/L  |
| 83-32-9    | Acenaphthene               | 110   | UD        | 11.6 | 110        | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 210   | UD        | 130  | 210        | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 210   | UD        | 50.1 | 210        | ug/L  |
| 132-64-9   | Dibenzofuran               | 110   | UD        | 12.8 | 110        | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 110   | UD        | 25.7 | 110        | ug/L  |
| 84-66-2    | Diethylphthalate           | 110   | UD        | 14.5 | 110        | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 110   | UD        | 14.3 | 110        | ug/L  |
| 86-73-7    | Fluorene                   | 110   | UD        | 13.3 | 110        | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 110   | UD        | 31.6 | 110        | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 210   | UD        | 60.6 | 210        | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 110   | UD        | 12.2 | 110        | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 110   | UD        | 8.40 | 110        | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 110   | UD        | 10.9 | 110        | ug/L  |
| 1912-24-9  | Atrazine                   | 110   | UD        | 21.3 | 110        | ug/L  |
| 87-86-5    | Pentachlorophenol          | 210   | UD        | 33.3 | 210        | ug/L  |
| 85-01-8    | Phenanthrene               | 110   | UD        | 10.5 | 110        | ug/L  |
| 120-12-7   | Anthracene                 | 110   | UD        | 12.8 | 110        | ug/L  |
| 86-74-8    | Carbazole                  | 110   | UD        | 15.2 | 110        | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 110   | UD        | 25.7 | 110        | ug/L  |
| 206-44-0   | Fluoranthene               | 110   | UD        | 17.3 | 110        | ug/L  |
| 129-00-0   | Pyrene                     | 110   | UD        | 10.5 | 110        | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 110   | UD        | 40.6 | 110        | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 210   | UD        | 19.6 | 210        | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 110   | UD        | 9.50 | 110        | ug/L  |
| 218-01-9   | Chrysene                   | 110   | UD        | 9.30 | 110        | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 110   | UD        | 33.7 | 110        | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 210   | UD        | 49.3 | 210        | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 110   | UD        | 10.3 | 110        | ug/L  |

## Report of Analysis

|                    |                                    |                 |                  |
|--------------------|------------------------------------|-----------------|------------------|
| Client:            | AECOM                              | Date Collected: | 08/20/25         |
| Project:           | National Grid Equity - Brooklyn NY | Date Received:  | 08/20/25         |
| Client Sample ID:  | MW-201-082025DL                    | SDG No.:        | Q2924            |
| Lab Sample ID:     | Q2924-02DL                         | Matrix:         | Water            |
| Analytical Method: | 8270E                              | % Solid:        | 0                |
| Sample Wt/Vol:     | 950 Units: mL                      | Final Vol:      | 1000 uL          |
| Soil Aliquot Vol:  | uL                                 | Test:           | SVOC-TCL BNA -20 |
| Extraction Type :  | Decanted : N                       | Level :         | LOW              |
| Injection Volume : | GPC Factor : 1.0                   | GPC Cleanup :   | N PH :           |
| Prep Method :      |                                    |                 |                  |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025607.D        | 20        | 08/22/25 10:34 | 08/27/25 17:44 | PB169362      |

| CAS Number                | Parameter                  | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------|----------------------------|---------|-----------|----------|------------|----------|
| 207-08-9                  | Benzo(k)fluoranthene       | 110     | UD        | 10.1     | 110        | ug/L     |
| 50-32-8                   | Benzo(a)pyrene             | 110     | UD        | 11.6     | 110        | ug/L     |
| 193-39-5                  | Indeno(1,2,3-cd)pyrene     | 110     | UD        | 12.4     | 110        | ug/L     |
| 53-70-3                   | Dibenzo(a,h)anthracene     | 110     | UD        | 14.1     | 110        | ug/L     |
| 191-24-2                  | Benzo(g,h,i)perylene       | 110     | UD        | 14.5     | 110        | ug/L     |
| 95-94-3                   | 1,2,4,5-Tetrachlorobenzene | 110     | UD        | 10.9     | 110        | ug/L     |
| 123-91-1                  | 1,4-Dioxane                | 110     | UD        | 21.1     | 110        | ug/L     |
| 58-90-2                   | 2,3,4,6-Tetrachlorophenol  | 110     | UD        | 15.2     | 110        | ug/L     |
| <b>SURROGATES</b>         |                            |         |           |          |            |          |
| 367-12-4                  | 2-Fluorophenol             | 63.7    |           | 23 - 138 | 42%        | SPK: 150 |
| 13127-88-3                | Phenol-d6                  | 31.8    |           | 10 - 134 | 21%        | SPK: 150 |
| 4165-60-0                 | Nitrobenzene-d5            | 77.1    |           | 67 - 132 | 77%        | SPK: 100 |
| 321-60-8                  | 2-Fluorobiphenyl           | 83.0    |           | 52 - 132 | 83%        | SPK: 100 |
| 118-79-6                  | 2,4,6-Tribromophenol       | 114     |           | 44 - 137 | 76%        | SPK: 150 |
| 1718-51-0                 | Terphenyl-d14              | 77.0    |           | 42 - 152 | 77%        | SPK: 100 |
| <b>INTERNAL STANDARDS</b> |                            |         |           |          |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4     | 147000  | 7.772     |          |            |          |
| 1146-65-2                 | Naphthalene-d8             | 575000  | 10.543    |          |            |          |
| 15067-26-2                | Acenaphthene-d10           | 353000  | 14.389    |          |            |          |
| 1517-22-2                 | Phenanthrene-d10           | 694000  | 17.189    |          |            |          |
| 1719-03-5                 | Chrysene-d12               | 837000  | 21.636    |          |            |          |
| 1520-96-3                 | Perylene-d12               | 1050000 | 25.001    |          |            |          |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |                 |                  |
|--------------------|------------------------------------|-----------------|------------------|
| Client:            | AECOM                              | Date Collected: | 08/20/25         |
| Project:           | National Grid Equity - Brooklyn NY | Date Received:  | 08/20/25         |
| Client Sample ID:  | MW-201-082025DL2                   | SDG No.:        | Q2924            |
| Lab Sample ID:     | Q2924-02DL2                        | Matrix:         | Water            |
| Analytical Method: | 8270E                              | % Solid:        | 0                |
| Sample Wt/Vol:     | 950      Units:    mL              | Final Vol:      | 1000      uL     |
| Soil Aliquot Vol:  | uL                                 | Test:           | SVOC-TCL BNA -20 |
| Extraction Type :  | Decanted :    N                    | Level :         | LOW              |
| Injection Volume : | GPC Factor :    1.0                | GPC Cleanup :   | N      PH :      |
| Prep Method :      |                                    |                 |                  |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025608.D        | 200       | 08/22/25 10:34 | 08/27/25 18:26 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|-----|------------|-------|
| <b>TARGETS</b> |                             |       |           |     |            |       |
| 100-52-7       | Benzaldehyde                | 2100  | UD        | 820 | 2100       | ug/L  |
| 108-95-2       | Phenol                      | 1100  | UD        | 190 | 1100       | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 1100  | UD        | 170 | 1100       | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 1100  | UD        | 120 | 1100       | ug/L  |
| 95-48-7        | 2-Methylphenol              | 1100  | UD        | 240 | 1100       | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 1100  | UD        | 270 | 1100       | ug/L  |
| 98-86-2        | Acetophenone                | 1100  | UD        | 160 | 1100       | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 2100  | UD        | 230 | 2100       | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 530   | UD        | 300 | 530        | ug/L  |
| 67-72-1        | Hexachloroethane            | 1100  | UD        | 140 | 1100       | ug/L  |
| 98-95-3        | Nitrobenzene                | 1100  | UD        | 160 | 1100       | ug/L  |
| 78-59-1        | Isophorone                  | 1100  | UD        | 160 | 1100       | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 1100  | UD        | 370 | 1100       | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 1100  | UD        | 390 | 1100       | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 1100  | UD        | 140 | 1100       | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 1100  | UD        | 110 | 1100       | ug/L  |
| 91-20-3        | Naphthalene                 | 6800  | D         | 110 | 1100       | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 1100  | UD        | 180 | 1100       | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 1100  | UD        | 110 | 1100       | ug/L  |
| 105-60-2       | Caprolactam                 | 2100  | UD        | 240 | 2100       | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 1100  | UD        | 120 | 1100       | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 570   | JD        | 120 | 1100       | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 2100  | UD        | 760 | 2100       | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 1100  | UD        | 110 | 1100       | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 1100  | UD        | 130 | 1100       | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 1100  | UD        | 110 | 1100       | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 1100  | UD        | 130 | 1100       | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 1100  | UD        | 270 | 1100       | ug/L  |
| 131-11-3       | Dimethylphthalate           | 1100  | UD        | 130 | 1100       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-201-082025DL2                   |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-02DL2                        |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 950                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025608.D        | 200       | 08/22/25 10:34 | 08/27/25 18:26 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 1100  | UD        | 160  | 1100       | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 1100  | UD        | 190  | 1100       | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 1100  | UD        | 220  | 1100       | ug/L  |
| 83-32-9    | Acenaphthene               | 1100  | UD        | 120  | 1100       | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 2100  | UD        | 1300 | 2100       | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 2100  | UD        | 500  | 2100       | ug/L  |
| 132-64-9   | Dibenzofuran               | 1100  | UD        | 130  | 1100       | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 1100  | UD        | 260  | 1100       | ug/L  |
| 84-66-2    | Diethylphthalate           | 1100  | UD        | 150  | 1100       | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 1100  | UD        | 140  | 1100       | ug/L  |
| 86-73-7    | Fluorene                   | 1100  | UD        | 130  | 1100       | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 1100  | UD        | 320  | 1100       | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 2100  | UD        | 610  | 2100       | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 1100  | UD        | 120  | 1100       | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 1100  | UD        | 84.2 | 1100       | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 1100  | UD        | 110  | 1100       | ug/L  |
| 1912-24-9  | Atrazine                   | 1100  | UD        | 210  | 1100       | ug/L  |
| 87-86-5    | Pentachlorophenol          | 2100  | UD        | 330  | 2100       | ug/L  |
| 85-01-8    | Phenanthrene               | 1100  | UD        | 110  | 1100       | ug/L  |
| 120-12-7   | Anthracene                 | 1100  | UD        | 130  | 1100       | ug/L  |
| 86-74-8    | Carbazole                  | 1100  | UD        | 150  | 1100       | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 1100  | UD        | 260  | 1100       | ug/L  |
| 206-44-0   | Fluoranthene               | 1100  | UD        | 170  | 1100       | ug/L  |
| 129-00-0   | Pyrene                     | 1100  | UD        | 110  | 1100       | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 1100  | UD        | 410  | 1100       | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 2100  | UD        | 200  | 2100       | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 1100  | UD        | 94.7 | 1100       | ug/L  |
| 218-01-9   | Chrysene                   | 1100  | UD        | 92.6 | 1100       | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 1100  | UD        | 340  | 1100       | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 2100  | UD        | 490  | 2100       | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 1100  | UD        | 100  | 1100       | ug/L  |

## Report of Analysis

|                    |                                    |                 |                  |
|--------------------|------------------------------------|-----------------|------------------|
| Client:            | AECOM                              | Date Collected: | 08/20/25         |
| Project:           | National Grid Equity - Brooklyn NY | Date Received:  | 08/20/25         |
| Client Sample ID:  | MW-201-082025DL2                   | SDG No.:        | Q2924            |
| Lab Sample ID:     | Q2924-02DL2                        | Matrix:         | Water            |
| Analytical Method: | 8270E                              | % Solid:        | 0                |
| Sample Wt/Vol:     | 950 Units: mL                      | Final Vol:      | 1000 uL          |
| Soil Aliquot Vol:  | uL                                 | Test:           | SVOC-TCL BNA -20 |
| Extraction Type :  | Decanted : N                       | Level :         | LOW              |
| Injection Volume : | GPC Factor : 1.0                   | GPC Cleanup :   | N PH :           |
| Prep Method :      |                                    |                 |                  |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025608.D        | 200       | 08/22/25 10:34 | 08/27/25 18:26 | PB169362      |

| CAS Number                | Parameter                  | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------|----------------------------|---------|-----------|----------|------------|----------|
| 207-08-9                  | Benzo(k)fluoranthene       | 1100    | UD        | 100      | 1100       | ug/L     |
| 50-32-8                   | Benzo(a)pyrene             | 1100    | UD        | 120      | 1100       | ug/L     |
| 193-39-5                  | Indeno(1,2,3-cd)pyrene     | 1100    | UD        | 120      | 1100       | ug/L     |
| 53-70-3                   | Dibenzo(a,h)anthracene     | 1100    | UD        | 140      | 1100       | ug/L     |
| 191-24-2                  | Benzo(g,h,i)perylene       | 1100    | UD        | 150      | 1100       | ug/L     |
| 95-94-3                   | 1,2,4,5-Tetrachlorobenzene | 1100    | UD        | 110      | 1100       | ug/L     |
| 123-91-1                  | 1,4-Dioxane                | 1100    | UD        | 210      | 1100       | ug/L     |
| 58-90-2                   | 2,3,4,6-Tetrachlorophenol  | 1100    | UD        | 150      | 1100       | ug/L     |
| <b>SURROGATES</b>         |                            |         |           |          |            |          |
| 367-12-4                  | 2-Fluorophenol             | 67.8    |           | 23 - 138 | 45%        | SPK: 150 |
| 13127-88-3                | Phenol-d6                  | 14.0    | *         | 10 - 134 | 9%         | SPK: 150 |
| 4165-60-0                 | Nitrobenzene-d5            | 69.0    |           | 67 - 132 | 69%        | SPK: 100 |
| 321-60-8                  | 2-Fluorobiphenyl           | 76.2    |           | 52 - 132 | 76%        | SPK: 100 |
| 118-79-6                  | 2,4,6-Tribromophenol       | 91.2    |           | 44 - 137 | 61%        | SPK: 150 |
| 1718-51-0                 | Terphenyl-d14              | 73.4    |           | 42 - 152 | 73%        | SPK: 100 |
| <b>INTERNAL STANDARDS</b> |                            |         |           |          |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4     | 151000  | 7.772     |          |            |          |
| 1146-65-2                 | Naphthalene-d8             | 582000  | 10.542    |          |            |          |
| 15067-26-2                | Acenaphthene-d10           | 354000  | 14.389    |          |            |          |
| 1517-22-2                 | Phenanthrene-d10           | 660000  | 17.183    |          |            |          |
| 1719-03-5                 | Chrysene-d12               | 857000  | 21.636    |          |            |          |
| 1520-96-3                 | Perylene-d12               | 1040000 | 25.006    |          |            |          |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-2B-082025                       |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-05                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 960                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BF143582.D        | 1         | 08/22/25 10:34 | 08/22/25 19:50 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 100-52-7       | Benzaldehyde                | 10.4  | U         | 4.10 | 10.4       | ug/L  |
| 108-95-2       | Phenol                      | 5.10  | J         | 0.95 | 5.20       | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 5.20  | U         | 0.84 | 5.20       | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 5.20  | U         | 0.60 | 5.20       | ug/L  |
| 95-48-7        | 2-Methylphenol              | 5.20  | U         | 1.20 | 5.20       | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 5.20  | U         | 1.30 | 5.20       | ug/L  |
| 98-86-2        | Acetophenone                | 5.20  | U         | 0.77 | 5.20       | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 10.4  | U         | 1.10 | 10.4       | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 2.60  | U         | 1.50 | 2.60       | ug/L  |
| 67-72-1        | Hexachloroethane            | 5.20  | U         | 0.68 | 5.20       | ug/L  |
| 98-95-3        | Nitrobenzene                | 5.20  | U         | 0.79 | 5.20       | ug/L  |
| 78-59-1        | Isophorone                  | 5.20  | U         | 0.78 | 5.20       | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 5.20  | U         | 1.80 | 5.20       | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 5.20  | U         | 1.90 | 5.20       | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 5.20  | U         | 0.71 | 5.20       | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 5.20  | U         | 0.54 | 5.20       | ug/L  |
| 91-20-3        | Naphthalene                 | 930   | E         | 0.52 | 5.20       | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 5.20  | U         | 0.88 | 5.20       | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 5.20  | U         | 0.56 | 5.20       | ug/L  |
| 105-60-2       | Caprolactam                 | 10.4  | U         | 1.20 | 10.4       | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 5.20  | U         | 0.61 | 5.20       | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 160   | E         | 0.58 | 5.20       | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 10.4  | U         | 3.80 | 10.4       | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 5.20  | U         | 0.53 | 5.20       | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 5.20  | U         | 0.65 | 5.20       | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 12.7  |           | 0.55 | 5.20       | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 5.20  | U         | 0.64 | 5.20       | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 5.20  | U         | 1.30 | 5.20       | ug/L  |
| 131-11-3       | Dimethylphthalate           | 5.20  | U         | 0.64 | 5.20       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-2B-082025                       |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-05                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 960                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BF143582.D        | 1         | 08/22/25 10:34 | 08/22/25 19:50 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 5.20  | U         | 0.78 | 5.20       | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 5.20  | U         | 0.96 | 5.20       | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 5.20  | U         | 1.10 | 5.20       | ug/L  |
| 83-32-9    | Acenaphthene               | 61.2  |           | 0.57 | 5.20       | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 10.4  | U         | 6.20 | 10.4       | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 10.4  | U         | 2.50 | 10.4       | ug/L  |
| 132-64-9   | Dibenzofuran               | 2.60  | J         | 0.64 | 5.20       | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 5.20  | U         | 1.30 | 5.20       | ug/L  |
| 84-66-2    | Diethylphthalate           | 5.20  | U         | 0.72 | 5.20       | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 5.20  | U         | 0.71 | 5.20       | ug/L  |
| 86-73-7    | Fluorene                   | 13.0  |           | 0.66 | 5.20       | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 5.20  | U         | 1.60 | 5.20       | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 10.4  | U         | 3.00 | 10.4       | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 5.20  | U         | 0.60 | 5.20       | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 5.20  | U         | 0.42 | 5.20       | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 5.20  | U         | 0.54 | 5.20       | ug/L  |
| 1912-24-9  | Atrazine                   | 5.20  | U         | 1.10 | 5.20       | ug/L  |
| 87-86-5    | Pentachlorophenol          | 10.4  | U         | 1.60 | 10.4       | ug/L  |
| 85-01-8    | Phenanthrene               | 17.2  |           | 0.52 | 5.20       | ug/L  |
| 120-12-7   | Anthracene                 | 2.60  | J         | 0.64 | 5.20       | ug/L  |
| 86-74-8    | Carbazole                  | 5.80  |           | 0.75 | 5.20       | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 5.20  | U         | 1.30 | 5.20       | ug/L  |
| 206-44-0   | Fluoranthene               | 5.20  | U         | 0.85 | 5.20       | ug/L  |
| 129-00-0   | Pyrene                     | 5.20  | U         | 0.52 | 5.20       | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 5.20  | U         | 2.00 | 5.20       | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 10.4  | U         | 0.97 | 10.4       | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 5.20  | U         | 0.47 | 5.20       | ug/L  |
| 218-01-9   | Chrysene                   | 5.20  | U         | 0.46 | 5.20       | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 5.20  | U         | 1.70 | 5.20       | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 10.4  | U         | 2.40 | 10.4       | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 5.20  | U         | 0.51 | 5.20       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-2B-082025                       |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-05                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 960                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BF143582.D        | 1         | 08/22/25 10:34 | 08/22/25 19:50 | PB169362      |

| CAS Number                            | Parameter                          | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------------------|------------------------------------|--------|-----------|----------|------------|----------|
| 207-08-9                              | Benzo(k)fluoranthene               | 5.20   | U         | 0.50     | 5.20       | ug/L     |
| 50-32-8                               | Benzo(a)pyrene                     | 5.20   | U         | 0.57     | 5.20       | ug/L     |
| 193-39-5                              | Indeno(1,2,3-cd)pyrene             | 5.20   | U         | 0.61     | 5.20       | ug/L     |
| 53-70-3                               | Dibenzo(a,h)anthracene             | 5.20   | U         | 0.70     | 5.20       | ug/L     |
| 191-24-2                              | Benzo(g,h,i)perylene               | 5.20   | U         | 0.72     | 5.20       | ug/L     |
| 95-94-3                               | 1,2,4,5-Tetrachlorobenzene         | 5.20   | U         | 0.54     | 5.20       | ug/L     |
| 123-91-1                              | 1,4-Dioxane                        | 5.20   | U         | 1.00     | 5.20       | ug/L     |
| 58-90-2                               | 2,3,4,6-Tetrachlorophenol          | 5.20   | U         | 0.75     | 5.20       | ug/L     |
| <b>SURROGATES</b>                     |                                    |        |           |          |            |          |
| 367-12-4                              | 2-Fluorophenol                     | 73.5   |           | 23 - 138 | 49%        | SPK: 150 |
| 13127-88-3                            | Phenol-d6                          | 53.5   |           | 10 - 134 | 36%        | SPK: 150 |
| 4165-60-0                             | Nitrobenzene-d5                    | 118    |           | 67 - 132 | 118%       | SPK: 100 |
| 321-60-8                              | 2-Fluorobiphenyl                   | 87.5   |           | 52 - 132 | 87%        | SPK: 100 |
| 118-79-6                              | 2,4,6-Tribromophenol               | 141    |           | 44 - 137 | 94%        | SPK: 150 |
| 1718-51-0                             | Terphenyl-d14                      | 101    |           | 42 - 152 | 101%       | SPK: 100 |
| <b>INTERNAL STANDARDS</b>             |                                    |        |           |          |            |          |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4             | 98000  | 6.928     |          |            |          |
| 1146-65-2                             | Naphthalene-d8                     | 310000 | 8.228     |          |            |          |
| 15067-26-2                            | Acenaphthene-d10                   | 221000 | 9.963     |          |            |          |
| 1517-22-2                             | Phenanthrene-d10                   | 347000 | 11.451    |          |            |          |
| 1719-03-5                             | Chrysene-d12                       | 201000 | 14.092    |          |            |          |
| 1520-96-3                             | Perylene-d12                       | 183000 | 15.592    |          |            |          |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                                    |        |           |          |            |          |
| 000994-05-8                           | Butane, 2-methoxy-2-methyl-        | 110    | J         |          | 2.25       | ug/L     |
| 002175-91-9                           | 1,3-Cyclopentadiene, 5-(1-methylet | 290    | J         |          | 5.44       | ug/L     |
| 000108-38-3                           | Benzene, 1,3-dimethyl-             | 28.6   | J         |          | 5.79       | ug/L     |
|                                       | unknown6.387                       | 14.0   | J         |          | 6.39       | ug/L     |
|                                       | unknown6.751                       | 40.5   | J         |          | 6.75       | ug/L     |
|                                       | unknown6.987                       | 76.5   | J         |          | 6.99       | ug/L     |
|                                       | unknown7.034                       | 14.1   | J         |          | 7.03       | ug/L     |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-2B-082025                       |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-05                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 960                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BF143582.D        | 1         | 08/22/25 10:34 | 08/22/25 19:50 | PB169362      |

| CAS Number  | Parameter                  | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|-------------|----------------------------|-------|-----------|-----|------------|-------|
|             | unknown7.134               | 770   | J         |     | 7.13       | ug/L  |
|             | unknown7.193               | 35.4  | J         |     | 7.19       | ug/L  |
|             | unknown7.951               | 2.30  | J         |     | 7.95       | ug/L  |
| 90-12-0     | 1-Methylnaphthalene        | 240   | J         |     | 9.03       | ug/L  |
| 001127-76-0 | Naphthalene, 1-ethyl-      | 8.40  | J         |     | 9.47       | ug/L  |
| 000575-43-9 | Naphthalene, 1,6-dimethyl- | 15.6  | J         |     | 9.54       | ug/L  |
| 000581-40-8 | Naphthalene, 2,3-dimethyl- | 18.7  | J         |     | 9.62       | ug/L  |
| 000581-42-0 | Naphthalene, 2,6-dimethyl- | 5.90  | J         |     | 9.64       | ug/L  |
| 000571-61-9 | Naphthalene, 1,5-dimethyl- | 8.90  | J         |     | 9.74       | ug/L  |
| 000090-15-3 | 1-Naphthalenol             | 7.10  | J         |     | 10.1       | ug/L  |
| 007469-77-4 | 1-Naphthalenol, 2-methyl-  | 8.80  | J         |     | 10.6       | ug/L  |
| 001484-12-4 | 9H-Carbazole, 9-methyl-    | 7.70  | J         |     | 12.0       | ug/L  |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-2B-082025DL                     |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-05DL                         |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 960                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025583.D        | 20        | 08/22/25 10:34 | 08/26/25 13:10 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 100-52-7       | Benzaldehyde                | 210   | UD        | 81.5 | 210        | ug/L  |
| 108-95-2       | Phenol                      | 100   | UD        | 19.0 | 100        | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 100   | UD        | 16.9 | 100        | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 100   | UD        | 12.1 | 100        | ug/L  |
| 95-48-7        | 2-Methylphenol              | 100   | UD        | 23.3 | 100        | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 100   | UD        | 26.7 | 100        | ug/L  |
| 98-86-2        | Acetophenone                | 100   | UD        | 15.4 | 100        | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 210   | UD        | 22.9 | 210        | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 52.1  | UD        | 29.4 | 52.1       | ug/L  |
| 67-72-1        | Hexachloroethane            | 100   | UD        | 13.5 | 100        | ug/L  |
| 98-95-3        | Nitrobenzene                | 100   | UD        | 15.8 | 100        | ug/L  |
| 78-59-1        | Isophorone                  | 100   | UD        | 15.6 | 100        | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 100   | UD        | 36.7 | 100        | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 100   | UD        | 38.5 | 100        | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 100   | UD        | 14.2 | 100        | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 100   | UD        | 10.8 | 100        | ug/L  |
| 91-20-3        | Naphthalene                 | 1800  | ED        | 10.4 | 100        | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 100   | UD        | 17.5 | 100        | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 100   | UD        | 11.3 | 100        | ug/L  |
| 105-60-2       | Caprolactam                 | 210   | UD        | 23.5 | 210        | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 100   | UD        | 12.3 | 100        | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 180   | D         | 11.7 | 100        | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 210   | UD        | 75.6 | 210        | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 100   | UD        | 10.6 | 100        | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 100   | UD        | 12.9 | 100        | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 100   | UD        | 11.0 | 100        | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 100   | UD        | 12.7 | 100        | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 100   | UD        | 26.3 | 100        | ug/L  |
| 131-11-3       | Dimethylphthalate           | 100   | UD        | 12.7 | 100        | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-2B-082025DL                     |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-05DL                         |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 960                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025583.D        | 20        | 08/22/25 10:34 | 08/26/25 13:10 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 100   | UD        | 15.6 | 100        | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 100   | UD        | 19.2 | 100        | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 100   | UD        | 21.9 | 100        | ug/L  |
| 83-32-9    | Acenaphthene               | 72.5  | JD        | 11.5 | 100        | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 210   | UD        | 120  | 210        | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 210   | UD        | 49.6 | 210        | ug/L  |
| 132-64-9   | Dibenzofuran               | 100   | UD        | 12.7 | 100        | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 100   | UD        | 25.4 | 100        | ug/L  |
| 84-66-2    | Diethylphthalate           | 100   | UD        | 14.4 | 100        | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 100   | UD        | 14.2 | 100        | ug/L  |
| 86-73-7    | Fluorene                   | 100   | UD        | 13.1 | 100        | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 100   | UD        | 31.3 | 100        | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 210   | UD        | 60.0 | 210        | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 100   | UD        | 12.1 | 100        | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 100   | UD        | 8.30 | 100        | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 100   | UD        | 10.8 | 100        | ug/L  |
| 1912-24-9  | Atrazine                   | 100   | UD        | 21.0 | 100        | ug/L  |
| 87-86-5    | Pentachlorophenol          | 210   | UD        | 32.9 | 210        | ug/L  |
| 85-01-8    | Phenanthrene               | 100   | UD        | 10.4 | 100        | ug/L  |
| 120-12-7   | Anthracene                 | 100   | UD        | 12.7 | 100        | ug/L  |
| 86-74-8    | Carbazole                  | 100   | UD        | 15.0 | 100        | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 100   | UD        | 25.4 | 100        | ug/L  |
| 206-44-0   | Fluoranthene               | 100   | UD        | 17.1 | 100        | ug/L  |
| 129-00-0   | Pyrene                     | 100   | UD        | 10.4 | 100        | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 100   | UD        | 40.2 | 100        | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 210   | UD        | 19.4 | 210        | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 100   | UD        | 9.40 | 100        | ug/L  |
| 218-01-9   | Chrysene                   | 100   | UD        | 9.20 | 100        | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 100   | UD        | 33.3 | 100        | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 210   | UD        | 48.8 | 210        | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 100   | UD        | 10.2 | 100        | ug/L  |

## Report of Analysis

|                    |                                    |                 |                  |
|--------------------|------------------------------------|-----------------|------------------|
| Client:            | AECOM                              | Date Collected: | 08/20/25         |
| Project:           | National Grid Equity - Brooklyn NY | Date Received:  | 08/20/25         |
| Client Sample ID:  | MW-2B-082025DL                     | SDG No.:        | Q2924            |
| Lab Sample ID:     | Q2924-05DL                         | Matrix:         | Water            |
| Analytical Method: | 8270E                              | % Solid:        | 0                |
| Sample Wt/Vol:     | 960 Units: mL                      | Final Vol:      | 1000 uL          |
| Soil Aliquot Vol:  | uL                                 | Test:           | SVOC-TCL BNA -20 |
| Extraction Type :  | Decanted : N                       | Level :         | LOW              |
| Injection Volume : | GPC Factor : 1.0                   | GPC Cleanup :   | N PH :           |
| Prep Method :      |                                    |                 |                  |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025583.D        | 20        | 08/22/25 10:34 | 08/26/25 13:10 | PB169362      |

| CAS Number                | Parameter                  | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------|----------------------------|---------|-----------|----------|------------|----------|
| 207-08-9                  | Benzo(k)fluoranthene       | 100     | UD        | 10.0     | 100        | ug/L     |
| 50-32-8                   | Benzo(a)pyrene             | 100     | UD        | 11.5     | 100        | ug/L     |
| 193-39-5                  | Indeno(1,2,3-cd)pyrene     | 100     | UD        | 12.3     | 100        | ug/L     |
| 53-70-3                   | Dibenzo(a,h)anthracene     | 100     | UD        | 14.0     | 100        | ug/L     |
| 191-24-2                  | Benzo(g,h,i)perylene       | 100     | UD        | 14.4     | 100        | ug/L     |
| 95-94-3                   | 1,2,4,5-Tetrachlorobenzene | 100     | UD        | 10.8     | 100        | ug/L     |
| 123-91-1                  | 1,4-Dioxane                | 100     | UD        | 20.8     | 100        | ug/L     |
| 58-90-2                   | 2,3,4,6-Tetrachlorophenol  | 100     | UD        | 15.0     | 100        | ug/L     |
| <b>SURROGATES</b>         |                            |         |           |          |            |          |
| 367-12-4                  | 2-Fluorophenol             | 80.0    |           | 23 - 138 | 53%        | SPK: 150 |
| 13127-88-3                | Phenol-d6                  | 51.4    |           | 10 - 134 | 34%        | SPK: 150 |
| 4165-60-0                 | Nitrobenzene-d5            | 85.6    |           | 67 - 132 | 86%        | SPK: 100 |
| 321-60-8                  | 2-Fluorobiphenyl           | 90.6    |           | 52 - 132 | 91%        | SPK: 100 |
| 118-79-6                  | 2,4,6-Tribromophenol       | 150     |           | 44 - 137 | 100%       | SPK: 150 |
| 1718-51-0                 | Terphenyl-d14              | 93.2    |           | 42 - 152 | 93%        | SPK: 100 |
| <b>INTERNAL STANDARDS</b> |                            |         |           |          |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4     | 167000  | 7.772     |          |            |          |
| 1146-65-2                 | Naphthalene-d8             | 645000  | 10.549    |          |            |          |
| 15067-26-2                | Acenaphthene-d10           | 404000  | 14.39     |          |            |          |
| 1517-22-2                 | Phenanthrene-d10           | 820000  | 17.189    |          |            |          |
| 1719-03-5                 | Chrysene-d12               | 923000  | 21.63     |          |            |          |
| 1520-96-3                 | Perylene-d12               | 1100000 | 24.995    |          |            |          |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-2B-082025DL2                    |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-05DL2                        |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 960                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025586.D        | 40        | 08/22/25 10:34 | 08/26/25 15:14 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 100-52-7       | Benzaldehyde                | 420   | UD        | 160  | 420        | ug/L  |
| 108-95-2       | Phenol                      | 210   | UD        | 37.9 | 210        | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 210   | UD        | 33.8 | 210        | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 210   | UD        | 24.2 | 210        | ug/L  |
| 95-48-7        | 2-Methylphenol              | 210   | UD        | 46.7 | 210        | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 210   | UD        | 53.3 | 210        | ug/L  |
| 98-86-2        | Acetophenone                | 210   | UD        | 30.8 | 210        | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 420   | UD        | 45.8 | 420        | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 100   | UD        | 58.8 | 100        | ug/L  |
| 67-72-1        | Hexachloroethane            | 210   | UD        | 27.1 | 210        | ug/L  |
| 98-95-3        | Nitrobenzene                | 210   | UD        | 31.7 | 210        | ug/L  |
| 78-59-1        | Isophorone                  | 210   | UD        | 31.3 | 210        | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 210   | UD        | 73.3 | 210        | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 210   | UD        | 77.1 | 210        | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 210   | UD        | 28.3 | 210        | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 210   | UD        | 21.7 | 210        | ug/L  |
| 91-20-3        | Naphthalene                 | 1800  | D         | 20.8 | 210        | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 210   | UD        | 35.0 | 210        | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 210   | UD        | 22.5 | 210        | ug/L  |
| 105-60-2       | Caprolactam                 | 420   | UD        | 47.1 | 420        | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 210   | UD        | 24.6 | 210        | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 180   | JD        | 23.3 | 210        | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 420   | UD        | 150  | 420        | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 210   | UD        | 21.3 | 210        | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 210   | UD        | 25.8 | 210        | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 210   | UD        | 22.1 | 210        | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 210   | UD        | 25.4 | 210        | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 210   | UD        | 52.5 | 210        | ug/L  |
| 131-11-3       | Dimethylphthalate           | 210   | UD        | 25.4 | 210        | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-2B-082025DL2                    |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-05DL2                        |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 960                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025586.D        | 40        | 08/22/25 10:34 | 08/26/25 15:14 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 210   | UD        | 31.3 | 210        | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 210   | UD        | 38.3 | 210        | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 210   | UD        | 43.8 | 210        | ug/L  |
| 83-32-9    | Acenaphthene               | 210   | UD        | 22.9 | 210        | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 420   | UD        | 250  | 420        | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 420   | UD        | 99.2 | 420        | ug/L  |
| 132-64-9   | Dibenzofuran               | 210   | UD        | 25.4 | 210        | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 210   | UD        | 50.8 | 210        | ug/L  |
| 84-66-2    | Diethylphthalate           | 210   | UD        | 28.8 | 210        | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 210   | UD        | 28.3 | 210        | ug/L  |
| 86-73-7    | Fluorene                   | 210   | UD        | 26.3 | 210        | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 210   | UD        | 62.5 | 210        | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 420   | UD        | 120  | 420        | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 210   | UD        | 24.2 | 210        | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 210   | UD        | 16.7 | 210        | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 210   | UD        | 21.7 | 210        | ug/L  |
| 1912-24-9  | Atrazine                   | 210   | UD        | 42.1 | 210        | ug/L  |
| 87-86-5    | Pentachlorophenol          | 420   | UD        | 65.8 | 420        | ug/L  |
| 85-01-8    | Phenanthrene               | 210   | UD        | 20.8 | 210        | ug/L  |
| 120-12-7   | Anthracene                 | 210   | UD        | 25.4 | 210        | ug/L  |
| 86-74-8    | Carbazole                  | 210   | UD        | 30.0 | 210        | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 210   | UD        | 50.8 | 210        | ug/L  |
| 206-44-0   | Fluoranthene               | 210   | UD        | 34.2 | 210        | ug/L  |
| 129-00-0   | Pyrene                     | 210   | UD        | 20.8 | 210        | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 210   | UD        | 80.4 | 210        | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 420   | UD        | 38.8 | 420        | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 210   | UD        | 18.8 | 210        | ug/L  |
| 218-01-9   | Chrysene                   | 210   | UD        | 18.3 | 210        | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 210   | UD        | 66.7 | 210        | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 420   | UD        | 97.5 | 420        | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 210   | UD        | 20.4 | 210        | ug/L  |

## Report of Analysis

|                    |                                    |                 |                  |
|--------------------|------------------------------------|-----------------|------------------|
| Client:            | AECOM                              | Date Collected: | 08/20/25         |
| Project:           | National Grid Equity - Brooklyn NY | Date Received:  | 08/20/25         |
| Client Sample ID:  | MW-2B-082025DL2                    | SDG No.:        | Q2924            |
| Lab Sample ID:     | Q2924-05DL2                        | Matrix:         | Water            |
| Analytical Method: | 8270E                              | % Solid:        | 0                |
| Sample Wt/Vol:     | 960 Units: mL                      | Final Vol:      | 1000 uL          |
| Soil Aliquot Vol:  | uL                                 | Test:           | SVOC-TCL BNA -20 |
| Extraction Type :  | Decanted : N                       | Level :         | LOW              |
| Injection Volume : | GPC Factor : 1.0                   | GPC Cleanup :   | N PH :           |
| Prep Method :      |                                    |                 |                  |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025586.D        | 40        | 08/22/25 10:34 | 08/26/25 15:14 | PB169362      |

| CAS Number                | Parameter                  | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------|----------------------------|---------|-----------|----------|------------|----------|
| 207-08-9                  | Benzo(k)fluoranthene       | 210     | UD        | 20.0     | 210        | ug/L     |
| 50-32-8                   | Benzo(a)pyrene             | 210     | UD        | 22.9     | 210        | ug/L     |
| 193-39-5                  | Indeno(1,2,3-cd)pyrene     | 210     | UD        | 24.6     | 210        | ug/L     |
| 53-70-3                   | Dibenzo(a,h)anthracene     | 210     | UD        | 27.9     | 210        | ug/L     |
| 191-24-2                  | Benzo(g,h,i)perylene       | 210     | UD        | 28.8     | 210        | ug/L     |
| 95-94-3                   | 1,2,4,5-Tetrachlorobenzene | 210     | UD        | 21.7     | 210        | ug/L     |
| 123-91-1                  | 1,4-Dioxane                | 210     | UD        | 41.7     | 210        | ug/L     |
| 58-90-2                   | 2,3,4,6-Tetrachlorophenol  | 210     | UD        | 30.0     | 210        | ug/L     |
| <b>SURROGATES</b>         |                            |         |           |          |            |          |
| 367-12-4                  | 2-Fluorophenol             | 79.8    |           | 23 - 138 | 53%        | SPK: 150 |
| 13127-88-3                | Phenol-d6                  | 48.0    |           | 10 - 134 | 32%        | SPK: 150 |
| 4165-60-0                 | Nitrobenzene-d5            | 80.4    |           | 67 - 132 | 80%        | SPK: 100 |
| 321-60-8                  | 2-Fluorobiphenyl           | 91.2    |           | 52 - 132 | 91%        | SPK: 100 |
| 118-79-6                  | 2,4,6-Tribromophenol       | 139     |           | 44 - 137 | 93%        | SPK: 150 |
| 1718-51-0                 | Terphenyl-d14              | 101     |           | 42 - 152 | 101%       | SPK: 100 |
| <b>INTERNAL STANDARDS</b> |                            |         |           |          |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4     | 173000  | 7.778     |          |            |          |
| 1146-65-2                 | Naphthalene-d8             | 701000  | 10.543    |          |            |          |
| 15067-26-2                | Acenaphthene-d10           | 444000  | 14.389    |          |            |          |
| 1517-22-2                 | Phenanthrene-d10           | 857000  | 17.189    |          |            |          |
| 1719-03-5                 | Chrysene-d12               | 880000  | 21.63     |          |            |          |
| 1520-96-3                 | Perylene-d12               | 1060000 | 24.989    |          |            |          |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-2A-082025                       |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-06                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 1000                               | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025590.D        | 1         | 08/22/25 10:34 | 08/26/25 17:59 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 100-52-7       | Benzaldehyde                | 10.0  | U         | 3.90 | 10.0       | ug/L  |
| 108-95-2       | Phenol                      | 5.00  | U         | 0.91 | 5.00       | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 5.00  | U         | 0.81 | 5.00       | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 5.00  | U         | 0.58 | 5.00       | ug/L  |
| 95-48-7        | 2-Methylphenol              | 5.00  | U         | 1.10 | 5.00       | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 5.00  | U         | 1.30 | 5.00       | ug/L  |
| 98-86-2        | Acetophenone                | 5.00  | U         | 0.74 | 5.00       | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 10.0  | U         | 1.10 | 10.0       | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 2.50  | U         | 1.40 | 2.50       | ug/L  |
| 67-72-1        | Hexachloroethane            | 5.00  | U         | 0.65 | 5.00       | ug/L  |
| 98-95-3        | Nitrobenzene                | 5.00  | U         | 0.76 | 5.00       | ug/L  |
| 78-59-1        | Isophorone                  | 5.00  | U         | 0.75 | 5.00       | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 5.00  | U         | 1.80 | 5.00       | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 5.00  | U         | 1.90 | 5.00       | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 5.00  | U         | 0.68 | 5.00       | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 5.00  | U         | 0.52 | 5.00       | ug/L  |
| 91-20-3        | Naphthalene                 | 5.00  | U         | 0.50 | 5.00       | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 5.00  | U         | 0.84 | 5.00       | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 5.00  | U         | 0.54 | 5.00       | ug/L  |
| 105-60-2       | Caprolactam                 | 10.0  | U         | 1.10 | 10.0       | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 5.00  | U         | 0.59 | 5.00       | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 5.00  | U         | 0.56 | 5.00       | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 10.0  | U         | 3.60 | 10.0       | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 5.00  | U         | 0.51 | 5.00       | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 5.00  | U         | 0.62 | 5.00       | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 5.00  | U         | 0.53 | 5.00       | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 5.00  | U         | 0.61 | 5.00       | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 5.00  | U         | 1.30 | 5.00       | ug/L  |
| 131-11-3       | Dimethylphthalate           | 5.00  | U         | 0.61 | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-2A-082025                       |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-06                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 1000                               | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025590.D        | 1         | 08/22/25 10:34 | 08/26/25 17:59 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 5.00  | U         | 0.75 | 5.00       | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 5.00  | U         | 0.92 | 5.00       | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 5.00  | U         | 1.10 | 5.00       | ug/L  |
| 83-32-9    | Acenaphthene               | 5.00  | U         | 0.55 | 5.00       | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 10.0  | U         | 6.00 | 10.0       | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 10.0  | U         | 2.40 | 10.0       | ug/L  |
| 132-64-9   | Dibenzofuran               | 5.00  | U         | 0.61 | 5.00       | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 5.00  | U         | 1.20 | 5.00       | ug/L  |
| 84-66-2    | Diethylphthalate           | 5.00  | U         | 0.69 | 5.00       | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 5.00  | U         | 0.68 | 5.00       | ug/L  |
| 86-73-7    | Fluorene                   | 5.00  | U         | 0.63 | 5.00       | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 5.00  | U         | 1.50 | 5.00       | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 10.0  | U         | 2.90 | 10.0       | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 5.00  | U         | 0.58 | 5.00       | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 5.00  | U         | 0.40 | 5.00       | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 5.00  | U         | 0.52 | 5.00       | ug/L  |
| 1912-24-9  | Atrazine                   | 5.00  | U         | 1.00 | 5.00       | ug/L  |
| 87-86-5    | Pentachlorophenol          | 10.0  | U         | 1.60 | 10.0       | ug/L  |
| 85-01-8    | Phenanthrene               | 5.00  | U         | 0.50 | 5.00       | ug/L  |
| 120-12-7   | Anthracene                 | 5.00  | U         | 0.61 | 5.00       | ug/L  |
| 86-74-8    | Carbazole                  | 5.00  | U         | 0.72 | 5.00       | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 5.00  | U         | 1.20 | 5.00       | ug/L  |
| 206-44-0   | Fluoranthene               | 5.00  | U         | 0.82 | 5.00       | ug/L  |
| 129-00-0   | Pyrene                     | 5.00  | U         | 0.50 | 5.00       | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 5.00  | U         | 1.90 | 5.00       | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 10.0  | U         | 0.93 | 10.0       | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 5.00  | U         | 0.45 | 5.00       | ug/L  |
| 218-01-9   | Chrysene                   | 5.00  | U         | 0.44 | 5.00       | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 5.00  | U         | 1.60 | 5.00       | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 10.0  | U         | 2.30 | 10.0       | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 5.00  | U         | 0.49 | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-2A-082025                       |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-06                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 1000                               | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025590.D        | 1         | 08/22/25 10:34 | 08/26/25 17:59 | PB169362      |

| CAS Number                            | Parameter                          | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------------------|------------------------------------|---------|-----------|----------|------------|----------|
| 207-08-9                              | Benzo(k)fluoranthene               | 5.00    | U         | 0.48     | 5.00       | ug/L     |
| 50-32-8                               | Benzo(a)pyrene                     | 5.00    | U         | 0.55     | 5.00       | ug/L     |
| 193-39-5                              | Indeno(1,2,3-cd)pyrene             | 5.00    | U         | 0.59     | 5.00       | ug/L     |
| 53-70-3                               | Dibenzo(a,h)anthracene             | 5.00    | U         | 0.67     | 5.00       | ug/L     |
| 191-24-2                              | Benzo(g,h,i)perylene               | 5.00    | U         | 0.69     | 5.00       | ug/L     |
| 95-94-3                               | 1,2,4,5-Tetrachlorobenzene         | 5.00    | U         | 0.52     | 5.00       | ug/L     |
| 123-91-1                              | 1,4-Dioxane                        | 5.00    | U         | 1.00     | 5.00       | ug/L     |
| 58-90-2                               | 2,3,4,6-Tetrachlorophenol          | 5.00    | U         | 0.72     | 5.00       | ug/L     |
| <b>SURROGATES</b>                     |                                    |         |           |          |            |          |
| 367-12-4                              | 2-Fluorophenol                     | 64.6    |           | 23 - 138 | 43%        | SPK: 150 |
| 13127-88-3                            | Phenol-d6                          | 39.1    |           | 10 - 134 | 26%        | SPK: 150 |
| 4165-60-0                             | Nitrobenzene-d5                    | 83.4    |           | 67 - 132 | 83%        | SPK: 100 |
| 321-60-8                              | 2-Fluorobiphenyl                   | 78.0    |           | 52 - 132 | 78%        | SPK: 100 |
| 118-79-6                              | 2,4,6-Tribromophenol               | 141     |           | 44 - 137 | 94%        | SPK: 150 |
| 1718-51-0                             | Terphenyl-d14                      | 77.0    |           | 42 - 152 | 77%        | SPK: 100 |
| <b>INTERNAL STANDARDS</b>             |                                    |         |           |          |            |          |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4             | 209000  | 7.778     |          |            |          |
| 1146-65-2                             | Naphthalene-d8                     | 834000  | 10.543    |          |            |          |
| 15067-26-2                            | Acenaphthene-d10                   | 541000  | 14.384    |          |            |          |
| 1517-22-2                             | Phenanthrene-d10                   | 1090000 | 17.19     |          |            |          |
| 1719-03-5                             | Chrysene-d12                       | 1130000 | 21.642    |          |            |          |
| 1520-96-3                             | Perylene-d12                       | 1290000 | 25.001    |          |            |          |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                                    |         |           |          |            |          |
| 000994-05-8                           | Butane, 2-methoxy-2-methyl-        | 96.5    | J         |          | 3.04       | ug/L     |
| 000123-42-2                           | 2-Pentanone, 4-hydroxy-4-methyl-   | 4.90    | AB        |          | 4.94       | ug/L     |
| 001638-16-0                           | 2-Propanol, 1,1-[(1-methyl-1,2-et  | 4.20    | J         |          | 11.1       | ug/L     |
| 000106-62-7                           | 1-Propanol, 2-(2-hydroxypropoxy)-  | 3.70    | J         |          | 11.1       | ug/L     |
| 054966-46-0                           | Butanoic acid, 4-(2-methoxy-1-meth | 2.70    | J         |          | 11.3       | ug/L     |
|                                       | unknown12.196                      | 17.6    | J         |          | 12.2       | ug/L     |
| 004455-26-9                           | 1-Octanamine, N-methyl-N-octyl-    | 2.90    | J         |          | 16.4       | ug/L     |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-2A-082025                       |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-06                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 1000                               | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025590.D        | 1         | 08/22/25 10:34 | 08/26/25 17:59 | PB169362      |

| CAS Number  | Parameter                     | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|-------------|-------------------------------|-------|-----------|-----|------------|-------|
| 004606-07-9 | Ethyl cyclopropanecarboxylate | 17.2  | J         |     | 16.7       | ug/L  |
| 019437-26-4 | di-2-Pyridyl ketone           | 5.00  | J         |     | 17.9       | ug/L  |
| 000057-10-3 | n-Hexadecanoic acid           | 4.30  | J         |     | 18.1       | ug/L  |
|             | unknown19.683                 | 4.80  | J         |     | 19.7       | ug/L  |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-18C-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-07                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 990                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BF143603.D        | 1         | 08/22/25 10:34 | 08/27/25 12:52 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 100-52-7       | Benzaldehyde                | 10.1  | U         | 3.90 | 10.1       | ug/L  |
| 108-95-2       | Phenol                      | 5.10  | U         | 0.92 | 5.10       | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 5.10  | U         | 0.82 | 5.10       | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 5.10  | U         | 0.59 | 5.10       | ug/L  |
| 95-48-7        | 2-Methylphenol              | 5.10  | U         | 1.10 | 5.10       | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 5.10  | U         | 1.30 | 5.10       | ug/L  |
| 98-86-2        | Acetophenone                | 5.10  | U         | 0.75 | 5.10       | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 10.1  | U         | 1.10 | 10.1       | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 2.50  | U         | 1.40 | 2.50       | ug/L  |
| 67-72-1        | Hexachloroethane            | 5.10  | U         | 0.66 | 5.10       | ug/L  |
| 98-95-3        | Nitrobenzene                | 5.10  | U         | 0.77 | 5.10       | ug/L  |
| 78-59-1        | Isophorone                  | 5.10  | U         | 0.76 | 5.10       | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 5.10  | U         | 1.80 | 5.10       | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 5.10  | U         | 1.90 | 5.10       | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 5.10  | U         | 0.69 | 5.10       | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 5.10  | U         | 0.53 | 5.10       | ug/L  |
| 91-20-3        | Naphthalene                 | 5.10  | U         | 0.51 | 5.10       | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 5.10  | U         | 0.85 | 5.10       | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 5.10  | U         | 0.55 | 5.10       | ug/L  |
| 105-60-2       | Caprolactam                 | 10.1  | U         | 1.10 | 10.1       | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 5.10  | U         | 0.60 | 5.10       | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 5.10  | U         | 0.57 | 5.10       | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 10.1  | U         | 3.70 | 10.1       | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 5.10  | U         | 0.52 | 5.10       | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 5.10  | U         | 0.63 | 5.10       | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 5.10  | U         | 0.54 | 5.10       | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 5.10  | U         | 0.62 | 5.10       | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 5.10  | U         | 1.30 | 5.10       | ug/L  |
| 131-11-3       | Dimethylphthalate           | 5.10  | U         | 0.62 | 5.10       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-18C-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-07                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 990                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BF143603.D        | 1         | 08/22/25 10:34 | 08/27/25 12:52 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 5.10  | U         | 0.76 | 5.10       | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 5.10  | U         | 0.93 | 5.10       | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 5.10  | U         | 1.10 | 5.10       | ug/L  |
| 83-32-9    | Acenaphthene               | 5.10  | U         | 0.56 | 5.10       | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 10.1  | U         | 6.00 | 10.1       | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 10.1  | U         | 2.40 | 10.1       | ug/L  |
| 132-64-9   | Dibenzofuran               | 5.10  | U         | 0.62 | 5.10       | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 5.10  | U         | 1.20 | 5.10       | ug/L  |
| 84-66-2    | Diethylphthalate           | 5.10  | U         | 0.70 | 5.10       | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 5.10  | U         | 0.69 | 5.10       | ug/L  |
| 86-73-7    | Fluorene                   | 5.10  | U         | 0.64 | 5.10       | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 5.10  | U         | 1.50 | 5.10       | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 10.1  | U         | 2.90 | 10.1       | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 5.10  | U         | 0.59 | 5.10       | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 5.10  | U         | 0.40 | 5.10       | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 5.10  | U         | 0.53 | 5.10       | ug/L  |
| 1912-24-9  | Atrazine                   | 5.10  | U         | 1.00 | 5.10       | ug/L  |
| 87-86-5    | Pentachlorophenol          | 10.1  | U         | 1.60 | 10.1       | ug/L  |
| 85-01-8    | Phenanthrene               | 5.10  | U         | 0.51 | 5.10       | ug/L  |
| 120-12-7   | Anthracene                 | 5.10  | U         | 0.62 | 5.10       | ug/L  |
| 86-74-8    | Carbazole                  | 5.10  | U         | 0.73 | 5.10       | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 5.10  | U         | 1.20 | 5.10       | ug/L  |
| 206-44-0   | Fluoranthene               | 5.10  | U         | 0.83 | 5.10       | ug/L  |
| 129-00-0   | Pyrene                     | 5.10  | U         | 0.51 | 5.10       | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 5.10  | U         | 1.90 | 5.10       | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 10.1  | U         | 0.94 | 10.1       | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 5.10  | U         | 0.45 | 5.10       | ug/L  |
| 218-01-9   | Chrysene                   | 5.10  | U         | 0.44 | 5.10       | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 5.10  | U         | 1.60 | 5.10       | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 10.1  | U         | 2.40 | 10.1       | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 5.10  | U         | 0.49 | 5.10       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-18C-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-07                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 990                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BF143603.D        | 1         | 08/22/25 10:34 | 08/27/25 12:52 | PB169362      |

| CAS Number                            | Parameter                        | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------------------|----------------------------------|--------|-----------|----------|------------|----------|
| 207-08-9                              | Benzo(k)fluoranthene             | 5.10   | U         | 0.48     | 5.10       | ug/L     |
| 50-32-8                               | Benzo(a)pyrene                   | 5.10   | U         | 0.56     | 5.10       | ug/L     |
| 193-39-5                              | Indeno(1,2,3-cd)pyrene           | 5.10   | U         | 0.60     | 5.10       | ug/L     |
| 53-70-3                               | Dibenzo(a,h)anthracene           | 5.10   | U         | 0.68     | 5.10       | ug/L     |
| 191-24-2                              | Benzo(g,h,i)perylene             | 5.10   | U         | 0.70     | 5.10       | ug/L     |
| 95-94-3                               | 1,2,4,5-Tetrachlorobenzene       | 5.10   | U         | 0.53     | 5.10       | ug/L     |
| 123-91-1                              | 1,4-Dioxane                      | 5.10   | U         | 1.00     | 5.10       | ug/L     |
| 58-90-2                               | 2,3,4,6-Tetrachlorophenol        | 5.10   | U         | 0.73     | 5.10       | ug/L     |
| <b>SURROGATES</b>                     |                                  |        |           |          |            |          |
| 367-12-4                              | 2-Fluorophenol                   | 67.3   |           | 23 - 138 | 45%        | SPK: 150 |
| 13127-88-3                            | Phenol-d6                        | 41.9   |           | 10 - 134 | 28%        | SPK: 150 |
| 4165-60-0                             | Nitrobenzene-d5                  | 113    |           | 67 - 132 | 113%       | SPK: 100 |
| 321-60-8                              | 2-Fluorobiphenyl                 | 95.1   |           | 52 - 132 | 95%        | SPK: 100 |
| 118-79-6                              | 2,4,6-Tribromophenol             | 142    |           | 44 - 137 | 95%        | SPK: 150 |
| 1718-51-0                             | Terphenyl-d14                    | 96.9   |           | 42 - 152 | 97%        | SPK: 100 |
| <b>INTERNAL STANDARDS</b>             |                                  |        |           |          |            |          |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4           | 107000 | 6.892     |          |            |          |
| 1146-65-2                             | Naphthalene-d8                   | 403000 | 8.169     |          |            |          |
| 15067-26-2                            | Acenaphthene-d10                 | 213000 | 9.922     |          |            |          |
| 1517-22-2                             | Phenanthrene-d10                 | 298000 | 11.41     |          |            |          |
| 1719-03-5                             | Chrysene-d12                     | 161000 | 14.045    |          |            |          |
| 1520-96-3                             | Perylene-d12                     | 219000 | 15.521    |          |            |          |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                                  |        |           |          |            |          |
| 000123-42-2                           | 2-Pentanone, 4-hydroxy-4-methyl- | 5.80   | AB        |          | 5.11       | ug/L     |
| 000119-61-9                           | Benzophenone                     | 2.30   | J         |          | 10.6       | ug/L     |
| 000057-10-3                           | n-Hexadecanoic acid              | 2.70   | J         |          | 11.9       | ug/L     |
| 000373-49-9                           | Palmitoleic acid                 | 2.30   | J         |          | 12.6       | ug/L     |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-18C-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-07                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 990                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BF143603.D        | 1         | 08/22/25 10:34 | 08/27/25 12:52 | PB169362      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-16A-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-08                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 910                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025588.D        | 1         | 08/22/25 10:34 | 08/26/25 16:37 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 100-52-7       | Benzaldehyde                | 11.0  | U         | 4.30 | 11.0       | ug/L  |
| 108-95-2       | Phenol                      | 5.50  | U         | 1.00 | 5.50       | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 5.50  | U         | 0.89 | 5.50       | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 5.50  | U         | 0.64 | 5.50       | ug/L  |
| 95-48-7        | 2-Methylphenol              | 5.50  | U         | 1.20 | 5.50       | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 5.50  | U         | 1.40 | 5.50       | ug/L  |
| 98-86-2        | Acetophenone                | 5.50  | U         | 0.81 | 5.50       | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 11.0  | U         | 1.20 | 11.0       | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 2.70  | U         | 1.50 | 2.70       | ug/L  |
| 67-72-1        | Hexachloroethane            | 5.50  | U         | 0.71 | 5.50       | ug/L  |
| 98-95-3        | Nitrobenzene                | 5.50  | U         | 0.84 | 5.50       | ug/L  |
| 78-59-1        | Isophorone                  | 5.50  | U         | 0.82 | 5.50       | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 5.50  | U         | 1.90 | 5.50       | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 5.50  | U         | 2.00 | 5.50       | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 5.50  | U         | 0.75 | 5.50       | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 5.50  | U         | 0.57 | 5.50       | ug/L  |
| 91-20-3        | Naphthalene                 | 5.50  | U         | 0.55 | 5.50       | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 5.50  | U         | 0.92 | 5.50       | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 5.50  | U         | 0.59 | 5.50       | ug/L  |
| 105-60-2       | Caprolactam                 | 11.0  | U         | 1.20 | 11.0       | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 5.50  | U         | 0.65 | 5.50       | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 5.50  | U         | 0.62 | 5.50       | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 11.0  | U         | 4.00 | 11.0       | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 5.50  | U         | 0.56 | 5.50       | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 5.50  | U         | 0.68 | 5.50       | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 5.50  | U         | 0.58 | 5.50       | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 5.50  | U         | 0.67 | 5.50       | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 5.50  | U         | 1.40 | 5.50       | ug/L  |
| 131-11-3       | Dimethylphthalate           | 5.50  | U         | 0.67 | 5.50       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-16A-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-08                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 910                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025588.D        | 1         | 08/22/25 10:34 | 08/26/25 16:37 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 5.50  | U         | 0.82 | 5.50       | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 5.50  | U         | 1.00 | 5.50       | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 5.50  | U         | 1.20 | 5.50       | ug/L  |
| 83-32-9    | Acenaphthene               | 5.50  | U         | 0.60 | 5.50       | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 11.0  | U         | 6.60 | 11.0       | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 11.0  | U         | 2.60 | 11.0       | ug/L  |
| 132-64-9   | Dibenzofuran               | 5.50  | U         | 0.67 | 5.50       | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 5.50  | U         | 1.30 | 5.50       | ug/L  |
| 84-66-2    | Diethylphthalate           | 5.50  | U         | 0.76 | 5.50       | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 5.50  | U         | 0.75 | 5.50       | ug/L  |
| 86-73-7    | Fluorene                   | 5.50  | U         | 0.69 | 5.50       | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 5.50  | U         | 1.60 | 5.50       | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 11.0  | U         | 3.20 | 11.0       | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 5.50  | U         | 0.64 | 5.50       | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 5.50  | U         | 0.44 | 5.50       | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 5.50  | U         | 0.57 | 5.50       | ug/L  |
| 1912-24-9  | Atrazine                   | 5.50  | U         | 1.10 | 5.50       | ug/L  |
| 87-86-5    | Pentachlorophenol          | 11.0  | U         | 1.70 | 11.0       | ug/L  |
| 85-01-8    | Phenanthrene               | 5.50  | U         | 0.55 | 5.50       | ug/L  |
| 120-12-7   | Anthracene                 | 5.50  | U         | 0.67 | 5.50       | ug/L  |
| 86-74-8    | Carbazole                  | 5.50  | U         | 0.79 | 5.50       | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 5.50  | U         | 1.30 | 5.50       | ug/L  |
| 206-44-0   | Fluoranthene               | 5.50  | U         | 0.90 | 5.50       | ug/L  |
| 129-00-0   | Pyrene                     | 5.50  | U         | 0.55 | 5.50       | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 5.50  | U         | 2.10 | 5.50       | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 11.0  | U         | 1.00 | 11.0       | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 5.50  | U         | 0.49 | 5.50       | ug/L  |
| 218-01-9   | Chrysene                   | 5.50  | U         | 0.48 | 5.50       | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 5.50  | U         | 1.80 | 5.50       | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 11.0  | U         | 2.60 | 11.0       | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 5.50  | U         | 0.54 | 5.50       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-16A-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-08                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 910                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025588.D        | 1         | 08/22/25 10:34 | 08/26/25 16:37 | PB169362      |

| CAS Number                            | Parameter                        | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------------------|----------------------------------|---------|-----------|----------|------------|----------|
| 207-08-9                              | Benzo(k)fluoranthene             | 5.50    | U         | 0.53     | 5.50       | ug/L     |
| 50-32-8                               | Benzo(a)pyrene                   | 5.50    | U         | 0.60     | 5.50       | ug/L     |
| 193-39-5                              | Indeno(1,2,3-cd)pyrene           | 5.50    | U         | 0.65     | 5.50       | ug/L     |
| 53-70-3                               | Dibenzo(a,h)anthracene           | 5.50    | U         | 0.74     | 5.50       | ug/L     |
| 191-24-2                              | Benzo(g,h,i)perylene             | 5.50    | U         | 0.76     | 5.50       | ug/L     |
| 95-94-3                               | 1,2,4,5-Tetrachlorobenzene       | 5.50    | U         | 0.57     | 5.50       | ug/L     |
| 123-91-1                              | 1,4-Dioxane                      | 5.50    | U         | 1.10     | 5.50       | ug/L     |
| 58-90-2                               | 2,3,4,6-Tetrachlorophenol        | 5.50    | U         | 0.79     | 5.50       | ug/L     |
| <b>SURROGATES</b>                     |                                  |         |           |          |            |          |
| 367-12-4                              | 2-Fluorophenol                   | 59.0    |           | 23 - 138 | 39%        | SPK: 150 |
| 13127-88-3                            | Phenol-d6                        | 34.9    |           | 10 - 134 | 23%        | SPK: 150 |
| 4165-60-0                             | Nitrobenzene-d5                  | 84.0    |           | 67 - 132 | 84%        | SPK: 100 |
| 321-60-8                              | 2-Fluorobiphenyl                 | 77.7    |           | 52 - 132 | 78%        | SPK: 100 |
| 118-79-6                              | 2,4,6-Tribromophenol             | 145     |           | 44 - 137 | 97%        | SPK: 150 |
| 1718-51-0                             | Terphenyl-d14                    | 72.4    |           | 42 - 152 | 72%        | SPK: 100 |
| <b>INTERNAL STANDARDS</b>             |                                  |         |           |          |            |          |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4           | 195000  | 7.778     |          |            |          |
| 1146-65-2                             | Naphthalene-d8                   | 751000  | 10.549    |          |            |          |
| 15067-26-2                            | Acenaphthene-d10                 | 492000  | 14.389    |          |            |          |
| 1517-22-2                             | Phenanthrene-d10                 | 1020000 | 17.195    |          |            |          |
| 1719-03-5                             | Chrysene-d12                     | 1120000 | 21.642    |          |            |          |
| 1520-96-3                             | Perylene-d12                     | 1310000 | 25.018    |          |            |          |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                                  |         |           |          |            |          |
| 000994-05-8                           | Butane, 2-methoxy-2-methyl-      | 110     | J         |          | 3.03       | ug/L     |
| 000123-42-2                           | 2-Pentanone, 4-hydroxy-4-methyl- | 5.60    | AB        |          | 4.94       | ug/L     |
| 000057-10-3                           | n-Hexadecanoic acid              | 5.20    | J         |          | 18.1       | ug/L     |
| 013126-39-1                           | cis-13-Octadecenoic acid         | 7.30    | J         |          | 19.3       | ug/L     |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-16A-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-08                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 910                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025588.D        | 1         | 08/22/25 10:34 | 08/26/25 16:37 | PB169362      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-16C-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-09                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 940                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025589.D        | 1         | 08/22/25 10:34 | 08/26/25 17:18 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 100-52-7       | Benzaldehyde                | 10.6  | U         | 4.20 | 10.6       | ug/L  |
| 108-95-2       | Phenol                      | 5.30  | U         | 0.97 | 5.30       | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 5.30  | U         | 0.86 | 5.30       | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 5.30  | U         | 0.62 | 5.30       | ug/L  |
| 95-48-7        | 2-Methylphenol              | 5.30  | U         | 1.20 | 5.30       | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 5.30  | U         | 1.40 | 5.30       | ug/L  |
| 98-86-2        | Acetophenone                | 5.30  | U         | 0.79 | 5.30       | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 10.6  | U         | 1.20 | 10.6       | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 2.70  | U         | 1.50 | 2.70       | ug/L  |
| 67-72-1        | Hexachloroethane            | 5.30  | U         | 0.69 | 5.30       | ug/L  |
| 98-95-3        | Nitrobenzene                | 5.30  | U         | 0.81 | 5.30       | ug/L  |
| 78-59-1        | Isophorone                  | 5.30  | U         | 0.80 | 5.30       | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 5.30  | U         | 1.90 | 5.30       | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 5.30  | U         | 2.00 | 5.30       | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 5.30  | U         | 0.72 | 5.30       | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 5.30  | U         | 0.55 | 5.30       | ug/L  |
| 91-20-3        | Naphthalene                 | 5.30  | U         | 0.53 | 5.30       | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 5.30  | U         | 0.89 | 5.30       | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 5.30  | U         | 0.57 | 5.30       | ug/L  |
| 105-60-2       | Caprolactam                 | 10.6  | U         | 1.20 | 10.6       | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 5.30  | U         | 0.63 | 5.30       | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 5.30  | U         | 0.60 | 5.30       | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 10.6  | U         | 3.90 | 10.6       | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 5.30  | U         | 0.54 | 5.30       | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 5.30  | U         | 0.66 | 5.30       | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 5.30  | U         | 0.56 | 5.30       | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 5.30  | U         | 0.65 | 5.30       | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 5.30  | U         | 1.30 | 5.30       | ug/L  |
| 131-11-3       | Dimethylphthalate           | 5.30  | U         | 0.65 | 5.30       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-16C-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-09                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 940                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025589.D        | 1         | 08/22/25 10:34 | 08/26/25 17:18 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 5.30  | U         | 0.80 | 5.30       | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 5.30  | U         | 0.98 | 5.30       | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 5.30  | U         | 1.10 | 5.30       | ug/L  |
| 83-32-9    | Acenaphthene               | 5.30  | U         | 0.59 | 5.30       | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 10.6  | U         | 6.40 | 10.6       | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 10.6  | U         | 2.50 | 10.6       | ug/L  |
| 132-64-9   | Dibenzofuran               | 5.30  | U         | 0.65 | 5.30       | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 5.30  | U         | 1.30 | 5.30       | ug/L  |
| 84-66-2    | Diethylphthalate           | 5.30  | U         | 0.73 | 5.30       | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 5.30  | U         | 0.72 | 5.30       | ug/L  |
| 86-73-7    | Fluorene                   | 5.30  | U         | 0.67 | 5.30       | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 5.30  | U         | 1.60 | 5.30       | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 10.6  | U         | 3.10 | 10.6       | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 5.30  | U         | 0.62 | 5.30       | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 5.30  | U         | 0.43 | 5.30       | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 5.30  | U         | 0.55 | 5.30       | ug/L  |
| 1912-24-9  | Atrazine                   | 5.30  | U         | 1.10 | 5.30       | ug/L  |
| 87-86-5    | Pentachlorophenol          | 10.6  | U         | 1.70 | 10.6       | ug/L  |
| 85-01-8    | Phenanthrene               | 5.30  | U         | 0.53 | 5.30       | ug/L  |
| 120-12-7   | Anthracene                 | 5.30  | U         | 0.65 | 5.30       | ug/L  |
| 86-74-8    | Carbazole                  | 5.30  | U         | 0.77 | 5.30       | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 5.30  | U         | 1.30 | 5.30       | ug/L  |
| 206-44-0   | Fluoranthene               | 5.30  | U         | 0.87 | 5.30       | ug/L  |
| 129-00-0   | Pyrene                     | 5.30  | U         | 0.53 | 5.30       | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 5.30  | U         | 2.10 | 5.30       | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 10.6  | U         | 0.99 | 10.6       | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 5.30  | U         | 0.48 | 5.30       | ug/L  |
| 218-01-9   | Chrysene                   | 5.30  | U         | 0.47 | 5.30       | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 5.30  | U         | 1.70 | 5.30       | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 10.6  | U         | 2.50 | 10.6       | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 5.30  | U         | 0.52 | 5.30       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-16C-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-09                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 940                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025589.D        | 1         | 08/22/25 10:34 | 08/26/25 17:18 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 207-08-9   | Benzo(k)fluoranthene       | 5.30  | U         | 0.51 | 5.30       | ug/L  |
| 50-32-8    | Benzo(a)pyrene             | 5.30  | U         | 0.59 | 5.30       | ug/L  |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 5.30  | U         | 0.63 | 5.30       | ug/L  |
| 53-70-3    | Dibenzo(a,h)anthracene     | 5.30  | U         | 0.71 | 5.30       | ug/L  |
| 191-24-2   | Benzo(g,h,i)perylene       | 5.30  | U         | 0.73 | 5.30       | ug/L  |
| 95-94-3    | 1,2,4,5-Tetrachlorobenzene | 5.30  | U         | 0.55 | 5.30       | ug/L  |
| 123-91-1   | 1,4-Dioxane                | 5.30  | U         | 1.10 | 5.30       | ug/L  |
| 58-90-2    | 2,3,4,6-Tetrachlorophenol  | 5.30  | U         | 0.77 | 5.30       | ug/L  |

### SURROGATES

|            |                      |      |  |          |     |          |
|------------|----------------------|------|--|----------|-----|----------|
| 367-12-4   | 2-Fluorophenol       | 64.9 |  | 23 - 138 | 43% | SPK: 150 |
| 13127-88-3 | Phenol-d6            | 39.6 |  | 10 - 134 | 26% | SPK: 150 |
| 4165-60-0  | Nitrobenzene-d5      | 83.7 |  | 67 - 132 | 84% | SPK: 100 |
| 321-60-8   | 2-Fluorobiphenyl     | 79.7 |  | 52 - 132 | 80% | SPK: 100 |
| 118-79-6   | 2,4,6-Tribromophenol | 141  |  | 44 - 137 | 94% | SPK: 150 |
| 1718-51-0  | Terphenyl-d14        | 76.3 |  | 42 - 152 | 76% | SPK: 100 |

### INTERNAL STANDARDS

|            |                        |         |        |
|------------|------------------------|---------|--------|
| 3855-82-1  | 1,4-Dichlorobenzene-d4 | 201000  | 7.772  |
| 1146-65-2  | Naphthalene-d8         | 791000  | 10.549 |
| 15067-26-2 | Acenaphthene-d10       | 502000  | 14.395 |
| 1517-22-2  | Phenanthrene-d10       | 1000000 | 17.195 |
| 1719-03-5  | Chrysene-d12           | 1070000 | 21.63  |
| 1520-96-3  | Perylene-d12           | 1230000 | 25.007 |

### TENTATIVE IDENTIFIED COMPOUNDS

|              |                                    |      |    |      |      |
|--------------|------------------------------------|------|----|------|------|
| 000994-05-8  | Butane, 2-methoxy-2-methyl-        | 110  | J  | 3.03 | ug/L |
| 000123-42-2  | 2-Pentanone, 4-hydroxy-4-methyl-   | 5.30 | AB | 4.94 | ug/L |
| 000617-94-7  | Benzenemethanol, .alpha.,.alpha.-d | 2.90 | J  | 8.87 | ug/L |
| 1010189-31-0 | 2,3,4,5,6,7-Hexahydro-1H-cyclopent | 2.60 | J  | 10.6 | ug/L |
| 215057-28-6  | 3-(2,3-Dihydro-1-benzofuran-5-yl)p | 3.90 | J  | 11.1 | ug/L |
| 006682-71-9  | 1H-Indene, 2,3-dihydro-4,7-dimethy | 2.90 | J  | 11.5 | ug/L |
| 000057-10-3  | n-Hexadecanoic acid                | 4.00 | J  | 18.1 | ug/L |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-16C-082025                      |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-09                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 940                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025589.D        | 1         | 08/22/25 10:34 | 08/26/25 17:18 | PB169362      |

| CAS Number  | Parameter   | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|-------------|-------------|-------|-----------|-----|------------|-------|
| 000112-86-7 | Erucic acid | 2.70  | J         |     | 19.3       | ug/L  |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-8A-082025                       |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-10                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 930                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025552.D        | 1         | 08/22/25 10:34 | 08/22/25 16:58 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 100-52-7       | Benzaldehyde                | 10.8  | U         | 4.20 | 10.8       | ug/L  |
| 108-95-2       | Phenol                      | 5.40  | U         | 0.98 | 5.40       | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 5.40  | U         | 0.87 | 5.40       | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 5.40  | U         | 0.62 | 5.40       | ug/L  |
| 95-48-7        | 2-Methylphenol              | 5.40  | U         | 1.20 | 5.40       | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 5.40  | U         | 1.40 | 5.40       | ug/L  |
| 98-86-2        | Acetophenone                | 5.40  | U         | 0.80 | 5.40       | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 10.8  | U         | 1.20 | 10.8       | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 2.70  | U         | 1.50 | 2.70       | ug/L  |
| 67-72-1        | Hexachloroethane            | 5.40  | U         | 0.70 | 5.40       | ug/L  |
| 98-95-3        | Nitrobenzene                | 5.40  | U         | 0.82 | 5.40       | ug/L  |
| 78-59-1        | Isophorone                  | 5.40  | U         | 0.81 | 5.40       | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 5.40  | U         | 1.90 | 5.40       | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 5.40  | U         | 2.00 | 5.40       | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 5.40  | U         | 0.73 | 5.40       | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 5.40  | U         | 0.56 | 5.40       | ug/L  |
| 91-20-3        | Naphthalene                 | 5.40  | U         | 0.54 | 5.40       | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 5.40  | U         | 0.90 | 5.40       | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 5.40  | U         | 0.58 | 5.40       | ug/L  |
| 105-60-2       | Caprolactam                 | 10.8  | U         | 1.20 | 10.8       | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 5.40  | U         | 0.63 | 5.40       | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 5.40  | U         | 0.60 | 5.40       | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 10.8  | U         | 3.90 | 10.8       | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 5.40  | U         | 0.55 | 5.40       | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 5.40  | U         | 0.67 | 5.40       | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 5.40  | U         | 0.57 | 5.40       | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 5.40  | U         | 0.66 | 5.40       | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 5.40  | U         | 1.40 | 5.40       | ug/L  |
| 131-11-3       | Dimethylphthalate           | 5.40  | U         | 0.66 | 5.40       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-8A-082025                       |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-10                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 930                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025552.D        | 1         | 08/22/25 10:34 | 08/22/25 16:58 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 5.40  | U         | 0.81 | 5.40       | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 5.40  | U         | 0.99 | 5.40       | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 5.40  | U         | 1.10 | 5.40       | ug/L  |
| 83-32-9    | Acenaphthene               | 5.40  | U         | 0.59 | 5.40       | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 10.8  | U         | 6.40 | 10.8       | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 10.8  | U         | 2.60 | 10.8       | ug/L  |
| 132-64-9   | Dibenzofuran               | 5.40  | U         | 0.66 | 5.40       | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 5.40  | U         | 1.30 | 5.40       | ug/L  |
| 84-66-2    | Diethylphthalate           | 5.40  | U         | 0.74 | 5.40       | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 5.40  | U         | 0.73 | 5.40       | ug/L  |
| 86-73-7    | Fluorene                   | 5.40  | U         | 0.68 | 5.40       | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 5.40  | U         | 1.60 | 5.40       | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 10.8  | U         | 3.10 | 10.8       | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 5.40  | U         | 0.62 | 5.40       | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 5.40  | U         | 0.43 | 5.40       | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 5.40  | U         | 0.56 | 5.40       | ug/L  |
| 1912-24-9  | Atrazine                   | 5.40  | U         | 1.10 | 5.40       | ug/L  |
| 87-86-5    | Pentachlorophenol          | 10.8  | U         | 1.70 | 10.8       | ug/L  |
| 85-01-8    | Phenanthrene               | 5.40  | U         | 0.54 | 5.40       | ug/L  |
| 120-12-7   | Anthracene                 | 5.40  | U         | 0.66 | 5.40       | ug/L  |
| 86-74-8    | Carbazole                  | 5.40  | U         | 0.77 | 5.40       | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 5.40  | U         | 1.30 | 5.40       | ug/L  |
| 206-44-0   | Fluoranthene               | 5.40  | U         | 0.88 | 5.40       | ug/L  |
| 129-00-0   | Pyrene                     | 5.40  | U         | 0.54 | 5.40       | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 5.40  | U         | 2.10 | 5.40       | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 10.8  | U         | 1.00 | 10.8       | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 5.40  | U         | 0.48 | 5.40       | ug/L  |
| 218-01-9   | Chrysene                   | 5.40  | U         | 0.47 | 5.40       | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 5.40  | U         | 1.70 | 5.40       | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 10.8  | U         | 2.50 | 10.8       | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 5.40  | U         | 0.53 | 5.40       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-8A-082025                       |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-10                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 930                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025552.D        | 1         | 08/22/25 10:34 | 08/22/25 16:58 | PB169362      |

| CAS Number                            | Parameter                        | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------------------|----------------------------------|---------|-----------|----------|------------|----------|
| 207-08-9                              | Benzo(k)fluoranthene             | 5.40    | U         | 0.52     | 5.40       | ug/L     |
| 50-32-8                               | Benzo(a)pyrene                   | 5.40    | U         | 0.59     | 5.40       | ug/L     |
| 193-39-5                              | Indeno(1,2,3-cd)pyrene           | 5.40    | U         | 0.63     | 5.40       | ug/L     |
| 53-70-3                               | Dibenzo(a,h)anthracene           | 5.40    | U         | 0.72     | 5.40       | ug/L     |
| 191-24-2                              | Benzo(g,h,i)perylene             | 5.40    | U         | 0.74     | 5.40       | ug/L     |
| 95-94-3                               | 1,2,4,5-Tetrachlorobenzene       | 5.40    | U         | 0.56     | 5.40       | ug/L     |
| 123-91-1                              | 1,4-Dioxane                      | 5.40    | U         | 1.10     | 5.40       | ug/L     |
| 58-90-2                               | 2,3,4,6-Tetrachlorophenol        | 5.40    | U         | 0.77     | 5.40       | ug/L     |
| <b>SURROGATES</b>                     |                                  |         |           |          |            |          |
| 367-12-4                              | 2-Fluorophenol                   | 64.1    |           | 23 - 138 | 43%        | SPK: 150 |
| 13127-88-3                            | Phenol-d6                        | 39.6    |           | 10 - 134 | 26%        | SPK: 150 |
| 4165-60-0                             | Nitrobenzene-d5                  | 84.7    |           | 67 - 132 | 85%        | SPK: 100 |
| 321-60-8                              | 2-Fluorobiphenyl                 | 77.8    |           | 52 - 132 | 78%        | SPK: 100 |
| 118-79-6                              | 2,4,6-Tribromophenol             | 138     |           | 44 - 137 | 92%        | SPK: 150 |
| 1718-51-0                             | Terphenyl-d14                    | 74.9    |           | 42 - 152 | 75%        | SPK: 100 |
| <b>INTERNAL STANDARDS</b>             |                                  |         |           |          |            |          |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4           | 178000  | 7.778     |          |            |          |
| 1146-65-2                             | Naphthalene-d8                   | 716000  | 10.543    |          |            |          |
| 15067-26-2                            | Acenaphthene-d10                 | 464000  | 14.384    |          |            |          |
| 1517-22-2                             | Phenanthrene-d10                 | 972000  | 17.183    |          |            |          |
| 1719-03-5                             | Chrysene-d12                     | 995000  | 21.63     |          |            |          |
| 1520-96-3                             | Perylene-d12                     | 1120000 | 24.989    |          |            |          |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                                  |         |           |          |            |          |
| 000994-05-8                           | Butane, 2-methoxy-2-methyl-      | 110     | J         |          | 3.03       | ug/L     |
| 000123-42-2                           | 2-Pentanone, 4-hydroxy-4-methyl- | 4.40    | AB        |          | 4.94       | ug/L     |
| 000119-61-9                           | Benzophenone                     | 2.80    | J         |          | 15.8       | ug/L     |
| 000057-10-3                           | n-Hexadecanoic acid              | 5.30    | J         |          | 18.1       | ug/L     |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-8A-082025                       |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-10                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 930                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025552.D        | 1         | 08/22/25 10:34 | 08/22/25 16:58 | PB169362      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | FB-01-082025                       |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-11                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 960                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BF143601.D        | 1         | 08/22/25 10:34 | 08/27/25 11:53 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 100-52-7       | Benzaldehyde                | 10.4  | U         | 4.10 | 10.4       | ug/L  |
| 108-95-2       | Phenol                      | 5.20  | U         | 0.95 | 5.20       | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 5.20  | U         | 0.84 | 5.20       | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 5.20  | U         | 0.60 | 5.20       | ug/L  |
| 95-48-7        | 2-Methylphenol              | 5.20  | U         | 1.20 | 5.20       | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 5.20  | U         | 1.30 | 5.20       | ug/L  |
| 98-86-2        | Acetophenone                | 5.20  | U         | 0.77 | 5.20       | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 10.4  | U         | 1.10 | 10.4       | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 2.60  | U         | 1.50 | 2.60       | ug/L  |
| 67-72-1        | Hexachloroethane            | 5.20  | U         | 0.68 | 5.20       | ug/L  |
| 98-95-3        | Nitrobenzene                | 5.20  | U         | 0.79 | 5.20       | ug/L  |
| 78-59-1        | Isophorone                  | 5.20  | U         | 0.78 | 5.20       | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 5.20  | U         | 1.80 | 5.20       | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 5.20  | U         | 1.90 | 5.20       | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 5.20  | U         | 0.71 | 5.20       | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 5.20  | U         | 0.54 | 5.20       | ug/L  |
| 91-20-3        | Naphthalene                 | 5.20  | U         | 0.52 | 5.20       | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 5.20  | U         | 0.88 | 5.20       | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 5.20  | U         | 0.56 | 5.20       | ug/L  |
| 105-60-2       | Caprolactam                 | 10.4  | U         | 1.20 | 10.4       | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 5.20  | U         | 0.61 | 5.20       | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 5.20  | U         | 0.58 | 5.20       | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 10.4  | U         | 3.80 | 10.4       | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 5.20  | U         | 0.53 | 5.20       | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 5.20  | U         | 0.65 | 5.20       | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 5.20  | U         | 0.55 | 5.20       | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 5.20  | U         | 0.64 | 5.20       | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 5.20  | U         | 1.30 | 5.20       | ug/L  |
| 131-11-3       | Dimethylphthalate           | 5.20  | U         | 0.64 | 5.20       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | FB-01-082025                       |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-11                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 960                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BF143601.D        | 1         | 08/22/25 10:34 | 08/27/25 11:53 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 5.20  | U         | 0.78 | 5.20       | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 5.20  | U         | 0.96 | 5.20       | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 5.20  | U         | 1.10 | 5.20       | ug/L  |
| 83-32-9    | Acenaphthene               | 5.20  | U         | 0.57 | 5.20       | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 10.4  | U         | 6.20 | 10.4       | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 10.4  | U         | 2.50 | 10.4       | ug/L  |
| 132-64-9   | Dibenzofuran               | 5.20  | U         | 0.64 | 5.20       | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 5.20  | U         | 1.30 | 5.20       | ug/L  |
| 84-66-2    | Diethylphthalate           | 5.20  | U         | 0.72 | 5.20       | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 5.20  | U         | 0.71 | 5.20       | ug/L  |
| 86-73-7    | Fluorene                   | 5.20  | U         | 0.66 | 5.20       | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 5.20  | U         | 1.60 | 5.20       | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 10.4  | U         | 3.00 | 10.4       | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 5.20  | U         | 0.60 | 5.20       | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 5.20  | U         | 0.42 | 5.20       | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 5.20  | U         | 0.54 | 5.20       | ug/L  |
| 1912-24-9  | Atrazine                   | 5.20  | U         | 1.10 | 5.20       | ug/L  |
| 87-86-5    | Pentachlorophenol          | 10.4  | U         | 1.60 | 10.4       | ug/L  |
| 85-01-8    | Phenanthrene               | 5.20  | U         | 0.52 | 5.20       | ug/L  |
| 120-12-7   | Anthracene                 | 5.20  | U         | 0.64 | 5.20       | ug/L  |
| 86-74-8    | Carbazole                  | 5.20  | U         | 0.75 | 5.20       | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 5.20  | U         | 1.30 | 5.20       | ug/L  |
| 206-44-0   | Fluoranthene               | 5.20  | U         | 0.85 | 5.20       | ug/L  |
| 129-00-0   | Pyrene                     | 5.20  | U         | 0.52 | 5.20       | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 5.20  | U         | 2.00 | 5.20       | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 10.4  | U         | 0.97 | 10.4       | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 5.20  | U         | 0.47 | 5.20       | ug/L  |
| 218-01-9   | Chrysene                   | 5.20  | U         | 0.46 | 5.20       | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 5.20  | U         | 1.70 | 5.20       | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 10.4  | U         | 2.40 | 10.4       | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 5.20  | U         | 0.51 | 5.20       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | FB-01-082025                       |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-11                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 960                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BF143601.D        | 1         | 08/22/25 10:34 | 08/27/25 11:53 | PB169362      |

| CAS Number                            | Parameter                        | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------------------|----------------------------------|--------|-----------|----------|------------|----------|
| 207-08-9                              | Benzo(k)fluoranthene             | 5.20   | U         | 0.50     | 5.20       | ug/L     |
| 50-32-8                               | Benzo(a)pyrene                   | 5.20   | U         | 0.57     | 5.20       | ug/L     |
| 193-39-5                              | Indeno(1,2,3-cd)pyrene           | 5.20   | U         | 0.61     | 5.20       | ug/L     |
| 53-70-3                               | Dibenzo(a,h)anthracene           | 5.20   | U         | 0.70     | 5.20       | ug/L     |
| 191-24-2                              | Benzo(g,h,i)perylene             | 5.20   | U         | 0.72     | 5.20       | ug/L     |
| 95-94-3                               | 1,2,4,5-Tetrachlorobenzene       | 5.20   | U         | 0.54     | 5.20       | ug/L     |
| 123-91-1                              | 1,4-Dioxane                      | 5.20   | U         | 1.00     | 5.20       | ug/L     |
| 58-90-2                               | 2,3,4,6-Tetrachlorophenol        | 5.20   | U         | 0.75     | 5.20       | ug/L     |
| <b>SURROGATES</b>                     |                                  |        |           |          |            |          |
| 367-12-4                              | 2-Fluorophenol                   | 66.7   |           | 23 - 138 | 44%        | SPK: 150 |
| 13127-88-3                            | Phenol-d6                        | 42.0   |           | 10 - 134 | 28%        | SPK: 150 |
| 4165-60-0                             | Nitrobenzene-d5                  | 114    |           | 67 - 132 | 114%       | SPK: 100 |
| 321-60-8                              | 2-Fluorobiphenyl                 | 93.8   |           | 52 - 132 | 94%        | SPK: 100 |
| 118-79-6                              | 2,4,6-Tribromophenol             | 147    |           | 44 - 137 | 98%        | SPK: 150 |
| 1718-51-0                             | Terphenyl-d14                    | 99.4   |           | 42 - 152 | 99%        | SPK: 100 |
| <b>INTERNAL STANDARDS</b>             |                                  |        |           |          |            |          |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4           | 111000 | 6.893     |          |            |          |
| 1146-65-2                             | Naphthalene-d8                   | 429000 | 8.169     |          |            |          |
| 15067-26-2                            | Acenaphthene-d10                 | 235000 | 9.922     |          |            |          |
| 1517-22-2                             | Phenanthrene-d10                 | 345000 | 11.41     |          |            |          |
| 1719-03-5                             | Chrysene-d12                     | 186000 | 14.045    |          |            |          |
| 1520-96-3                             | Perylene-d12                     | 227000 | 15.521    |          |            |          |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                                  |        |           |          |            |          |
| 000123-42-2                           | 2-Pentanone, 4-hydroxy-4-methyl- | 6.90   | AB        |          | 5.11       | ug/L     |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | FB-01-082025                       |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-11                           |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 960                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BF143601.D        | 1         | 08/22/25 10:34 | 08/27/25 11:53 | PB169362      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products



# QC SUMMARY

### Surrogate Summary

SW-846

SDG No.: Q2924

Client: AECOM

Analytical Method: 8270E

| Lab Sample ID | Client ID        | Parameter            | Spike (PPM) | Result (PPM) | Recovery (%) | Qual | Limits (%) |      |
|---------------|------------------|----------------------|-------------|--------------|--------------|------|------------|------|
|               |                  |                      |             |              |              |      | Low        | High |
| PB169362BL    | PB169362BL       | 2-Fluorophenol       | 150         | 125          | 83           |      | 23         | 138  |
|               |                  | Phenol-d6            | 150         | 120          | 80           |      | 10         | 134  |
|               |                  | Nitrobenzene-d5      | 100         | 75.8         | 76           |      | 67         | 132  |
|               |                  | 2-Fluorobiphenyl     | 100         | 73.3         | 73           |      | 52         | 132  |
|               |                  | 2,4,6-Tribromophenol | 150         | 123          | 82           |      | 44         | 137  |
|               |                  | Terphenyl-d14        | 100         | 70.1         | 70           |      | 42         | 152  |
| PB169362BS    | PB169362BS       | 2-Fluorophenol       | 150         | 116          | 77           |      | 23         | 138  |
|               |                  | Phenol-d6            | 150         | 113          | 75           |      | 10         | 134  |
|               |                  | Nitrobenzene-d5      | 100         | 72.1         | 72           |      | 67         | 132  |
|               |                  | 2-Fluorobiphenyl     | 100         | 70.9         | 71           |      | 52         | 132  |
|               |                  | 2,4,6-Tribromophenol | 150         | 119          | 80           |      | 44         | 137  |
|               |                  | Terphenyl-d14        | 100         | 75.5         | 75           |      | 42         | 152  |
| Q2924-01      | MW-10C-082025    | 2-Fluorophenol       | 150         | 59.1         | 39           |      | 23         | 138  |
|               |                  | Phenol-d6            | 150         | 38.3         | 26           |      | 10         | 134  |
|               |                  | Nitrobenzene-d5      | 100         | 99.8         | 100          |      | 67         | 132  |
|               |                  | 2-Fluorobiphenyl     | 100         | 88.3         | 88           |      | 52         | 132  |
|               |                  | 2,4,6-Tribromophenol | 150         | 151          | 101          |      | 44         | 137  |
|               |                  | Terphenyl-d14        | 100         | 102          | 102          |      | 42         | 152  |
| Q2924-02      | MW-201-082025    | 2-Fluorophenol       | 150         | 65.2         | 43           |      | 23         | 138  |
|               |                  | Phenol-d6            | 150         | 38.7         | 26           |      | 10         | 134  |
|               |                  | Nitrobenzene-d5      | 100         | 171          | 171          | *    | 67         | 132  |
|               |                  | 2-Fluorobiphenyl     | 100         | 78.8         | 79           |      | 52         | 132  |
|               |                  | 2,4,6-Tribromophenol | 150         | 136          | 91           |      | 44         | 137  |
|               |                  | Terphenyl-d14        | 100         | 74.0         | 74           |      | 42         | 152  |
| Q2924-02DL    | MW-201-082025DL  | 2-Fluorophenol       | 150         | 63.7         | 42           |      | 23         | 138  |
|               |                  | Phenol-d6            | 150         | 31.8         | 21           |      | 10         | 134  |
|               |                  | Nitrobenzene-d5      | 100         | 77.1         | 77           |      | 67         | 132  |
|               |                  | 2-Fluorobiphenyl     | 100         | 83.0         | 83           |      | 52         | 132  |
|               |                  | 2,4,6-Tribromophenol | 150         | 114          | 76           |      | 44         | 137  |
|               |                  | Terphenyl-d14        | 100         | 77.0         | 77           |      | 42         | 152  |
| Q2924-02DL2   | MW-201-082025DL2 | 2-Fluorophenol       | 150         | 67.8         | 45           |      | 23         | 138  |
|               |                  | Phenol-d6            | 150         | 14.0         | 9            | *    | 10         | 134  |
|               |                  | Nitrobenzene-d5      | 100         | 69.0         | 69           |      | 67         | 132  |
|               |                  | 2-Fluorobiphenyl     | 100         | 76.2         | 76           |      | 52         | 132  |
|               |                  | 2,4,6-Tribromophenol | 150         | 91.2         | 61           |      | 44         | 137  |
|               |                  | Terphenyl-d14        | 100         | 73.4         | 73           |      | 42         | 152  |
| Q2924-03MS    | MW-201-082025MS  | 2-Fluorophenol       | 150         | 67.6         | 45           |      | 23         | 138  |
|               |                  | Phenol-d6            | 150         | 41.4         | 28           |      | 10         | 134  |
|               |                  | Nitrobenzene-d5      | 100         | 148          | 148          | *    | 67         | 132  |
|               |                  | 2-Fluorobiphenyl     | 100         | 79.6         | 80           |      | 52         | 132  |
|               |                  | 2,4,6-Tribromophenol | 150         | 138          | 92           |      | 44         | 137  |
|               |                  | Terphenyl-d14        | 100         | 76.9         | 77           |      | 42         | 152  |
| Q2924-04MSD   | MW-201-082025MSD | 2-Fluorophenol       | 150         | 67.4         | 45           |      | 23         | 138  |
|               |                  | Phenol-d6            | 150         | 41.4         | 28           |      | 10         | 134  |
|               |                  | Nitrobenzene-d5      | 100         | 149          | 149          | *    | 67         | 132  |
|               |                  | 2-Fluorobiphenyl     | 100         | 78.1         | 78           |      | 52         | 132  |
|               |                  | 2,4,6-Tribromophenol | 150         | 135          | 90           |      | 44         | 137  |
|               |                  | Terphenyl-d14        | 100         | 76.2         | 76           |      | 42         | 152  |
| Q2924-05      | MW-2B-082025     | 2-Fluorophenol       | 150         | 73.5         | 49           |      | 23         | 138  |
|               |                  | Phenol-d6            | 150         | 53.5         | 36           |      | 10         | 134  |
|               |                  | Nitrobenzene-d5      | 100         | 118          | 118          |      | 67         | 132  |

### Surrogate Summary

SW-846

SDG No.: Q2924

Client: AECOM

Analytical Method: 8270E

| Lab Sample ID | Client ID       | Parameter            | Spike (PPM) | Result (PPM) | Recovery (%) | Qual | Limits (%) |      |
|---------------|-----------------|----------------------|-------------|--------------|--------------|------|------------|------|
|               |                 |                      |             |              |              |      | Low        | High |
| Q2924-05      | MW-2B-082025    | 2-Fluorobiphenyl     | 100         | 87.5         | 87           |      | 52         | 132  |
|               |                 | 2,4,6-Tribromophenol | 150         | 141          | 94           |      | 44         | 137  |
|               |                 | Terphenyl-d14        | 100         | 101          | 101          |      | 42         | 152  |
| Q2924-05DL    | MW-2B-082025DL  | 2-Fluorophenol       | 150         | 80.0         | 53           |      | 23         | 138  |
|               |                 | Phenol-d6            | 150         | 51.4         | 34           |      | 10         | 134  |
|               |                 | Nitrobenzene-d5      | 100         | 85.6         | 86           |      | 67         | 132  |
| Q2924-05DL2   | MW-2B-082025DL2 | 2-Fluorobiphenyl     | 100         | 90.6         | 91           |      | 52         | 132  |
|               |                 | 2,4,6-Tribromophenol | 150         | 150          | 100          |      | 44         | 137  |
|               |                 | Terphenyl-d14        | 100         | 93.2         | 93           |      | 42         | 152  |
| Q2924-06      | MW-2A-082025    | 2-Fluorophenol       | 150         | 79.8         | 53           |      | 23         | 138  |
|               |                 | Phenol-d6            | 150         | 48.0         | 32           |      | 10         | 134  |
|               |                 | Nitrobenzene-d5      | 100         | 80.4         | 80           |      | 67         | 132  |
| Q2924-07      | MW-18C-082025   | 2-Fluorobiphenyl     | 100         | 91.2         | 91           |      | 52         | 132  |
|               |                 | 2,4,6-Tribromophenol | 150         | 139          | 93           |      | 44         | 137  |
|               |                 | Terphenyl-d14        | 100         | 101          | 101          |      | 42         | 152  |
| Q2924-08      | MW-16A-082025   | 2-Fluorophenol       | 150         | 64.6         | 43           |      | 23         | 138  |
|               |                 | Phenol-d6            | 150         | 39.1         | 26           |      | 10         | 134  |
|               |                 | Nitrobenzene-d5      | 100         | 83.4         | 83           |      | 67         | 132  |
| Q2924-09      | MW-16C-082025   | 2-Fluorobiphenyl     | 100         | 78.0         | 78           |      | 52         | 132  |
|               |                 | 2,4,6-Tribromophenol | 150         | 141          | 94           |      | 44         | 137  |
|               |                 | Terphenyl-d14        | 100         | 77.0         | 77           |      | 42         | 152  |
| Q2924-10      | MW-8A-082025    | 2-Fluorophenol       | 150         | 67.3         | 45           |      | 23         | 138  |
|               |                 | Phenol-d6            | 150         | 41.9         | 28           |      | 10         | 134  |
|               |                 | Nitrobenzene-d5      | 100         | 113          | 113          |      | 67         | 132  |
| Q2924-11      | FB-01-082025    | 2-Fluorobiphenyl     | 100         | 95.1         | 95           |      | 52         | 132  |
|               |                 | 2,4,6-Tribromophenol | 150         | 142          | 95           |      | 44         | 137  |
|               |                 | Terphenyl-d14        | 100         | 96.9         | 97           |      | 42         | 152  |
| Q2924-12      | MW-16C-082025   | 2-Fluorophenol       | 150         | 59.0         | 39           |      | 23         | 138  |
|               |                 | Phenol-d6            | 150         | 34.9         | 23           |      | 10         | 134  |
|               |                 | Nitrobenzene-d5      | 100         | 84.0         | 84           |      | 67         | 132  |
| Q2924-13      | MW-16C-082025   | 2-Fluorobiphenyl     | 100         | 77.7         | 78           |      | 52         | 132  |
|               |                 | 2,4,6-Tribromophenol | 150         | 145          | 97           |      | 44         | 137  |
|               |                 | Terphenyl-d14        | 100         | 72.4         | 72           |      | 42         | 152  |
| Q2924-14      | MW-16C-082025   | 2-Fluorophenol       | 150         | 64.9         | 43           |      | 23         | 138  |
|               |                 | Phenol-d6            | 150         | 39.6         | 26           |      | 10         | 134  |
|               |                 | Nitrobenzene-d5      | 100         | 83.7         | 84           |      | 67         | 132  |
| Q2924-15      | MW-16C-082025   | 2-Fluorobiphenyl     | 100         | 79.7         | 80           |      | 52         | 132  |
|               |                 | 2,4,6-Tribromophenol | 150         | 141          | 94           |      | 44         | 137  |
|               |                 | Terphenyl-d14        | 100         | 76.3         | 76           |      | 42         | 152  |
| Q2924-16      | MW-8A-082025    | 2-Fluorophenol       | 150         | 64.1         | 43           |      | 23         | 138  |
|               |                 | Phenol-d6            | 150         | 39.6         | 26           |      | 10         | 134  |
|               |                 | Nitrobenzene-d5      | 100         | 84.7         | 85           |      | 67         | 132  |
| Q2924-17      | MW-8A-082025    | 2-Fluorobiphenyl     | 100         | 77.8         | 78           |      | 52         | 132  |
|               |                 | 2,4,6-Tribromophenol | 150         | 138          | 92           |      | 44         | 137  |
|               |                 | Terphenyl-d14        | 100         | 74.9         | 75           |      | 42         | 152  |
| Q2924-18      | FB-01-082025    | 2-Fluorophenol       | 150         | 66.7         | 44           |      | 23         | 138  |
|               |                 | Phenol-d6            | 150         | 42.0         | 28           |      | 10         | 134  |
|               |                 | Nitrobenzene-d5      | 100         | 114          | 114          |      | 67         | 132  |
| Q2924-19      | FB-01-082025    | 2-Fluorobiphenyl     | 100         | 93.8         | 94           |      | 52         | 132  |
|               |                 | 2,4,6-Tribromophenol | 150         | 147          | 98           |      | 44         | 137  |
|               |                 | Terphenyl-d14        | 100         | 99.4         | 99           |      | 42         | 152  |

**Matrix Spike/Matrix Spike Duplicate Summary**

**SW-846**

**SDG No.:** Q2924

**Analytical Method:** SW8270E

**Client:** AECOM

**DataFile:** BP025602.D

| Parameter                        | Spike | Sample Result                            | Result | Units | Rec  | Rec Qual | RPD | RPD Qual | Low | Limits High | RPD |
|----------------------------------|-------|------------------------------------------|--------|-------|------|----------|-----|----------|-----|-------------|-----|
| <b>Lab Sample ID: Q2924-03MS</b> |       | <b>Client Sample ID: MW-201-082025MS</b> |        |       |      |          |     |          |     |             |     |
| Benzaldehyde                     | 51.5  | 0                                        | 38.4   | ug/L  | 75   |          |     |          | 10  | 137         |     |
| Phenol                           | 51.5  | 0                                        | 16.0   | ug/L  | 31   |          |     |          | 10  | 130         |     |
| bis(2-Chloroethyl)ether          | 51.5  | 0                                        | 44.3   | ug/L  | 86   |          |     |          | 29  | 141         |     |
| 2-Chlorophenol                   | 51.5  | 0                                        | 39.6   | ug/L  | 77   |          |     |          | 23  | 127         |     |
| 2-Methylphenol                   | 51.5  | 0                                        | 33.8   | ug/L  | 66   |          |     |          | 60  | 131         |     |
| 2,2-oxybis(1-Chloropropane)      | 51.5  | 0                                        | 17.1   | ug/L  | 33   | *        |     |          | 36  | 141         |     |
| Acetophenone                     | 51.5  | 0                                        | 87.5   | ug/L  | 170  | *        |     |          | 31  | 164         |     |
| 3+4-Methylphenols                | 51.5  | 10.0                                     | 40.9   | ug/L  | 60   |          |     |          | 54  | 136         |     |
| N-Nitroso-di-n-propylamine       | 51.5  | 0                                        | 42.2   | ug/L  | 82   |          |     |          | 36  | 147         |     |
| Hexachloroethane                 | 51.5  | 0                                        | 93.2   | ug/L  | 181  | *        |     |          | 19  | 146         |     |
| Nitrobenzene                     | 51.5  | 0                                        | 85.8   | ug/L  | 167  | *        |     |          | 62  | 112         |     |
| Isophorone                       | 51.5  | 0                                        | 81.7   | ug/L  | 159  | *        |     |          | 39  | 146         |     |
| 2-Nitrophenol                    | 51.5  | 0                                        | 86.6   | ug/L  | 168  | *        |     |          | 30  | 148         |     |
| 2,4-Dimethylphenol               | 51.5  | 0                                        | 75.6   | ug/L  | 147  | *        |     |          | 17  | 143         |     |
| bis(2-Chloroethoxy)methane       | 51.5  | 0                                        | 70.3   | ug/L  | 137  |          |     |          | 39  | 143         |     |
| 2,4-Dichlorophenol               | 51.5  | 0                                        | 80.6   | ug/L  | 157  | *        |     |          | 22  | 146         |     |
| Naphthalene                      | 51.5  | 3100                                     | 2700   | ug/L  | -777 | *        |     |          | 17  | 157         |     |
| 4-Chloroaniline                  | 51.5  | 0                                        | 28.5   | ug/L  | 55   |          |     |          | 10  | 95          |     |
| Hexachlorobutadiene              | 51.5  | 0                                        | 71.5   | ug/L  | 139  | *        |     |          | 52  | 125         |     |
| Caprolactam                      | 51.5  | 0                                        | 16.7   | ug/L  | 32   |          |     |          | 10  | 130         |     |
| 4-Chloro-3-methylphenol          | 51.5  | 0                                        | 76.5   | ug/L  | 149  | *        |     |          | 17  | 148         |     |
| 2-Methylnaphthalene              | 51.5  | 920                                      | 800    | ug/L  | -233 | *        |     |          | 38  | 146         |     |
| Hexachlorocyclopentadiene        | 100   | 0                                        | 79.7   | ug/L  | 80   |          |     |          | 20  | 153         |     |
| 2,4,6-Trichlorophenol            | 51.5  | 0                                        | 50.2   | ug/L  | 97   |          |     |          | 78  | 112         |     |
| 2,4,5-Trichlorophenol            | 51.5  | 0                                        | 49.4   | ug/L  | 96   |          |     |          | 71  | 111         |     |
| 1,1-Biphenyl                     | 51.5  | 26.0                                     | 69.6   | ug/L  | 85   |          |     |          | 38  | 154         |     |
| 2-Chloronaphthalene              | 51.5  | 0                                        | 48.9   | ug/L  | 95   |          |     |          | 41  | 145         |     |
| 2-Nitroaniline                   | 51.5  | 0                                        | 54.1   | ug/L  | 105  |          |     |          | 39  | 151         |     |
| Dimethylphthalate                | 51.5  | 0                                        | 48.7   | ug/L  | 95   |          |     |          | 42  | 147         |     |
| Acenaphthylene                   | 51.5  | 160                                      | 190    | ug/L  | 58   |          |     |          | 40  | 141         |     |
| 2,6-Dinitrotoluene               | 51.5  | 0                                        | 51.4   | ug/L  | 100  |          |     |          | 43  | 148         |     |
| 3-Nitroaniline                   | 51.5  | 0                                        | 31.7   | ug/L  | 62   |          |     |          | 10  | 111         |     |
| Acenaphthene                     | 51.5  | 8.70                                     | 54.7   | ug/L  | 89   |          |     |          | 37  | 146         |     |
| 2,4-Dinitrophenol                | 100   | 0                                        | 120    | ug/L  | 120  |          |     |          | 14  | 167         |     |
| 4-Nitrophenol                    | 100   | 0                                        | 42.0   | ug/L  | 42   |          |     |          | 10  | 130         |     |
| Dibenzofuran                     | 51.5  | 4.00                                     | 51.3   | ug/L  | 92   |          |     |          | 41  | 145         |     |
| 2,4-Dinitrotoluene               | 51.5  | 0                                        | 53.3   | ug/L  | 103  |          |     |          | 74  | 137         |     |
| Diethylphthalate                 | 51.5  | 2.10                                     | 50.3   | ug/L  | 94   |          |     |          | 41  | 148         |     |
| 4-Chlorophenyl-phenylether       | 51.5  | 0                                        | 48.5   | ug/L  | 94   |          |     |          | 38  | 149         |     |
| Fluorene                         | 51.5  | 27.4                                     | 67.1   | ug/L  | 77   |          |     |          | 39  | 144         |     |
| 4-Nitroaniline                   | 51.5  | 0                                        | 48.5   | ug/L  | 94   |          |     |          | 27  | 138         |     |
| 4,6-Dinitro-2-methylphenol       | 51.5  | 0                                        | 55.1   | ug/L  | 107  |          |     |          | 32  | 175         |     |
| N-Nitrosodiphenylamine           | 51.5  | 0                                        | 51.1   | ug/L  | 99   |          |     |          | 40  | 150         |     |
| 4-Bromophenyl-phenylether        | 51.5  | 0                                        | 51.5   | ug/L  | 100  |          |     |          | 42  | 151         |     |
| Hexachlorobenzene                | 51.5  | 0                                        | 52.5   | ug/L  | 102  |          |     |          | 72  | 115         |     |
| Atrazine                         | 51.5  | 0                                        | 47.8   | ug/L  | 93   |          |     |          | 20  | 162         |     |

**Matrix Spike/Matrix Spike Duplicate Summary**

**SW-846**

**SDG No.:** Q2924

**Analytical Method:** SW8270E

**Client:** AECOM

**DataFile:** BP025602.D

| Parameter                  | Spike | Sample<br>Result | Result | Units | Rec | Rec<br>Qual | RPD | RPD<br>Qual | Low | Limits<br>High | RPD |
|----------------------------|-------|------------------|--------|-------|-----|-------------|-----|-------------|-----|----------------|-----|
| Pentachlorophenol          | 100   | 0                | 110    | ug/L  | 110 |             |     |             | 52  | 162            |     |
| Phenanthrene               | 51.5  | 29.6             | 65.4   | ug/L  | 70  |             |     |             | 40  | 147            |     |
| Anthracene                 | 51.5  | 6.10             | 53.4   | ug/L  | 92  |             |     |             | 41  | 146            |     |
| Carbazole                  | 51.5  | 13.8             | 61.7   | ug/L  | 93  |             |     |             | 37  | 154            |     |
| Di-n-butylphthalate        | 51.5  | 2.70             | 51.3   | ug/L  | 94  |             |     |             | 40  | 151            |     |
| Fluoranthene               | 51.5  | 2.30             | 51.0   | ug/L  | 95  |             |     |             | 42  | 146            |     |
| Pyrene                     | 51.5  | 3.00             | 50.5   | ug/L  | 92  |             |     |             | 41  | 149            |     |
| Butylbenzylphthalate       | 51.5  | 0                | 51.2   | ug/L  | 99  |             |     |             | 39  | 155            |     |
| 3,3-Dichlorobenzidine      | 51.5  | 0                | 41.5   | ug/L  | 81  |             |     |             | 10  | 114            |     |
| Benzo(a)anthracene         | 51.5  | 0                | 49.5   | ug/L  | 96  |             |     |             | 41  | 147            |     |
| Chrysene                   | 51.5  | 0                | 50.0   | ug/L  | 97  |             |     |             | 44  | 144            |     |
| bis(2-Ethylhexyl)phthalate | 51.5  | 0                | 48.4   | ug/L  | 94  |             |     |             | 33  | 160            |     |
| Di-n-octyl phthalate       | 51.5  | 0                | 51.0   | ug/L  | 99  |             |     |             | 36  | 158            |     |
| Benzo(b)fluoranthene       | 51.5  | 0                | 50.0   | ug/L  | 97  |             |     |             | 40  | 150            |     |
| Benzo(k)fluoranthene       | 51.5  | 0                | 49.1   | ug/L  | 95  |             |     |             | 40  | 147            |     |
| Benzo(a)pyrene             | 51.5  | 0                | 49.3   | ug/L  | 96  |             |     |             | 42  | 147            |     |
| Indeno(1,2,3-cd)pyrene     | 51.5  | 0                | 50.1   | ug/L  | 97  |             |     |             | 30  | 166            |     |
| Dibenz(a,h)anthracene      | 51.5  | 0                | 51.0   | ug/L  | 99  |             |     |             | 23  | 172            |     |
| Benzo(g,h,i)perylene       | 51.5  | 0                | 50.2   | ug/L  | 97  |             |     |             | 27  | 167            |     |
| 1,2,4,5-Tetrachlorobenzene | 51.5  | 0                | 50.1   | ug/L  | 97  |             |     |             | 89  | 102            |     |
| 1,4-Dioxane                | 51.5  | 0                | 21.2   | ug/L  | 41  |             |     |             | 38  | 130            |     |
| 2,3,4,6-Tetrachlorophenol  | 51.5  | 0                | 51.1   | ug/L  | 99  |             |     |             | 91  | 111            |     |

**Matrix Spike/Matrix Spike Duplicate Summary**

**SW-846**

**SDG No.:** Q2924

**Analytical Method:** SW8270E

**Client:** AECOM

**DataFile:** BP025603.D

| Parameter                   | Spike              | Sample Result            | Result                  | Units | Rec  | Rec Qual | RPD | RPD Qual | Low | Limits High | RPD |
|-----------------------------|--------------------|--------------------------|-------------------------|-------|------|----------|-----|----------|-----|-------------|-----|
| <b>Lab Sample ID:</b>       | <b>Q2924-04MSD</b> | <b>Client Sample ID:</b> | <b>MW-201-082025MSD</b> |       |      |          |     |          |     |             |     |
| Benzaldehyde                | 52.6               | 0                        | 37.7                    | ug/L  | 72   |          | 4   |          | 10  | 137         | 20  |
| Phenol                      | 52.6               | 0                        | 16.4                    | ug/L  | 31   |          | 0   |          | 10  | 130         | 20  |
| bis(2-Chloroethyl)ether     | 52.6               | 0                        | 44.2                    | ug/L  | 84   |          | 2   |          | 29  | 141         | 20  |
| 2-Chlorophenol              | 52.6               | 0                        | 39.6                    | ug/L  | 75   |          | 3   |          | 23  | 127         | 20  |
| 2-Methylphenol              | 52.6               | 0                        | 34.5                    | ug/L  | 66   |          | 0   |          | 60  | 131         | 20  |
| 2,2-oxybis(1-Chloropropane) | 52.6               | 0                        | 18.4                    | ug/L  | 35   | *        | 6   |          | 36  | 141         | 20  |
| Acetophenone                | 52.6               | 0                        | 90.6                    | ug/L  | 172  | *        | 1   |          | 31  | 164         | 20  |
| 3+4-Methylphenols           | 52.6               | 10.0                     | 41.7                    | ug/L  | 60   |          | 0   |          | 54  | 136         | 20  |
| N-Nitroso-di-n-propylamine  | 52.6               | 0                        | 43.2                    | ug/L  | 82   |          | 0   |          | 36  | 147         | 20  |
| Hexachloroethane            | 52.6               | 0                        | 96.9                    | ug/L  | 184  | *        | 2   |          | 19  | 146         | 20  |
| Nitrobenzene                | 52.6               | 0                        | 89.1                    | ug/L  | 169  | *        | 1   |          | 62  | 112         | 20  |
| Isophorone                  | 52.6               | 0                        | 85.0                    | ug/L  | 162  | *        | 2   |          | 39  | 146         | 20  |
| 2-Nitrophenol               | 52.6               | 0                        | 90.4                    | ug/L  | 172  | *        | 2   |          | 30  | 148         | 20  |
| 2,4-Dimethylphenol          | 52.6               | 0                        | 78.4                    | ug/L  | 149  | *        | 1   |          | 17  | 143         | 20  |
| bis(2-Chloroethoxy)methane  | 52.6               | 0                        | 75.2                    | ug/L  | 143  |          | 4   |          | 39  | 143         | 20  |
| 2,4-Dichlorophenol          | 52.6               | 0                        | 85.1                    | ug/L  | 162  | *        | 3   |          | 22  | 146         | 20  |
| Naphthalene                 | 52.6               | 3100                     | 3000                    | ug/L  | -190 | *        | 121 | *        | 17  | 157         | 20  |
| 4-Chloroaniline             | 52.6               | 0                        | 31.3                    | ug/L  | 60   |          | 9   |          | 10  | 95          | 20  |
| Hexachlorobutadiene         | 52.6               | 0                        | 74.8                    | ug/L  | 142  | *        | 2   |          | 52  | 125         | 20  |
| Caprolactam                 | 52.6               | 0                        | 17.0                    | ug/L  | 32   |          | 0   |          | 10  | 130         | 20  |
| 4-Chloro-3-methylphenol     | 52.6               | 0                        | 79.4                    | ug/L  | 151  | *        | 1   |          | 17  | 148         | 20  |
| 2-Methylnaphthalene         | 52.6               | 920                      | 850                     | ug/L  | -133 | *        | 55  | *        | 38  | 146         | 20  |
| Hexachlorocyclopentadiene   | 110                | 0                        | 76.9                    | ug/L  | 70   |          | 13  |          | 20  | 153         | 20  |
| 2,4,6-Trichlorophenol       | 52.6               | 0                        | 50.2                    | ug/L  | 95   |          | 2   |          | 78  | 112         | 20  |
| 2,4,5-Trichlorophenol       | 52.6               | 0                        | 49.9                    | ug/L  | 95   |          | 1   |          | 71  | 111         | 20  |
| 1,1-Biphenyl                | 52.6               | 26.0                     | 70.0                    | ug/L  | 84   |          | 1   |          | 38  | 154         | 20  |
| 2-Chloronaphthalene         | 52.6               | 0                        | 48.9                    | ug/L  | 93   |          | 2   |          | 41  | 145         | 20  |
| 2-Nitroaniline              | 52.6               | 0                        | 53.9                    | ug/L  | 102  |          | 3   |          | 39  | 151         | 20  |
| Dimethylphthalate           | 52.6               | 0                        | 47.8                    | ug/L  | 91   |          | 4   |          | 42  | 147         | 20  |
| Acenaphthylene              | 52.6               | 160                      | 190                     | ug/L  | 57   |          | 2   |          | 40  | 141         | 20  |
| 2,6-Dinitrotoluene          | 52.6               | 0                        | 51.3                    | ug/L  | 98   |          | 2   |          | 43  | 148         | 20  |
| 3-Nitroaniline              | 52.6               | 0                        | 32.4                    | ug/L  | 62   |          | 0   |          | 10  | 111         | 20  |
| Acenaphthene                | 52.6               | 8.70                     | 53.5                    | ug/L  | 85   |          | 5   |          | 37  | 146         | 20  |
| 2,4-Dinitrophenol           | 110                | 0                        | 120                     | ug/L  | 109  |          | 10  |          | 14  | 167         | 20  |
| 4-Nitrophenol               | 110                | 0                        | 42.7                    | ug/L  | 39   |          | 7   |          | 10  | 130         | 20  |
| Dibenzofuran                | 52.6               | 4.00                     | 51.2                    | ug/L  | 90   |          | 2   |          | 41  | 145         | 20  |
| 2,4-Dinitrotoluene          | 52.6               | 0                        | 52.5                    | ug/L  | 100  |          | 3   |          | 74  | 137         | 20  |
| Diethylphthalate            | 52.6               | 2.10                     | 48.9                    | ug/L  | 89   |          | 5   |          | 41  | 148         | 20  |
| 4-Chlorophenyl-phenylether  | 52.6               | 0                        | 48.5                    | ug/L  | 92   |          | 2   |          | 38  | 149         | 20  |
| Fluorene                    | 52.6               | 27.4                     | 68.0                    | ug/L  | 77   |          | 0   |          | 39  | 144         | 20  |
| 4-Nitroaniline              | 52.6               | 0                        | 49.4                    | ug/L  | 94   |          | 0   |          | 27  | 138         | 20  |
| 4,6-Dinitro-2-methylphenol  | 52.6               | 0                        | 55.7                    | ug/L  | 106  |          | 1   |          | 32  | 175         | 20  |
| N-Nitrosodiphenylamine      | 52.6               | 0                        | 50.5                    | ug/L  | 96   |          | 3   |          | 40  | 150         | 20  |
| 4-Bromophenyl-phenylether   | 52.6               | 0                        | 52.0                    | ug/L  | 99   |          | 1   |          | 42  | 151         | 20  |
| Hexachlorobenzene           | 52.6               | 0                        | 52.3                    | ug/L  | 99   |          | 3   |          | 72  | 115         | 20  |
| Atrazine                    | 52.6               | 0                        | 48.5                    | ug/L  | 92   |          | 1   |          | 20  | 162         | 20  |

**Matrix Spike/Matrix Spike Duplicate Summary**

**SW-846**

**SDG No.:** Q2924

**Analytical Method:** SW8270E

**Client:** AECOM

**DataFile:** BP025603.D

| Parameter                  | Spike | Sample<br>Result | Result | Units | Rec | Rec<br>Qual | RPD | RPD<br>Qual | Low | Limits<br>High | RPD |
|----------------------------|-------|------------------|--------|-------|-----|-------------|-----|-------------|-----|----------------|-----|
| Pentachlorophenol          | 110   | 0                | 110    | ug/L  | 100 |             | 10  |             | 52  | 162            | 20  |
| Phenanthrene               | 52.6  | 29.6             | 67.2   | ug/L  | 71  |             | 1   |             | 40  | 147            | 20  |
| Anthracene                 | 52.6  | 6.10             | 53.2   | ug/L  | 90  |             | 2   |             | 41  | 146            | 20  |
| Carbazole                  | 52.6  | 13.8             | 63.5   | ug/L  | 94  |             | 1   |             | 37  | 154            | 20  |
| Di-n-butylphthalate        | 52.6  | 2.70             | 52.6   | ug/L  | 95  |             | 1   |             | 40  | 151            | 20  |
| Fluoranthene               | 52.6  | 2.30             | 51.6   | ug/L  | 94  |             | 1   |             | 42  | 146            | 20  |
| Pyrene                     | 52.6  | 3.00             | 50.1   | ug/L  | 90  |             | 2   |             | 41  | 149            | 20  |
| Butylbenzylphthalate       | 52.6  | 0                | 52.0   | ug/L  | 99  |             | 0   |             | 39  | 155            | 20  |
| 3,3-Dichlorobenzidine      | 52.6  | 0                | 43.4   | ug/L  | 83  |             | 2   |             | 10  | 114            | 20  |
| Benzo(a)anthracene         | 52.6  | 0                | 49.9   | ug/L  | 95  |             | 1   |             | 41  | 147            | 20  |
| Chrysene                   | 52.6  | 0                | 51.2   | ug/L  | 97  |             | 0   |             | 44  | 144            | 20  |
| bis(2-Ethylhexyl)phthalate | 52.6  | 0                | 49.7   | ug/L  | 94  |             | 0   |             | 33  | 160            | 20  |
| Di-n-octyl phthalate       | 52.6  | 0                | 52.0   | ug/L  | 99  |             | 0   |             | 36  | 158            | 20  |
| Benzo(b)fluoranthene       | 52.6  | 0                | 50.4   | ug/L  | 96  |             | 1   |             | 40  | 150            | 20  |
| Benzo(k)fluoranthene       | 52.6  | 0                | 49.6   | ug/L  | 94  |             | 1   |             | 40  | 147            | 20  |
| Benzo(a)pyrene             | 52.6  | 0                | 49.8   | ug/L  | 95  |             | 1   |             | 42  | 147            | 20  |
| Indeno(1,2,3-cd)pyrene     | 52.6  | 0                | 50.9   | ug/L  | 97  |             | 0   |             | 30  | 166            | 20  |
| Dibenz(a,h)anthracene      | 52.6  | 0                | 51.6   | ug/L  | 98  |             | 1   |             | 23  | 172            | 20  |
| Benzo(g,h,i)perylene       | 52.6  | 0                | 50.6   | ug/L  | 96  |             | 1   |             | 27  | 167            | 20  |
| 1,2,4,5-Tetrachlorobenzene | 52.6  | 0                | 49.8   | ug/L  | 95  |             | 2   |             | 89  | 102            | 20  |
| 1,4-Dioxane                | 52.6  | 0                | 21.8   | ug/L  | 41  |             | 0   |             | 38  | 130            | 20  |
| 2,3,4,6-Tetrachlorophenol  | 52.6  | 0                | 50.6   | ug/L  | 96  |             | 3   |             | 91  | 111            | 20  |

## Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q2924

Analytical Method: 8270E

Client: AECOM

DataFile: BP025581.D

| Lab Sample ID | Parameter                   | Spike | Result | Unit | Rec | RPD | Qual | RPD  | Limits |      | RPD |
|---------------|-----------------------------|-------|--------|------|-----|-----|------|------|--------|------|-----|
|               |                             |       |        |      |     |     |      | Qual | Low    | High |     |
| PB169362BS    | Benzaldehyde                | 50    | 22.9   | ug/L | 46  |     |      |      | 10     | 162  |     |
|               | Phenol                      | 50    | 41.7   | ug/L | 83  |     |      |      | 66     | 118  |     |
|               | bis(2-Chloroethyl)ether     | 50    | 40.1   | ug/L | 80  |     |      |      | 62     | 103  |     |
|               | 2-Chlorophenol              | 50    | 41.6   | ug/L | 83  |     |      |      | 70     | 117  |     |
|               | 2-Methylphenol              | 50    | 41.0   | ug/L | 82  |     |      |      | 69     | 109  |     |
|               | 2,2-oxybis(1-Chloropropane) | 50    | 39.7   | ug/L | 79  |     |      |      | 65     | 100  |     |
|               | Acetophenone                | 50    | 39.8   | ug/L | 80  |     |      |      | 60     | 104  |     |
|               | 3+4-Methylphenols           | 50    | 39.8   | ug/L | 80  |     |      |      | 67     | 106  |     |
|               | N-Nitroso-di-n-propylamine  | 50    | 36.5   | ug/L | 73  |     |      |      | 57     | 107  |     |
|               | Hexachloroethane            | 50    | 40.3   | ug/L | 81  |     |      |      | 76     | 118  |     |
|               | Nitrobenzene                | 50    | 41.1   | ug/L | 82  |     |      |      | 58     | 106  |     |
|               | Isophorone                  | 50    | 39.2   | ug/L | 78  |     |      |      | 61     | 102  |     |
|               | 2-Nitrophenol               | 50    | 42.7   | ug/L | 85  |     |      |      | 70     | 115  |     |
|               | 2,4-Dimethylphenol          | 50    | 42.3   | ug/L | 85  |     |      |      | 42     | 142  |     |
|               | bis(2-Chloroethoxy)methane  | 50    | 40.1   | ug/L | 80  |     |      |      | 58     | 109  |     |
|               | 2,4-Dichlorophenol          | 50    | 43.3   | ug/L | 87  |     |      |      | 66     | 115  |     |
|               | Naphthalene                 | 50    | 39.7   | ug/L | 79  |     |      |      | 64     | 107  |     |
|               | 4-Chloroaniline             | 50    | 23.9   | ug/L | 48  |     |      |      | 10     | 85   |     |
|               | Hexachlorobutadiene         | 50    | 40.8   | ug/L | 82  |     |      |      | 69     | 101  |     |
|               | Caprolactam                 | 50    | 39.6   | ug/L | 79  |     |      |      | 58     | 128  |     |
|               | 4-Chloro-3-methylphenol     | 50    | 42.3   | ug/L | 85  |     |      |      | 65     | 114  |     |
|               | 2-Methylnaphthalene         | 50    | 39.4   | ug/L | 79  |     |      |      | 64     | 107  |     |
|               | Hexachlorocyclopentadiene   | 100   | 81.0   | ug/L | 81  |     |      |      | 36     | 160  |     |
|               | 2,4,6-Trichlorophenol       | 50    | 44.1   | ug/L | 88  |     |      |      | 61     | 110  |     |
|               | 2,4,5-Trichlorophenol       | 50    | 44.6   | ug/L | 89  |     |      |      | 70     | 106  |     |
|               | 1,1-Biphenyl                | 50    | 41.8   | ug/L | 84  |     |      |      | 72     | 98   |     |
|               | 2-Chloronaphthalene         | 50    | 40.9   | ug/L | 82  |     |      |      | 59     | 106  |     |
|               | 2-Nitroaniline              | 50    | 44.0   | ug/L | 88  |     |      |      | 73     | 114  |     |
|               | Dimethylphthalate           | 50    | 41.0   | ug/L | 82  |     |      |      | 64     | 103  |     |
|               | Acenaphthylene              | 50    | 40.9   | ug/L | 82  |     |      |      | 79     | 103  |     |
|               | 2,6-Dinitrotoluene          | 50    | 43.4   | ug/L | 87  |     |      |      | 64     | 110  |     |
|               | 3-Nitroaniline              | 50    | 31.2   | ug/L | 62  |     |      |      | 28     | 100  |     |
|               | Acenaphthene                | 50    | 39.5   | ug/L | 79  |     |      |      | 59     | 113  |     |
|               | 2,4-Dinitrophenol           | 100   | 100    | ug/L | 100 |     |      |      | 36     | 166  |     |
|               | 4-Nitrophenol               | 100   | 82.1   | ug/L | 82  |     |      |      | 45     | 147  |     |
|               | Dibenzofuran                | 50    | 40.2   | ug/L | 80  |     |      |      | 65     | 106  |     |
|               | 2,4-Dinitrotoluene          | 50    | 43.7   | ug/L | 87  |     |      |      | 60     | 115  |     |
|               | Diethylphthalate            | 50    | 41.5   | ug/L | 83  |     |      |      | 63     | 105  |     |
|               | 4-Chlorophenyl-phenylether  | 50    | 40.1   | ug/L | 80  |     |      |      | 61     | 104  |     |
|               | Fluorene                    | 50    | 40.2   | ug/L | 80  |     |      |      | 64     | 107  |     |
|               | 4-Nitroaniline              | 50    | 41.1   | ug/L | 82  |     |      |      | 55     | 125  |     |
|               | 4,6-Dinitro-2-methylphenol  | 50    | 46.7   | ug/L | 93  |     |      |      | 62     | 132  |     |
|               | N-Nitrosodiphenylamine      | 50    | 42.8   | ug/L | 86  |     |      |      | 61     | 109  |     |
|               | 4-Bromophenyl-phenylether   | 50    | 42.2   | ug/L | 84  |     |      |      | 73     | 103  |     |

## Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q2924

Analytical Method: 8270E

Client: AECOM

DataFile: BP025581.D

| Lab Sample ID | Parameter                  | Spike | Result | Unit | Rec | RPD | RPD  |      | Limits |      | RPD |
|---------------|----------------------------|-------|--------|------|-----|-----|------|------|--------|------|-----|
|               |                            |       |        |      |     |     | Qual | Qual | Low    | High |     |
| PB169362BS    | Hexachlorobenzene          | 50    | 42.0   | ug/L | 84  |     |      |      | 73     | 106  |     |
|               | Atrazine                   | 50    | 41.7   | ug/L | 83  |     |      |      | 76     | 120  |     |
|               | Pentachlorophenol          | 100   | 86.9   | ug/L | 87  |     |      |      | 47     | 114  |     |
|               | Phenanthrene               | 50    | 40.9   | ug/L | 82  |     |      |      | 62     | 109  |     |
|               | Anthracene                 | 50    | 41.4   | ug/L | 83  |     |      |      | 65     | 110  |     |
|               | Carbazole                  | 50    | 40.8   | ug/L | 82  |     |      |      | 62     | 106  |     |
|               | Di-n-butylphthalate        | 50    | 41.6   | ug/L | 83  |     |      |      | 64     | 106  |     |
|               | Fluoranthene               | 50    | 40.0   | ug/L | 80  |     |      |      | 64     | 110  |     |
|               | Pyrene                     | 50    | 43.4   | ug/L | 87  |     |      |      | 71     | 103  |     |
|               | Butylbenzylphthalate       | 50    | 44.3   | ug/L | 89  |     |      |      | 61     | 105  |     |
|               | 3,3-Dichlorobenzidine      | 50    | 31.9   | ug/L | 64  |     |      |      | 43     | 108  |     |
|               | Benzo(a)anthracene         | 50    | 42.2   | ug/L | 84  |     |      |      | 62     | 107  |     |
|               | Chrysene                   | 50    | 41.7   | ug/L | 83  |     |      |      | 61     | 108  |     |
|               | bis(2-Ethylhexyl)phthalate | 50    | 42.7   | ug/L | 85  |     |      |      | 59     | 110  |     |
|               | Di-n-octyl phthalate       | 50    | 43.7   | ug/L | 87  |     |      |      | 52     | 139  |     |
|               | Benzo(b)fluoranthene       | 50    | 41.7   | ug/L | 83  |     |      |      | 77     | 113  |     |
|               | Benzo(k)fluoranthene       | 50    | 41.8   | ug/L | 84  |     |      |      | 77     | 105  |     |
|               | Benzo(a)pyrene             | 50    | 41.6   | ug/L | 83  |     |      |      | 72     | 131  |     |
|               | Indeno(1,2,3-cd)pyrene     | 50    | 41.4   | ug/L | 83  |     |      |      | 72     | 105  |     |
|               | Dibenz(a,h)anthracene      | 50    | 41.8   | ug/L | 84  |     |      |      | 78     | 115  |     |
|               | Benzo(g,h,i)perylene       | 50    | 41.4   | ug/L | 83  |     |      |      | 75     | 118  |     |
|               | 1,2,4,5-Tetrachlorobenzene | 50    | 42.0   | ug/L | 84  |     |      |      | 72     | 101  |     |
|               | 1,4-Dioxane                | 50    | 34.9   | ug/L | 70  |     |      |      | 38     | 125  |     |
|               | 2,3,4,6-Tetrachlorophenol  | 50    | 43.0   | ug/L | 86  |     |      |      | 63     | 116  |     |

4B

SEMIVOLATILE METHOD BLANK SUMMARY

Client ID

PB169362BL

Lab Name: Alliance

Contract: AEC002

Lab Code: ACE

SDG NO.: Q2924

Lab File ID: BP025562.D

Lab Sample ID: PB169362BL

Instrument ID: BNA\_P

Date Extracted: 08/22/2025

Matrix: (soil/water) Water

Date Analyzed: 08/25/2025

Level: (low/med) LOW

Time Analyzed: 12:54

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

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED |
|-------------------|------------------|----------------|------------------|
| PB169362BS        | PB169362BS       | BP025581.D     | 08/26/2025       |
| MW-16A-082025     | Q2924-08         | BP025588.D     | 08/26/2025       |
| MW-16C-082025     | Q2924-09         | BP025589.D     | 08/26/2025       |
| MW-2A-082025      | Q2924-06         | BP025590.D     | 08/26/2025       |
| MW-201-082025     | Q2924-02         | BP025601.D     | 08/27/2025       |
| MW-201-082025MS   | Q2924-03MS       | BP025602.D     | 08/27/2025       |
| MW-201-082025MSD  | Q2924-04MSD      | BP025603.D     | 08/27/2025       |
| MW-10C-082025     | Q2924-01         | BF143581.D     | 08/22/2025       |
| MW-2B-082025      | Q2924-05         | BF143582.D     | 08/22/2025       |
| FB-01-082025      | Q2924-11         | BF143601.D     | 08/27/2025       |
| MW-18C-082025     | Q2924-07         | BF143603.D     | 08/27/2025       |
| MW-8A-082025      | Q2924-10         | BP025552.D     | 08/22/2025       |

COMMENTS: \_\_\_\_\_

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BF143476.D  
Instrument ID: BNA\_F

Contract: AECO02  
SDG NO.: Q2924  
DFTPP Injection Date: 08/20/2025  
DFTPP Injection Time: 10:26

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.5 ( 1.9 ) 1        |
| 69  | Mass 69 relative abundance         | 27.4                 |
| 70  | Less than 2.0% of mass 69          | 0.1 ( 0.5 ) 1        |
| 197 | Less than 2.0% of mass 198         | 0.0                  |
| 198 | Base Peak, 100% relative abundance | 100                  |
| 199 | 5.0 to 9.0% of mass 198            | 5.5                  |
| 365 | Greater than 1% of mass 198        | 3.1                  |
| 441 | Present, but less than mass 443    | 15.4                 |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 19.3 ( 19.3 ) 2      |

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|-------------------|------------------|----------------|------------------|------------------|
| SSTDICC2.5        | SSTDICC2.5       | BF143477.D     | 08/20/2025       | 10:55            |
| SSTDICC005        | SSTDICC005       | BF143478.D     | 08/20/2025       | 11:25            |
| SSTDICC010        | SSTDICC010       | BF143479.D     | 08/20/2025       | 11:54            |
| SSTDICC020        | SSTDICC020       | BF143480.D     | 08/20/2025       | 12:23            |
| SSTDICCC040       | SSTDICCC040      | BF143481.D     | 08/20/2025       | 12:53            |
| SSTDICC050        | SSTDICC050       | BF143482.D     | 08/20/2025       | 13:22            |
| SSTDICC060        | SSTDICC060       | BF143483.D     | 08/20/2025       | 13:51            |
| SSTDICC080        | SSTDICC080       | BF143484.D     | 08/20/2025       | 14:20            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BF143562.D  
Instrument ID: BNA\_F

Contract: AECO02  
SDG NO.: Q2924  
DFTPP Injection Date: 08/22/2025  
DFTPP Injection Time: 09:30

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.6 ( 2 ) 1          |
| 69  | Mass 69 relative abundance         | 29.5                 |
| 70  | Less than 2.0% of mass 69          | 0.2 ( 0.6 ) 1        |
| 197 | Less than 2.0% of mass 198         | 0.0                  |
| 198 | Base Peak, 100% relative abundance | 100                  |
| 199 | 5.0 to 9.0% of mass 198            | 5.6                  |
| 365 | Greater than 1% of mass 198        | 3.1                  |
| 441 | Present, but less than mass 443    | 15.3                 |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 19.3 ( 19.3 ) 2      |

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|-------------------|------------------|----------------|------------------|------------------|
| SSTDCCC040        | SSTDCCC040       | BF143563.D     | 08/22/2025       | 10:29            |
| MW-10C-082025     | Q2924-01         | BF143581.D     | 08/22/2025       | 19:21            |
| MW-2B-082025      | Q2924-05         | BF143582.D     | 08/22/2025       | 19:50            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BF143584.D  
Instrument ID: BNA\_F

Contract: AEC002  
SDG NO.: Q2924  
DFTPP Injection Date: 08/26/2025  
DFTPP Injection Time: 08:36

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.7 ( 1.9 ) 1        |
| 69  | Mass 69 relative abundance         | 35.3                 |
| 70  | Less than 2.0% of mass 69          | 0.2 ( 0.6 ) 1        |
| 197 | Less than 2.0% of mass 198         | 0.0                  |
| 198 | Base Peak, 100% relative abundance | 100                  |
| 199 | 5.0 to 9.0% of mass 198            | 6.7                  |
| 365 | Greater than 1% of mass 198        | 2.9                  |
| 441 | Present, but less than mass 443    | 14.4                 |
| 442 | Greater than 50% of mass 198       | 91.9                 |
| 443 | 15.0 - 24.0% of mass 442           | 16.5 ( 17.9 ) 2      |

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|-------------------|------------------|----------------|------------------|------------------|
| SSTDICC2.5        | SSTDICC2.5       | BF143585.D     | 08/26/2025       | 09:11            |
| SSTDICC005        | SSTDICC005       | BF143586.D     | 08/26/2025       | 09:39            |
| SSTDICC010        | SSTDICC010       | BF143587.D     | 08/26/2025       | 10:08            |
| SSTDICC020        | SSTDICC020       | BF143588.D     | 08/26/2025       | 10:37            |
| SSTDICCC040       | SSTDICCC040      | BF143589.D     | 08/26/2025       | 11:06            |
| SSTDICC050        | SSTDICC050       | BF143590.D     | 08/26/2025       | 11:35            |
| SSTDICC060        | SSTDICC060       | BF143591.D     | 08/26/2025       | 12:03            |
| SSTDICC080        | SSTDICC080       | BF143592.D     | 08/26/2025       | 12:32            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BF143595.D  
Instrument ID: BNA\_F

Contract: AECO02  
SDG NO.: Q2924  
DFTPP Injection Date: 08/27/2025  
DFTPP Injection Time: 08:48

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.6 ( 1.6 ) 1        |
| 69  | Mass 69 relative abundance         | 33.7                 |
| 70  | Less than 2.0% of mass 69          | 0.2 ( 0.6 ) 1        |
| 197 | Less than 2.0% of mass 198         | 0.0                  |
| 198 | Base Peak, 100% relative abundance | 100                  |
| 199 | 5.0 to 9.0% of mass 198            | 6.6                  |
| 365 | Greater than 1% of mass 198        | 3.2                  |
| 441 | Present, but less than mass 443    | 16                   |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 19.2 ( 19.2 ) 2      |

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|-------------------|------------------|----------------|------------------|------------------|
| SSTDCCC040        | SSTDCCC040       | BF143596.D     | 08/27/2025       | 09:16            |
| FB-01-082025      | Q2924-11         | BF143601.D     | 08/27/2025       | 11:53            |
| MW-18C-082025     | Q2924-07         | BF143603.D     | 08/27/2025       | 12:52            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BP025390.D  
Instrument ID: BNA\_P

Contract: AECO02  
SDG NO.: Q2924  
DFTPP Injection Date: 08/14/2025  
DFTPP Injection Time: 10:30

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.4 ( 1.6 ) 1        |
| 69  | Mass 69 relative abundance         | 23.1                 |
| 70  | Less than 2.0% of mass 69          | 0.1 ( 0.4 ) 1        |
| 197 | Less than 2.0% of mass 198         | 0.0                  |
| 198 | Base Peak, 100% relative abundance | 100                  |
| 199 | 5.0 to 9.0% of mass 198            | 5.2                  |
| 365 | Greater than 1% of mass 198        | 3.5                  |
| 441 | Present, but less than mass 443    | 15.6                 |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 19.2 ( 19.2 ) 2      |

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|-------------------|------------------|----------------|------------------|------------------|
| SSTDICC2.5        | SSTDICC2.5       | BP025392.D     | 08/14/2025       | 12:33            |
| SSTDICC005        | SSTDICC005       | BP025393.D     | 08/14/2025       | 13:15            |
| SSTDICC010        | SSTDICC010       | BP025394.D     | 08/14/2025       | 13:56            |
| SSTDICC020        | SSTDICC020       | BP025395.D     | 08/14/2025       | 14:38            |
| SSTDICCC040       | SSTDICCC040      | BP025396.D     | 08/14/2025       | 15:19            |
| SSTDICC050        | SSTDICC050       | BP025397.D     | 08/14/2025       | 16:01            |
| SSTDICC060        | SSTDICC060       | BP025398.D     | 08/14/2025       | 16:42            |
| SSTDICC080        | SSTDICC080       | BP025399.D     | 08/14/2025       | 17:24            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BP025543.D  
Instrument ID: BNA\_P

Contract: AECO02  
SDG NO.: Q2924  
DFTPP Injection Date: 08/22/2025  
DFTPP Injection Time: 10:08

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.5 ( 1.6 ) 1        |
| 69  | Mass 69 relative abundance         | 32.9                 |
| 70  | Less than 2.0% of mass 69          | 0.1 ( 0.4 ) 1        |
| 197 | Less than 2.0% of mass 198         | 0.2                  |
| 198 | Base Peak, 100% relative abundance | 100                  |
| 199 | 5.0 to 9.0% of mass 198            | 6.8                  |
| 365 | Greater than 1% of mass 198        | 3.8                  |
| 441 | Present, but less than mass 443    | 15.5                 |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 19.7 ( 19.7 ) 2      |

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|-------------------|------------------|----------------|------------------|------------------|
| SSTDCCC040        | SSTDCCC040       | BP025544.D     | 08/22/2025       | 10:49            |
| MW-8A-082025      | Q2924-10         | BP025552.D     | 08/22/2025       | 16:58            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance

Contract: AECO02

Lab Code: ACE

SDG NO.: Q2924

Lab File ID: BP025560.D

DFTPP Injection Date: 08/25/2025

Instrument ID: BNA\_P

DFTPP Injection Time: 10:24

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.5 ( 1.7 ) 1        |
| 69  | Mass 69 relative abundance         | 30.5                 |
| 70  | Less than 2.0% of mass 69          | 0.1 ( 0.3 ) 1        |
| 197 | Less than 2.0% of mass 198         | 0.0                  |
| 198 | Base Peak, 100% relative abundance | 100                  |
| 199 | 5.0 to 9.0% of mass 198            | 6.3                  |
| 365 | Greater than 1% of mass 198        | 3.6                  |
| 441 | Present, but less than mass 443    | 15.6                 |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 19.3 ( 19.3 ) 2      |

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|-------------------|------------------|----------------|------------------|------------------|
| SSTDCCC040        | SSTDCCC040       | BP025561.D     | 08/25/2025       | 12:13            |
| PB169362BL        | PB169362BL       | BP025562.D     | 08/25/2025       | 12:54            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BP025578.D  
Instrument ID: BNA\_P

Contract: AECO02  
SDG NO.: Q2924  
DFTPP Injection Date: 08/26/2025  
DFTPP Injection Time: 08:59

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.5 ( 1.6 ) 1        |
| 69  | Mass 69 relative abundance         | 28.9                 |
| 70  | Less than 2.0% of mass 69          | 0.2 ( 0.6 ) 1        |
| 197 | Less than 2.0% of mass 198         | 0.0                  |
| 198 | Base Peak, 100% relative abundance | 100                  |
| 199 | 5.0 to 9.0% of mass 198            | 6                    |
| 365 | Greater than 1% of mass 198        | 3.4                  |
| 441 | Present, but less than mass 443    | 15.3                 |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 19.6 ( 19.6 ) 2      |

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|-------------------|------------------|----------------|------------------|------------------|
| SSTDCCC040        | SSTDCCC040       | BP025579.D     | 08/26/2025       | 10:21            |
| PB169362BS        | PB169362BS       | BP025581.D     | 08/26/2025       | 11:44            |
| MW-2B-082025DL    | Q2924-05DL       | BP025583.D     | 08/26/2025       | 13:10            |
| MW-2B-082025DL2   | Q2924-05DL2      | BP025586.D     | 08/26/2025       | 15:14            |
| MW-16A-082025     | Q2924-08         | BP025588.D     | 08/26/2025       | 16:37            |
| MW-16C-082025     | Q2924-09         | BP025589.D     | 08/26/2025       | 17:18            |
| MW-2A-082025      | Q2924-06         | BP025590.D     | 08/26/2025       | 17:59            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BP025595.D  
Instrument ID: BNA\_P

Contract: AECO02  
SDG NO.: Q2924  
DFTPP Injection Date: 08/27/2025  
DFTPP Injection Time: 08:58

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.5 ( 1.7 ) 1        |
| 69  | Mass 69 relative abundance         | 31.5                 |
| 70  | Less than 2.0% of mass 69          | 0.2 ( 0.6 ) 1        |
| 197 | Less than 2.0% of mass 198         | 0.0                  |
| 198 | Base Peak, 100% relative abundance | 100                  |
| 199 | 5.0 to 9.0% of mass 198            | 6.5                  |
| 365 | Greater than 1% of mass 198        | 3.6                  |
| 441 | Present, but less than mass 443    | 15.5                 |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 19.6 ( 19.6 ) 2      |

1-Value is % mass 69

2-Value is % mass 442

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED | TIME<br>ANALYZED |
|-------------------|------------------|----------------|------------------|------------------|
| SSTDCCC040        | SSTDCCC040       | BP025596.D     | 08/27/2025       | 09:39            |
| MW-201-082025     | Q2924-02         | BP025601.D     | 08/27/2025       | 13:37            |
| MW-201-082025MS   | Q2924-03MS       | BP025602.D     | 08/27/2025       | 14:18            |
| MW-201-082025MSD  | Q2924-04MSD      | BP025603.D     | 08/27/2025       | 15:00            |
| MW-201-082025DL   | Q2924-02DL       | BP025607.D     | 08/27/2025       | 17:44            |
| MW-201-082025DL2  | Q2924-02DL2      | BP025608.D     | 08/27/2025       | 18:26            |

8B

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q2924

Client ID : SSTDCCC040

Date Analyzed: 08/22/2025

Lab File ID: BF143563.D

Time Analyzed: 10:29

Instrument ID: BNA\_F

GC Column: DB-UI ID: 0.18 (mm)

|                  | IS1 (DCB)<br>AREA # | RT #  | IS2 (NPT)<br>AREA # | RT # | IS3 (ANT)<br>AREA # | RT #   |
|------------------|---------------------|-------|---------------------|------|---------------------|--------|
| 12 HOUR STD      | 94271               | 6.928 | 366222              | 8.21 | 205977              | 9.96   |
| UPPER LIMIT      | 188542              | 7.428 | 732444              | 8.71 | 411954              | 10.463 |
| LOWER LIMIT      | 47135.5             | 6.428 | 183111              | 7.71 | 102989              | 9.463  |
| EPA SAMPLE NO.   |                     |       |                     |      |                     |        |
| 01 MW-10C-082025 | 104622              | 6.92  | 416287              | 8.20 | 252150              | 9.96   |
| 02 MW-2B-082025  | 97960               | 6.93  | 310044              | 8.23 | 220938              | 9.96   |

IS1 (DCB) = 1,4-Dichlorobenzene-d4

IS2 (NPT) = Naphthalene-d8

IS3 (ANT) = Acenaphthene-d10

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8C

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q2924

Client ID: SSTDCCC040

Date Analyzed: 08/22/2025

Lab File ID: BF143563.D

Time Analyzed: 10:29

Instrument ID: BNA\_F

GC Column: DB-UI ID: 0.18 (mm)

|                  | IS4 (PHN)<br>AREA # | RT #   | IS5 (CRY)<br>AREA # | RT #   | IS6 (PRY)<br>AREA # | RT #   |
|------------------|---------------------|--------|---------------------|--------|---------------------|--------|
| 12 HOUR STD      | 328132              | 11.451 | 185804              | 14.098 | 191413              | 15.598 |
| UPPER LIMIT      | 656264              | 11.951 | 371608              | 14.598 | 382826              | 16.098 |
| LOWER LIMIT      | 164066              | 10.951 | 92902               | 13.598 | 95706.5             | 15.098 |
| EPA SAMPLE NO.   |                     |        |                     |        |                     |        |
| 01 MW-10C-082025 | 422746              | 11.45  | 233264              | 14.09  | 203920              | 15.59  |
| 02 MW-2B-082025  | 347033              | 11.45  | 201002              | 14.09  | 183367              | 15.59  |

IS4 (PHN) = Phenanthrene-d10

IS5 (CRY) = Chrysene-d12

IS6 (PRY) = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8B

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q2924

Client ID : SSTDCCC040

Date Analyzed: 08/27/2025

Lab File ID: BF143596.D

Time Analyzed: 09:16

Instrument ID: BNA\_F

GC Column: DB-UI ID: 0.18 (mm)

|                  | IS1 (DCB)<br>AREA # | RT #  | IS2 (NPT)<br>AREA # | RT #  | IS3 (ANT)<br>AREA # | RT #   |
|------------------|---------------------|-------|---------------------|-------|---------------------|--------|
| 12 HOUR STD      | 115923              | 6.893 | 446099              | 8.18  | 249607              | 9.93   |
| UPPER LIMIT      | 231846              | 7.393 | 892198              | 8.675 | 499214              | 10.428 |
| LOWER LIMIT      | 57961.5             | 6.393 | 223050              | 7.675 | 124804              | 9.428  |
| EPA SAMPLE NO.   |                     |       |                     |       |                     |        |
| 01 MW-18C-082025 | 107445              | 6.89  | 403245              | 8.17  | 213409              | 9.92   |
| 02 FB-01-082025  | 111189              | 6.89  | 428717              | 8.17  | 235375              | 9.92   |

IS1 (DCB) = 1,4-Dichlorobenzene-d4

IS2 (NPT) = Naphthalene-d8

IS3 (ANT) = Acenaphthene-d10

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8C

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q2924

Client ID: SSTDCCC040

Date Analyzed: 08/27/2025

Lab File ID: BF143596.D

Time Analyzed: 09:16

Instrument ID: BNA\_F

GC Column: DB-UI ID: 0.18 (mm)

|                  | IS4 (PHN)<br>AREA # | RT #   | IS5 (CRY)<br>AREA # | RT #   | IS6 (PRY)<br>AREA # | RT #   |
|------------------|---------------------|--------|---------------------|--------|---------------------|--------|
| 12 HOUR STD      | 402587              | 11.416 | 235836              | 14.051 | 227679              | 15.521 |
| UPPER LIMIT      | 805174              | 11.916 | 471672              | 14.551 | 455358              | 16.021 |
| LOWER LIMIT      | 201294              | 10.916 | 117918              | 13.551 | 113840              | 15.021 |
| EPA SAMPLE NO.   |                     |        |                     |        |                     |        |
| 01 MW-18C-082025 | 297699              | 11.41  | 161336              | 14.05  | 218565              | 15.52  |
| 02 FB-01-082025  | 345422              | 11.41  | 186452              | 14.05  | 227431              | 15.52  |

IS4 (PHN) = Phenanthrene-d10

IS5 (CRY) = Chrysene-d12

IS6 (PRY) = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8B

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q2924

Client ID : SSTDCCC040

Date Analyzed: 08/22/2025

Lab File ID: BP025544.D

Time Analyzed: 10:49

Instrument ID: BNA\_P

GC Column: ZB-GR ID: 0.25 (mm)

|                 | IS1 (DCB)<br>AREA # | RT #  | IS2 (NPT)<br>AREA # | RT #   | IS3 (ANT)<br>AREA # | RT #   |
|-----------------|---------------------|-------|---------------------|--------|---------------------|--------|
| 12 HOUR STD     | 137158              | 7.778 | 591885              | 10.55  | 400784              | 14.39  |
| UPPER LIMIT     | 274316              | 8.278 | 1183770             | 11.048 | 801568              | 14.889 |
| LOWER LIMIT     | 68579               | 7.278 | 295943              | 10.048 | 200392              | 13.889 |
| EPA SAMPLE NO.  |                     |       |                     |        |                     |        |
| 01 MW-8A-082025 | 178406              | 7.78  | 716017              | 10.54  | 464364              | 14.38  |

IS1 (DCB) = 1,4-Dichlorobenzene-d4

IS2 (NPT) = Naphthalene-d8

IS3 (ANT) = Acenaphthene-d10

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8C

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q2924

Client ID: SSTDCCC040

Date Analyzed: 08/22/2025

Lab File ID: BP025544.D

Time Analyzed: 10:49

Instrument ID: BNA\_P

GC Column: ZB-GR ID: 0.25 (mm)

|                 | IS4 (PHN)<br>AREA # | RT #   | IS5 (CRY)<br>AREA # | RT #   | IS6 (PRY)<br>AREA # | RT #   |
|-----------------|---------------------|--------|---------------------|--------|---------------------|--------|
| 12 HOUR STD     | 833863              | 17.189 | 897419              | 21.624 | 1034530             | 24.989 |
| UPPER LIMIT     | 1667730             | 17.689 | 1794840             | 22.124 | 2069060             | 25.489 |
| LOWER LIMIT     | 416932              | 16.689 | 448710              | 21.124 | 517265              | 24.489 |
| EPA SAMPLE NO.  |                     |        |                     |        |                     |        |
| 01 MW-8A-082025 | 972343              | 17.18  | 994558              | 21.63  | 1124630             | 24.99  |

IS4 (PHN) = Phenanthrene-d10

IS5 (CRY) = Chrysene-d12

IS6 (PRY) = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8B

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q2924

Client ID : SSTDCCC040

Date Analyzed: 08/25/2025

Lab File ID: BP025561.D

Time Analyzed: 12:13

Instrument ID: BNA\_P

GC Column: ZB-GR ID: 0.25 (mm)

|                | IS1 (DCB)<br>AREA # | RT #  | IS2 (NPT)<br>AREA # | RT #   | IS3 (ANT)<br>AREA # | RT #  |
|----------------|---------------------|-------|---------------------|--------|---------------------|-------|
| 12 HOUR STD    | 176050              | 7.766 | 715069              | 10.54  | 441846              | 14.39 |
| UPPER LIMIT    | 352100              | 8.266 | 1430140             | 11.037 | 883692              | 14.89 |
| LOWER LIMIT    | 88025               | 7.266 | 357535              | 10.037 | 220923              | 13.89 |
| EPA SAMPLE NO. |                     |       |                     |        |                     |       |
| 01 PB169362BL  | 173347              | 7.78  | 665603              | 10.54  | 424932              | 14.39 |

IS1 (DCB) = 1,4-Dichlorobenzene-d4

IS2 (NPT) = Naphthalene-d8

IS3 (ANT) = Acenaphthene-d10

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8C

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q2924

Client ID: SSTDCCC040

Date Analyzed: 08/25/2025

Lab File ID: BP025561.D

Time Analyzed: 12:13

Instrument ID: BNA\_P

GC Column: ZB-GR ID: 0.25 (mm)

|                | IS4 (PHN)<br>AREA # | RT #   | IS5 (CRY)<br>AREA # | RT #  | IS6 (PRY)<br>AREA # | RT #   |
|----------------|---------------------|--------|---------------------|-------|---------------------|--------|
| 12 HOUR STD    | 900532              | 17.189 | 945335              | 21.63 | 1072630             | 25.001 |
| UPPER LIMIT    | 1801060             | 17.689 | 1890670             | 22.13 | 2145260             | 25.501 |
| LOWER LIMIT    | 450266              | 16.689 | 472668              | 21.13 | 536315              | 24.501 |
| EPA SAMPLE NO. |                     |        |                     |       |                     |        |
| 01 PB169362BL  | 890876              | 17.19  | 1139400             | 21.63 | 1287990             | 25.00  |

IS4 (PHN) = Phenanthrene-d10

IS5 (CRY) = Chrysene-d12

IS6 (PRY) = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8B

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q2924

Client ID : SSTDCCC040

Date Analyzed: 08/26/2025

Lab File ID: BP025579.D

Time Analyzed: 10:21

Instrument ID: BNA\_P

GC Column: ZB-GR ID: 0.25 (mm)

|                    | IS1 (DCB)<br>AREA # | RT #  | IS2 (NPT)<br>AREA # | RT #   | IS3 (ANT)<br>AREA # | RT #  |
|--------------------|---------------------|-------|---------------------|--------|---------------------|-------|
| 12 HOUR STD        | 235051              | 7.778 | 932609              | 10.55  | 572065              | 14.39 |
| UPPER LIMIT        | 470102              | 8.278 | 1865220             | 11.049 | 1144130             | 14.89 |
| LOWER LIMIT        | 117526              | 7.278 | 466305              | 10.049 | 286033              | 13.89 |
| EPA SAMPLE NO.     |                     |       |                     |        |                     |       |
| 01 PB169362BS      | 256194              | 7.78  | 1004580             | 10.55  | 636349              | 14.41 |
| 02 MW-2B-082025DL  | 167152              | 7.77  | 645100              | 10.55  | 403608              | 14.39 |
| 03 MW-2B-082025DL2 | 173400              | 7.78  | 700981              | 10.54  | 444336              | 14.39 |
| 04 MW-2A-082025    | 209304              | 7.78  | 834432              | 10.54  | 540709              | 14.38 |
| 05 MW-16A-082025   | 194953              | 7.78  | 750992              | 10.55  | 492017              | 14.39 |
| 06 MW-16C-082025   | 200652              | 7.77  | 790738              | 10.55  | 502369              | 14.40 |

IS1 (DCB) = 1,4-Dichlorobenzene-d4

IS2 (NPT) = Naphthalene-d8

IS3 (ANT) = Acenaphthene-d10

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8C

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q2924

Client ID: SSTDCCC040

Date Analyzed: 08/26/2025

Lab File ID: BP025579.D

Time Analyzed: 10:21

Instrument ID: BNA\_P

GC Column: ZB-GR ID: 0.25 (mm)

|                    | IS4 (PHN)<br>AREA # | RT #   | IS5 (CRY)<br>AREA # | RT #   | IS6 (PRY)<br>AREA # | RT #   |
|--------------------|---------------------|--------|---------------------|--------|---------------------|--------|
| 12 HOUR STD        | 1130980             | 17.195 | 1110360             | 21.636 | 1292920             | 24.995 |
| UPPER LIMIT        | 2261960             | 17.695 | 2220720             | 22.136 | 2585840             | 25.495 |
| LOWER LIMIT        | 565490              | 16.695 | 555180              | 21.136 | 646460              | 24.495 |
| EPA SAMPLE NO.     |                     |        |                     |        |                     |        |
| 01 PB169362BS      | 1239650             | 17.20  | 1177570             | 21.63  | 1318340             | 24.99  |
| 02 MW-2B-082025DL  | 820492              | 17.19  | 922833              | 21.63  | 1102540             | 25.00  |
| 03 MW-2B-082025DL2 | 857309              | 17.19  | 880360              | 21.63  | 1060520             | 24.99  |
| 04 MW-2A-082025    | 1091600             | 17.19  | 1128820             | 21.64  | 1291270             | 25.00  |
| 05 MW-16A-082025   | 1024760             | 17.20  | 1121360             | 21.64  | 1313750             | 25.02  |
| 06 MW-16C-082025   | 1002260             | 17.20  | 1066650             | 21.63  | 1227490             | 25.01  |

IS4 (PHN) = Phenanthrene-d10

IS5 (CRY) = Chrysene-d12

IS6 (PRY) = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8B

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q2924

Client ID : SSTDCCC040

Date Analyzed: 08/27/2025

Lab File ID: BP025596.D

Time Analyzed: 09:39

Instrument ID: BNA\_P

GC Column: ZB-GR ID: 0.25 (mm)

|                     | IS1 (DCB)<br>AREA # | RT #  | IS2 (NPT)<br>AREA # | RT #   | IS3 (ANT)<br>AREA # | RT #   |
|---------------------|---------------------|-------|---------------------|--------|---------------------|--------|
| 12 HOUR STD         | 161611              | 7.778 | 681785              | 10.54  | 463868              | 14.40  |
| UPPER LIMIT         | 323222              | 8.278 | 1363570             | 11.043 | 927736              | 14.896 |
| LOWER LIMIT         | 80805.5             | 7.278 | 340893              | 10.043 | 231934              | 13.896 |
| EPA SAMPLE NO.      |                     |       |                     |        |                     |        |
| 01 MW-201-082025    | 185057              | 7.78  | 347077              | 10.57  | 427017              | 14.39  |
| 02 MW-201-082025MS  | 177033              | 7.78  | 389081              | 10.57  | 432672              | 14.39  |
| 03 MW-201-082025MSD | 165573              | 7.78  | 358938              | 10.57  | 417829              | 14.40  |
| 04 MW-201-082025DL  | 147420              | 7.77  | 575398              | 10.54  | 352604              | 14.39  |
| 05 MW-201-082025DL2 | 150824              | 7.77  | 581758              | 10.54  | 353815              | 14.39  |

IS1 (DCB) = 1,4-Dichlorobenzene-d4

IS2 (NPT) = Naphthalene-d8

IS3 (ANT) = Acenaphthene-d10

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.

8C

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q2924

Client ID: SSTDCCC040

Date Analyzed: 08/27/2025

Lab File ID: BP025596.D

Time Analyzed: 09:39

Instrument ID: BNA\_P

GC Column: ZB-GR ID: 0.25 (mm)

|                     | IS4 (PHN)<br>AREA # | RT #   | IS5 (CRY)<br>AREA # | RT #   | IS6 (PRY)<br>AREA # | RT #   |
|---------------------|---------------------|--------|---------------------|--------|---------------------|--------|
| 12 HOUR STD         | 978179              | 17.207 | 1036720             | 21.642 | 1177640             | 25.001 |
| UPPER LIMIT         | 1956360             | 17.707 | 2073440             | 22.142 | 2355280             | 25.501 |
| LOWER LIMIT         | 489090              | 16.707 | 518360              | 21.142 | 588820              | 24.501 |
| EPA SAMPLE NO.      |                     |        |                     |        |                     |        |
| 01 MW-201-082025    | 842894              | 17.20  | 908516              | 21.64  | 1056970             | 25.00  |
| 02 MW-201-082025MS  | 850262              | 17.19  | 903685              | 21.64  | 1056320             | 25.00  |
| 03 MW-201-082025MSD | 814298              | 17.20  | 887974              | 21.64  | 1057490             | 25.01  |
| 04 MW-201-082025DL  | 694303              | 17.19  | 837318              | 21.64  | 1049340             | 25.00  |
| 05 MW-201-082025DL2 | 659654              | 17.18  | 856697              | 21.64  | 1043580             | 25.01  |

IS4 (PHN) = Phenanthrene-d10

IS5 (CRY) = Chrysene-d12

IS6 (PRY) = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT UPPER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.



# QC SAMPLE DATA

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: |                  |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  |                  |      |
| Client Sample ID:  | PB169362BL                         |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | PB169362BL                         |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 1000                               | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025562.D        | 1         | 08/22/25 10:34 | 08/25/25 12:54 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 100-52-7       | Benzaldehyde                | 10.0  | U         | 3.90 | 10.0       | ug/L  |
| 108-95-2       | Phenol                      | 5.00  | U         | 0.91 | 5.00       | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 5.00  | U         | 0.81 | 5.00       | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 5.00  | U         | 0.58 | 5.00       | ug/L  |
| 95-48-7        | 2-Methylphenol              | 5.00  | U         | 1.10 | 5.00       | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 5.00  | U         | 1.30 | 5.00       | ug/L  |
| 98-86-2        | Acetophenone                | 5.00  | U         | 0.74 | 5.00       | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 10.0  | U         | 1.10 | 10.0       | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 2.50  | U         | 1.40 | 2.50       | ug/L  |
| 67-72-1        | Hexachloroethane            | 5.00  | U         | 0.65 | 5.00       | ug/L  |
| 98-95-3        | Nitrobenzene                | 5.00  | U         | 0.76 | 5.00       | ug/L  |
| 78-59-1        | Isophorone                  | 5.00  | U         | 0.75 | 5.00       | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 5.00  | U         | 1.80 | 5.00       | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 5.00  | U         | 1.90 | 5.00       | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 5.00  | U         | 0.68 | 5.00       | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 5.00  | U         | 0.52 | 5.00       | ug/L  |
| 91-20-3        | Naphthalene                 | 5.00  | U         | 0.50 | 5.00       | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 5.00  | U         | 0.84 | 5.00       | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 5.00  | U         | 0.54 | 5.00       | ug/L  |
| 105-60-2       | Caprolactam                 | 10.0  | U         | 1.10 | 10.0       | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 5.00  | U         | 0.59 | 5.00       | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 5.00  | U         | 0.56 | 5.00       | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 10.0  | U         | 3.60 | 10.0       | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 5.00  | U         | 0.51 | 5.00       | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 5.00  | U         | 0.62 | 5.00       | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 5.00  | U         | 0.53 | 5.00       | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 5.00  | U         | 0.61 | 5.00       | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 5.00  | U         | 1.30 | 5.00       | ug/L  |
| 131-11-3       | Dimethylphthalate           | 5.00  | U         | 0.61 | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |
|--------------------|------------------------------------|------------------|-----------------|------------------|
| Client:            | AECOM                              |                  | Date Collected: |                  |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  |                  |
| Client Sample ID:  | PB169362BL                         |                  | SDG No.:        | Q2924            |
| Lab Sample ID:     | PB169362BL                         |                  | Matrix:         | Water            |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |
| Sample Wt/Vol:     | 1000                               | Units: mL        | Final Vol:      | 1000 uL          |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N PH :           |
| Prep Method :      |                                    |                  |                 |                  |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025562.D        | 1         | 08/22/25 10:34 | 08/25/25 12:54 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 5.00  | U         | 0.75 | 5.00       | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 5.00  | U         | 0.92 | 5.00       | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 5.00  | U         | 1.10 | 5.00       | ug/L  |
| 83-32-9    | Acenaphthene               | 5.00  | U         | 0.55 | 5.00       | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 10.0  | U         | 6.00 | 10.0       | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 10.0  | U         | 2.40 | 10.0       | ug/L  |
| 132-64-9   | Dibenzofuran               | 5.00  | U         | 0.61 | 5.00       | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 5.00  | U         | 1.20 | 5.00       | ug/L  |
| 84-66-2    | Diethylphthalate           | 5.00  | U         | 0.69 | 5.00       | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 5.00  | U         | 0.68 | 5.00       | ug/L  |
| 86-73-7    | Fluorene                   | 5.00  | U         | 0.63 | 5.00       | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 5.00  | U         | 1.50 | 5.00       | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 10.0  | U         | 2.90 | 10.0       | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 5.00  | U         | 0.58 | 5.00       | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 5.00  | U         | 0.40 | 5.00       | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 5.00  | U         | 0.52 | 5.00       | ug/L  |
| 1912-24-9  | Atrazine                   | 5.00  | U         | 1.00 | 5.00       | ug/L  |
| 87-86-5    | Pentachlorophenol          | 10.0  | U         | 1.60 | 10.0       | ug/L  |
| 85-01-8    | Phenanthrene               | 5.00  | U         | 0.50 | 5.00       | ug/L  |
| 120-12-7   | Anthracene                 | 5.00  | U         | 0.61 | 5.00       | ug/L  |
| 86-74-8    | Carbazole                  | 5.00  | U         | 0.72 | 5.00       | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 5.00  | U         | 1.20 | 5.00       | ug/L  |
| 206-44-0   | Fluoranthene               | 5.00  | U         | 0.82 | 5.00       | ug/L  |
| 129-00-0   | Pyrene                     | 5.00  | U         | 0.50 | 5.00       | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 5.00  | U         | 1.90 | 5.00       | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 10.0  | U         | 0.93 | 10.0       | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 5.00  | U         | 0.45 | 5.00       | ug/L  |
| 218-01-9   | Chrysene                   | 5.00  | U         | 0.44 | 5.00       | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 5.00  | U         | 1.60 | 5.00       | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 10.0  | U         | 2.30 | 10.0       | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 5.00  | U         | 0.49 | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: |                  |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  |                  |      |
| Client Sample ID:  | PB169362BL                         |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | PB169362BL                         |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 1000                               | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025562.D        | 1         | 08/22/25 10:34 | 08/25/25 12:54 | PB169362      |

| CAS Number                            | Parameter                        | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------------------|----------------------------------|---------|-----------|----------|------------|----------|
| 207-08-9                              | Benzo(k)fluoranthene             | 5.00    | U         | 0.48     | 5.00       | ug/L     |
| 50-32-8                               | Benzo(a)pyrene                   | 5.00    | U         | 0.55     | 5.00       | ug/L     |
| 193-39-5                              | Indeno(1,2,3-cd)pyrene           | 5.00    | U         | 0.59     | 5.00       | ug/L     |
| 53-70-3                               | Dibenzo(a,h)anthracene           | 5.00    | U         | 0.67     | 5.00       | ug/L     |
| 191-24-2                              | Benzo(g,h,i)perylene             | 5.00    | U         | 0.69     | 5.00       | ug/L     |
| 95-94-3                               | 1,2,4,5-Tetrachlorobenzene       | 5.00    | U         | 0.52     | 5.00       | ug/L     |
| 123-91-1                              | 1,4-Dioxane                      | 5.00    | U         | 1.00     | 5.00       | ug/L     |
| 58-90-2                               | 2,3,4,6-Tetrachlorophenol        | 5.00    | U         | 0.72     | 5.00       | ug/L     |
| <b>SURROGATES</b>                     |                                  |         |           |          |            |          |
| 367-12-4                              | 2-Fluorophenol                   | 125     |           | 23 - 138 | 83%        | SPK: 150 |
| 13127-88-3                            | Phenol-d6                        | 120     |           | 10 - 134 | 80%        | SPK: 150 |
| 4165-60-0                             | Nitrobenzene-d5                  | 75.8    |           | 67 - 132 | 76%        | SPK: 100 |
| 321-60-8                              | 2-Fluorobiphenyl                 | 73.3    |           | 52 - 132 | 73%        | SPK: 100 |
| 118-79-6                              | 2,4,6-Tribromophenol             | 123     |           | 44 - 137 | 82%        | SPK: 150 |
| 1718-51-0                             | Terphenyl-d14                    | 70.1    |           | 42 - 152 | 70%        | SPK: 100 |
| <b>INTERNAL STANDARDS</b>             |                                  |         |           |          |            |          |
| 3855-82-1                             | 1,4-Dichlorobenzene-d4           | 173000  | 7.778     |          |            |          |
| 1146-65-2                             | Naphthalene-d8                   | 666000  | 10.543    |          |            |          |
| 15067-26-2                            | Acenaphthene-d10                 | 425000  | 14.389    |          |            |          |
| 1517-22-2                             | Phenanthrene-d10                 | 891000  | 17.189    |          |            |          |
| 1719-03-5                             | Chrysene-d12                     | 1140000 | 21.63     |          |            |          |
| 1520-96-3                             | Perylene-d12                     | 1290000 | 25.001    |          |            |          |
| <b>TENTATIVE IDENTIFIED COMPOUNDS</b> |                                  |         |           |          |            |          |
| 000123-42-2                           | 2-Pentanone, 4-hydroxy-4-methyl- | 7.70    | A         |          | 4.94       | ug/L     |

## Report of Analysis

|                    |                                    |                  |                 |                  |
|--------------------|------------------------------------|------------------|-----------------|------------------|
| Client:            | AECOM                              |                  | Date Collected: |                  |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  |                  |
| Client Sample ID:  | PB169362BL                         |                  | SDG No.:        | Q2924            |
| Lab Sample ID:     | PB169362BL                         |                  | Matrix:         | Water            |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |
| Sample Wt/Vol:     | 1000                               | Units: mL        | Final Vol:      | 1000 uL          |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N PH :           |
| Prep Method :      |                                    |                  |                 |                  |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025562.D        | 1         | 08/22/25 10:34 | 08/25/25 12:54 | PB169362      |

| CAS Number | Parameter | Conc. | Qualifier | MDL | LOQ / CRQL | Units |
|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

U = Not Detected  
 LOQ = Limit of Quantitation  
 MDL = Method Detection Limit  
 LOD = Limit of Detection  
 E = Value Exceeds Calibration Range  
 Q = indicates LCS control criteria did not meet requirements  
 M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value  
 B = Analyte Found in Associated Method Blank  
 N = Presumptive Evidence of a Compound  
 \* = Values outside of QC limits  
 D = Dilution  
 () = Laboratory InHouse Limit  
 A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: |                  |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  |                  |      |
| Client Sample ID:  | PB169362BS                         |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | PB169362BS                         |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 1000                               | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025581.D        | 1         | 08/22/25 10:34 | 08/26/25 11:44 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 100-52-7       | Benzaldehyde                | 22.9  |           | 3.90 | 10.0       | ug/L  |
| 108-95-2       | Phenol                      | 41.7  |           | 0.91 | 5.00       | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 40.1  |           | 0.81 | 5.00       | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 41.6  |           | 0.58 | 5.00       | ug/L  |
| 95-48-7        | 2-Methylphenol              | 41.0  |           | 1.10 | 5.00       | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 39.7  |           | 1.30 | 5.00       | ug/L  |
| 98-86-2        | Acetophenone                | 39.8  |           | 0.74 | 5.00       | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 39.8  |           | 1.10 | 10.0       | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 36.5  |           | 1.40 | 2.50       | ug/L  |
| 67-72-1        | Hexachloroethane            | 40.3  |           | 0.65 | 5.00       | ug/L  |
| 98-95-3        | Nitrobenzene                | 41.1  |           | 0.76 | 5.00       | ug/L  |
| 78-59-1        | Isophorone                  | 39.2  |           | 0.75 | 5.00       | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 42.7  |           | 1.80 | 5.00       | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 42.3  |           | 1.90 | 5.00       | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 40.1  |           | 0.68 | 5.00       | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 43.3  |           | 0.52 | 5.00       | ug/L  |
| 91-20-3        | Naphthalene                 | 39.7  |           | 0.50 | 5.00       | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 23.9  |           | 0.84 | 5.00       | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 40.8  |           | 0.54 | 5.00       | ug/L  |
| 105-60-2       | Caprolactam                 | 39.6  |           | 1.10 | 10.0       | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 42.3  |           | 0.59 | 5.00       | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 39.4  |           | 0.56 | 5.00       | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 81.0  | E         | 3.60 | 10.0       | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 44.1  |           | 0.51 | 5.00       | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 44.6  |           | 0.62 | 5.00       | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 41.8  |           | 0.53 | 5.00       | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 40.9  |           | 0.61 | 5.00       | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 44.0  |           | 1.30 | 5.00       | ug/L  |
| 131-11-3       | Dimethylphthalate           | 41.0  |           | 0.61 | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: |                  |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  |                  |      |
| Client Sample ID:  | PB169362BS                         |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | PB169362BS                         |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 1000                               | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025581.D        | 1         | 08/22/25 10:34 | 08/26/25 11:44 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 40.9  |           | 0.75 | 5.00       | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 43.4  |           | 0.92 | 5.00       | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 31.2  |           | 1.10 | 5.00       | ug/L  |
| 83-32-9    | Acenaphthene               | 39.5  |           | 0.55 | 5.00       | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 100   | E         | 6.00 | 10.0       | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 82.1  | E         | 2.40 | 10.0       | ug/L  |
| 132-64-9   | Dibenzofuran               | 40.2  |           | 0.61 | 5.00       | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 43.7  |           | 1.20 | 5.00       | ug/L  |
| 84-66-2    | Diethylphthalate           | 41.5  |           | 0.69 | 5.00       | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 40.1  |           | 0.68 | 5.00       | ug/L  |
| 86-73-7    | Fluorene                   | 40.2  |           | 0.63 | 5.00       | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 41.1  |           | 1.50 | 5.00       | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 46.7  |           | 2.90 | 10.0       | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 42.8  |           | 0.58 | 5.00       | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 42.2  |           | 0.40 | 5.00       | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 42.0  |           | 0.52 | 5.00       | ug/L  |
| 1912-24-9  | Atrazine                   | 41.7  |           | 1.00 | 5.00       | ug/L  |
| 87-86-5    | Pentachlorophenol          | 86.9  | E         | 1.60 | 10.0       | ug/L  |
| 85-01-8    | Phenanthrene               | 40.9  |           | 0.50 | 5.00       | ug/L  |
| 120-12-7   | Anthracene                 | 41.4  |           | 0.61 | 5.00       | ug/L  |
| 86-74-8    | Carbazole                  | 40.8  |           | 0.72 | 5.00       | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 41.6  |           | 1.20 | 5.00       | ug/L  |
| 206-44-0   | Fluoranthene               | 40.0  |           | 0.82 | 5.00       | ug/L  |
| 129-00-0   | Pyrene                     | 43.4  |           | 0.50 | 5.00       | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 44.3  |           | 1.90 | 5.00       | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 31.9  |           | 0.93 | 10.0       | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 42.2  |           | 0.45 | 5.00       | ug/L  |
| 218-01-9   | Chrysene                   | 41.7  |           | 0.44 | 5.00       | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 42.7  |           | 1.60 | 5.00       | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 43.7  |           | 2.30 | 10.0       | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 41.7  |           | 0.49 | 5.00       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: |                  |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  |                  |      |
| Client Sample ID:  | PB169362BS                         |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | PB169362BS                         |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 1000                               | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025581.D        | 1         | 08/22/25 10:34 | 08/26/25 11:44 | PB169362      |

| CAS Number                | Parameter                  | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------|----------------------------|---------|-----------|----------|------------|----------|
| 207-08-9                  | Benzo(k)fluoranthene       | 41.8    |           | 0.48     | 5.00       | ug/L     |
| 50-32-8                   | Benzo(a)pyrene             | 41.6    |           | 0.55     | 5.00       | ug/L     |
| 193-39-5                  | Indeno(1,2,3-cd)pyrene     | 41.4    |           | 0.59     | 5.00       | ug/L     |
| 53-70-3                   | Dibenzo(a,h)anthracene     | 41.8    |           | 0.67     | 5.00       | ug/L     |
| 191-24-2                  | Benzo(g,h,i)perylene       | 41.4    |           | 0.69     | 5.00       | ug/L     |
| 95-94-3                   | 1,2,4,5-Tetrachlorobenzene | 42.0    |           | 0.52     | 5.00       | ug/L     |
| 123-91-1                  | 1,4-Dioxane                | 34.9    |           | 1.00     | 5.00       | ug/L     |
| 58-90-2                   | 2,3,4,6-Tetrachlorophenol  | 43.0    |           | 0.72     | 5.00       | ug/L     |
| <b>SURROGATES</b>         |                            |         |           |          |            |          |
| 367-12-4                  | 2-Fluorophenol             | 116     |           | 23 - 138 | 77%        | SPK: 150 |
| 13127-88-3                | Phenol-d6                  | 113     |           | 10 - 134 | 75%        | SPK: 150 |
| 4165-60-0                 | Nitrobenzene-d5            | 72.1    |           | 67 - 132 | 72%        | SPK: 100 |
| 321-60-8                  | 2-Fluorobiphenyl           | 70.9    |           | 52 - 132 | 71%        | SPK: 100 |
| 118-79-6                  | 2,4,6-Tribromophenol       | 119     |           | 44 - 137 | 80%        | SPK: 150 |
| 1718-51-0                 | Terphenyl-d14              | 75.5    |           | 42 - 152 | 75%        | SPK: 100 |
| <b>INTERNAL STANDARDS</b> |                            |         |           |          |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4     | 256000  | 7.778     |          |            |          |
| 1146-65-2                 | Naphthalene-d8             | 1000000 | 10.549    |          |            |          |
| 15067-26-2                | Acenaphthene-d10           | 636000  | 14.407    |          |            |          |
| 1517-22-2                 | Phenanthrene-d10           | 1240000 | 17.195    |          |            |          |
| 1719-03-5                 | Chrysene-d12               | 1180000 | 21.63     |          |            |          |
| 1520-96-3                 | Perylene-d12               | 1320000 | 24.989    |          |            |          |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-201-082025MS                    |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-03MS                         |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 970                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025602.D        | 1         | 08/22/25 10:34 | 08/27/25 14:18 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 100-52-7       | Benzaldehyde                | 38.4  |           | 4.00 | 10.3       | ug/L  |
| 108-95-2       | Phenol                      | 16.0  |           | 0.94 | 5.20       | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 44.3  |           | 0.84 | 5.20       | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 39.6  |           | 0.60 | 5.20       | ug/L  |
| 95-48-7        | 2-Methylphenol              | 33.8  |           | 1.20 | 5.20       | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 17.1  |           | 1.30 | 5.20       | ug/L  |
| 98-86-2        | Acetophenone                | 87.5  | E         | 0.76 | 5.20       | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 40.9  |           | 1.10 | 10.3       | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 42.2  |           | 1.50 | 2.60       | ug/L  |
| 67-72-1        | Hexachloroethane            | 93.2  | E         | 0.67 | 5.20       | ug/L  |
| 98-95-3        | Nitrobenzene                | 85.8  | E         | 0.78 | 5.20       | ug/L  |
| 78-59-1        | Isophorone                  | 81.7  |           | 0.77 | 5.20       | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 86.6  | E         | 1.80 | 5.20       | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 75.6  |           | 1.90 | 5.20       | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 70.3  |           | 0.70 | 5.20       | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 80.6  |           | 0.54 | 5.20       | ug/L  |
| 91-20-3        | Naphthalene                 | 2700  | E         | 0.52 | 5.20       | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 28.5  |           | 0.87 | 5.20       | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 71.5  |           | 0.56 | 5.20       | ug/L  |
| 105-60-2       | Caprolactam                 | 16.7  |           | 1.20 | 10.3       | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 76.5  |           | 0.61 | 5.20       | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 800   | E         | 0.58 | 5.20       | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 79.7  |           | 3.70 | 10.3       | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 50.2  |           | 0.53 | 5.20       | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 49.4  |           | 0.64 | 5.20       | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 69.6  |           | 0.55 | 5.20       | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 48.9  |           | 0.63 | 5.20       | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 54.1  |           | 1.30 | 5.20       | ug/L  |
| 131-11-3       | Dimethylphthalate           | 48.7  |           | 0.63 | 5.20       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-201-082025MS                    |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-03MS                         |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 970                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025602.D        | 1         | 08/22/25 10:34 | 08/27/25 14:18 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 190   | E         | 0.77 | 5.20       | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 51.4  |           | 0.95 | 5.20       | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 31.7  |           | 1.10 | 5.20       | ug/L  |
| 83-32-9    | Acenaphthene               | 54.7  |           | 0.57 | 5.20       | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 120   | E         | 6.20 | 10.3       | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 42.0  |           | 2.50 | 10.3       | ug/L  |
| 132-64-9   | Dibenzofuran               | 51.3  |           | 0.63 | 5.20       | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 53.3  |           | 1.30 | 5.20       | ug/L  |
| 84-66-2    | Diethylphthalate           | 50.3  |           | 0.71 | 5.20       | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 48.5  |           | 0.70 | 5.20       | ug/L  |
| 86-73-7    | Fluorene                   | 67.1  |           | 0.65 | 5.20       | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 48.5  |           | 1.50 | 5.20       | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 55.1  |           | 3.00 | 10.3       | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 51.1  |           | 0.60 | 5.20       | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 51.5  |           | 0.41 | 5.20       | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 52.5  |           | 0.54 | 5.20       | ug/L  |
| 1912-24-9  | Atrazine                   | 47.8  |           | 1.00 | 5.20       | ug/L  |
| 87-86-5    | Pentachlorophenol          | 110   | E         | 1.60 | 10.3       | ug/L  |
| 85-01-8    | Phenanthrene               | 65.4  |           | 0.52 | 5.20       | ug/L  |
| 120-12-7   | Anthracene                 | 53.4  |           | 0.63 | 5.20       | ug/L  |
| 86-74-8    | Carbazole                  | 61.7  |           | 0.74 | 5.20       | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 51.3  |           | 1.30 | 5.20       | ug/L  |
| 206-44-0   | Fluoranthene               | 51.0  |           | 0.85 | 5.20       | ug/L  |
| 129-00-0   | Pyrene                     | 50.5  |           | 0.52 | 5.20       | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 51.2  |           | 2.00 | 5.20       | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 41.5  |           | 0.96 | 10.3       | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 49.5  |           | 0.46 | 5.20       | ug/L  |
| 218-01-9   | Chrysene                   | 50.0  |           | 0.45 | 5.20       | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 48.4  |           | 1.60 | 5.20       | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 51.0  |           | 2.40 | 10.3       | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 50.0  |           | 0.51 | 5.20       | ug/L  |

## Report of Analysis

|                    |                                    |                 |                  |
|--------------------|------------------------------------|-----------------|------------------|
| Client:            | AECOM                              | Date Collected: | 08/20/25         |
| Project:           | National Grid Equity - Brooklyn NY | Date Received:  | 08/20/25         |
| Client Sample ID:  | MW-201-082025MS                    | SDG No.:        | Q2924            |
| Lab Sample ID:     | Q2924-03MS                         | Matrix:         | Water            |
| Analytical Method: | 8270E                              | % Solid:        | 0                |
| Sample Wt/Vol:     | 970 Units: mL                      | Final Vol:      | 1000 uL          |
| Soil Aliquot Vol:  | uL                                 | Test:           | SVOC-TCL BNA -20 |
| Extraction Type :  | Decanted : N                       | Level :         | LOW              |
| Injection Volume : | GPC Factor : 1.0                   | GPC Cleanup :   | N PH :           |
| Prep Method :      |                                    |                 |                  |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025602.D        | 1         | 08/22/25 10:34 | 08/27/25 14:18 | PB169362      |

| CAS Number                | Parameter                  | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------|----------------------------|---------|-----------|----------|------------|----------|
| 207-08-9                  | Benzo(k)fluoranthene       | 49.1    |           | 0.49     | 5.20       | ug/L     |
| 50-32-8                   | Benzo(a)pyrene             | 49.3    |           | 0.57     | 5.20       | ug/L     |
| 193-39-5                  | Indeno(1,2,3-cd)pyrene     | 50.1    |           | 0.61     | 5.20       | ug/L     |
| 53-70-3                   | Dibenzo(a,h)anthracene     | 51.0    |           | 0.69     | 5.20       | ug/L     |
| 191-24-2                  | Benzo(g,h,i)perylene       | 50.2    |           | 0.71     | 5.20       | ug/L     |
| 95-94-3                   | 1,2,4,5-Tetrachlorobenzene | 50.1    |           | 0.54     | 5.20       | ug/L     |
| 123-91-1                  | 1,4-Dioxane                | 21.2    |           | 1.00     | 5.20       | ug/L     |
| 58-90-2                   | 2,3,4,6-Tetrachlorophenol  | 51.1    |           | 0.74     | 5.20       | ug/L     |
| <b>SURROGATES</b>         |                            |         |           |          |            |          |
| 367-12-4                  | 2-Fluorophenol             | 67.6    |           | 23 - 138 | 45%        | SPK: 150 |
| 13127-88-3                | Phenol-d6                  | 41.4    |           | 10 - 134 | 28%        | SPK: 150 |
| 4165-60-0                 | Nitrobenzene-d5            | 148     | *         | 67 - 132 | 148%       | SPK: 100 |
| 321-60-8                  | 2-Fluorobiphenyl           | 79.6    |           | 52 - 132 | 80%        | SPK: 100 |
| 118-79-6                  | 2,4,6-Tribromophenol       | 138     |           | 44 - 137 | 92%        | SPK: 150 |
| 1718-51-0                 | Terphenyl-d14              | 76.9    |           | 42 - 152 | 77%        | SPK: 100 |
| <b>INTERNAL STANDARDS</b> |                            |         |           |          |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4     | 177000  |           | 7.784    |            |          |
| 1146-65-2                 | Naphthalene-d8             | 389000  |           | 10.566   |            |          |
| 15067-26-2                | Acenaphthene-d10           | 433000  |           | 14.389   |            |          |
| 1517-22-2                 | Phenanthrene-d10           | 850000  |           | 17.189   |            |          |
| 1719-03-5                 | Chrysene-d12               | 904000  |           | 21.636   |            |          |
| 1520-96-3                 | Perylene-d12               | 1060000 |           | 25.001   |            |          |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-201-082025MSD                   |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-04MSD                        |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 950                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025603.D        | 1         | 08/22/25 10:34 | 08/27/25 15:00 | PB169362      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 100-52-7       | Benzaldehyde                | 37.7  |           | 4.10 | 10.5       | ug/L  |
| 108-95-2       | Phenol                      | 16.4  |           | 0.96 | 5.30       | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 44.2  |           | 0.85 | 5.30       | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 39.6  |           | 0.61 | 5.30       | ug/L  |
| 95-48-7        | 2-Methylphenol              | 34.5  |           | 1.20 | 5.30       | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 18.4  |           | 1.30 | 5.30       | ug/L  |
| 98-86-2        | Acetophenone                | 90.6  | E         | 0.78 | 5.30       | ug/L  |
| 65794-96-9     | 3+4-Methylphenols           | 41.7  |           | 1.20 | 10.5       | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 43.2  |           | 1.50 | 2.60       | ug/L  |
| 67-72-1        | Hexachloroethane            | 96.9  | E         | 0.68 | 5.30       | ug/L  |
| 98-95-3        | Nitrobenzene                | 89.1  | E         | 0.80 | 5.30       | ug/L  |
| 78-59-1        | Isophorone                  | 85.0  | E         | 0.79 | 5.30       | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 90.4  | E         | 1.90 | 5.30       | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 78.4  |           | 1.90 | 5.30       | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 75.2  |           | 0.72 | 5.30       | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 85.1  | E         | 0.55 | 5.30       | ug/L  |
| 91-20-3        | Naphthalene                 | 3000  | E         | 0.53 | 5.30       | ug/L  |
| 106-47-8       | 4-Chloroaniline             | 31.3  |           | 0.88 | 5.30       | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 74.8  |           | 0.57 | 5.30       | ug/L  |
| 105-60-2       | Caprolactam                 | 17.0  |           | 1.20 | 10.5       | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 79.4  |           | 0.62 | 5.30       | ug/L  |
| 91-57-6        | 2-Methylnaphthalene         | 850   | E         | 0.59 | 5.30       | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 76.9  |           | 3.80 | 10.5       | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 50.2  |           | 0.54 | 5.30       | ug/L  |
| 95-95-4        | 2,4,5-Trichlorophenol       | 49.9  |           | 0.65 | 5.30       | ug/L  |
| 92-52-4        | 1,1-Biphenyl                | 70.0  |           | 0.56 | 5.30       | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 48.9  |           | 0.64 | 5.30       | ug/L  |
| 88-74-4        | 2-Nitroaniline              | 53.9  |           | 1.30 | 5.30       | ug/L  |
| 131-11-3       | Dimethylphthalate           | 47.8  |           | 0.64 | 5.30       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-201-082025MSD                   |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-04MSD                        |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 950                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025603.D        | 1         | 08/22/25 10:34 | 08/27/25 15:00 | PB169362      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 208-96-8   | Acenaphthylene             | 190   | E         | 0.79 | 5.30       | ug/L  |
| 606-20-2   | 2,6-Dinitrotoluene         | 51.3  |           | 0.97 | 5.30       | ug/L  |
| 99-09-2    | 3-Nitroaniline             | 32.4  |           | 1.10 | 5.30       | ug/L  |
| 83-32-9    | Acenaphthene               | 53.5  |           | 0.58 | 5.30       | ug/L  |
| 51-28-5    | 2,4-Dinitrophenol          | 120   | E         | 6.30 | 10.5       | ug/L  |
| 100-02-7   | 4-Nitrophenol              | 42.7  |           | 2.50 | 10.5       | ug/L  |
| 132-64-9   | Dibenzofuran               | 51.2  |           | 0.64 | 5.30       | ug/L  |
| 121-14-2   | 2,4-Dinitrotoluene         | 52.5  |           | 1.30 | 5.30       | ug/L  |
| 84-66-2    | Diethylphthalate           | 48.9  |           | 0.73 | 5.30       | ug/L  |
| 7005-72-3  | 4-Chlorophenyl-phenylether | 48.5  |           | 0.72 | 5.30       | ug/L  |
| 86-73-7    | Fluorene                   | 68.0  |           | 0.66 | 5.30       | ug/L  |
| 100-01-6   | 4-Nitroaniline             | 49.4  |           | 1.60 | 5.30       | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 55.7  |           | 3.00 | 10.5       | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 50.5  |           | 0.61 | 5.30       | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 52.0  |           | 0.42 | 5.30       | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 52.3  |           | 0.55 | 5.30       | ug/L  |
| 1912-24-9  | Atrazine                   | 48.5  |           | 1.10 | 5.30       | ug/L  |
| 87-86-5    | Pentachlorophenol          | 110   | E         | 1.70 | 10.5       | ug/L  |
| 85-01-8    | Phenanthrene               | 67.2  |           | 0.53 | 5.30       | ug/L  |
| 120-12-7   | Anthracene                 | 53.2  |           | 0.64 | 5.30       | ug/L  |
| 86-74-8    | Carbazole                  | 63.5  |           | 0.76 | 5.30       | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 52.6  |           | 1.30 | 5.30       | ug/L  |
| 206-44-0   | Fluoranthene               | 51.6  |           | 0.86 | 5.30       | ug/L  |
| 129-00-0   | Pyrene                     | 50.1  |           | 0.53 | 5.30       | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 52.0  |           | 2.00 | 5.30       | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 43.4  |           | 0.98 | 10.5       | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 49.9  |           | 0.47 | 5.30       | ug/L  |
| 218-01-9   | Chrysene                   | 51.2  |           | 0.46 | 5.30       | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 49.7  |           | 1.70 | 5.30       | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 52.0  |           | 2.50 | 10.5       | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 50.4  |           | 0.52 | 5.30       | ug/L  |

## Report of Analysis

|                    |                                    |                  |                 |                  |      |
|--------------------|------------------------------------|------------------|-----------------|------------------|------|
| Client:            | AECOM                              |                  | Date Collected: | 08/20/25         |      |
| Project:           | National Grid Equity - Brooklyn NY |                  | Date Received:  | 08/20/25         |      |
| Client Sample ID:  | MW-201-082025MSD                   |                  | SDG No.:        | Q2924            |      |
| Lab Sample ID:     | Q2924-04MSD                        |                  | Matrix:         | Water            |      |
| Analytical Method: | 8270E                              |                  | % Solid:        | 0                |      |
| Sample Wt/Vol:     | 950                                | Units: mL        | Final Vol:      | 1000             | uL   |
| Soil Aliquot Vol:  |                                    | uL               | Test:           | SVOC-TCL BNA -20 |      |
| Extraction Type :  |                                    | Decanted : N     | Level :         | LOW              |      |
| Injection Volume : |                                    | GPC Factor : 1.0 | GPC Cleanup :   | N                | PH : |
| Prep Method :      |                                    |                  |                 |                  |      |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BP025603.D        | 1         | 08/22/25 10:34 | 08/27/25 15:00 | PB169362      |

| CAS Number                | Parameter                  | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------|----------------------------|---------|-----------|----------|------------|----------|
| 207-08-9                  | Benzo(k)fluoranthene       | 49.6    |           | 0.51     | 5.30       | ug/L     |
| 50-32-8                   | Benzo(a)pyrene             | 49.8    |           | 0.58     | 5.30       | ug/L     |
| 193-39-5                  | Indeno(1,2,3-cd)pyrene     | 50.9    |           | 0.62     | 5.30       | ug/L     |
| 53-70-3                   | Dibenzo(a,h)anthracene     | 51.6    |           | 0.71     | 5.30       | ug/L     |
| 191-24-2                  | Benzo(g,h,i)perylene       | 50.6    |           | 0.73     | 5.30       | ug/L     |
| 95-94-3                   | 1,2,4,5-Tetrachlorobenzene | 49.8    |           | 0.55     | 5.30       | ug/L     |
| 123-91-1                  | 1,4-Dioxane                | 21.8    |           | 1.10     | 5.30       | ug/L     |
| 58-90-2                   | 2,3,4,6-Tetrachlorophenol  | 50.6    |           | 0.76     | 5.30       | ug/L     |
| <b>SURROGATES</b>         |                            |         |           |          |            |          |
| 367-12-4                  | 2-Fluorophenol             | 67.4    |           | 23 - 138 | 45%        | SPK: 150 |
| 13127-88-3                | Phenol-d6                  | 41.4    |           | 10 - 134 | 28%        | SPK: 150 |
| 4165-60-0                 | Nitrobenzene-d5            | 149     | *         | 67 - 132 | 149%       | SPK: 100 |
| 321-60-8                  | 2-Fluorobiphenyl           | 78.1    |           | 52 - 132 | 78%        | SPK: 100 |
| 118-79-6                  | 2,4,6-Tribromophenol       | 135     |           | 44 - 137 | 90%        | SPK: 150 |
| 1718-51-0                 | Terphenyl-d14              | 76.2    |           | 42 - 152 | 76%        | SPK: 100 |
| <b>INTERNAL STANDARDS</b> |                            |         |           |          |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4     | 166000  | 7.778     |          |            |          |
| 1146-65-2                 | Naphthalene-d8             | 359000  | 10.566    |          |            |          |
| 15067-26-2                | Acenaphthene-d10           | 418000  | 14.401    |          |            |          |
| 1517-22-2                 | Phenanthrene-d10           | 814000  | 17.201    |          |            |          |
| 1719-03-5                 | Chrysene-d12               | 888000  | 21.636    |          |            |          |
| 1520-96-3                 | Perylene-d12               | 1060000 | 25.007    |          |            |          |

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products



# CALIBRATION SUMMARY

Method Path : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\  
 Method File : 8270-BF082025.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 Last Update : Thu Aug 21 06:12:21 2025  
 Response Via : Initial Calibration

## Calibration Files

2.5 =BF143477.D 5 =BF143478.D 10 =BF143479.D 20 =BF143480.D 40 =BF143481.D 50 =BF143482.D 60 =BF143483.D 80 =BF143484.D

| Compound                   | 2.5            | 5     | 10    | 20    | 40    | 50    | 60    | 80    | Avg   | %RSD  |
|----------------------------|----------------|-------|-------|-------|-------|-------|-------|-------|-------|-------|
| -----                      |                |       |       |       |       |       |       |       |       |       |
| 1) I 1,4-Dichlorobenzen... | -----ISTD----- |       |       |       |       |       |       |       |       |       |
| 2) 1,4-Dioxane             |                | 0.528 | 0.529 | 0.539 | 0.536 | 0.519 | 0.542 | 0.522 | 0.531 | 1.62  |
| 3) Pyridine                |                | 1.339 | 1.392 | 1.435 | 1.438 | 1.389 | 1.477 | 1.421 | 1.413 | 3.13  |
| 4) n-Nitrosodimet...       |                |       | 0.604 | 0.619 | 0.641 | 0.633 | 0.666 | 0.648 | 0.635 | 3.44  |
| 5) S 2-Fluorophenol        |                | 1.203 | 1.198 | 1.194 | 1.138 | 1.087 | 1.137 | 1.061 | 1.145 | 4.92  |
| 6) Aniline                 |                | 2.034 | 1.969 | 1.959 | 1.890 | 1.800 | 1.864 | 1.714 | 1.890 | 5.78  |
| 7) S Phenol-d6             |                | 1.536 | 1.460 | 1.471 | 1.407 | 1.342 | 1.384 | 1.299 | 1.414 | 5.76  |
| 8) 2-Chlorophenol          |                | 1.301 | 1.284 | 1.302 | 1.268 | 1.229 | 1.270 | 1.202 | 1.265 | 2.94  |
| 9) Benzaldehyde            |                |       | 1.008 | 0.981 | 1.200 | 1.133 | 1.402 |       | 1.145 | 14.80 |
| 10) C Phenol               |                | 1.741 | 1.695 | 1.725 | 1.646 | 1.554 | 1.593 | 1.449 | 1.629 | 6.43  |
| 11) bis(2-Chloroet...      |                | 1.296 | 1.230 | 1.235 | 1.213 | 1.170 | 1.221 | 1.156 | 1.217 | 3.78  |
| 12) 1,3-Dichlorobe...      |                | 1.558 | 1.451 | 1.466 | 1.414 | 1.345 | 1.399 | 1.317 | 1.421 | 5.65  |
| 13) C 1,4-Dichlorobe...    |                | 1.531 | 1.487 | 1.483 | 1.411 | 1.355 | 1.409 | 1.304 | 1.426 | 5.62  |
| 14) 1,2-Dichlorobe...      |                | 1.448 | 1.413 | 1.386 | 1.348 | 1.289 | 1.333 | 1.233 | 1.350 | 5.45  |
| 15) Benzyl Alcohol         |                |       | 1.037 | 1.063 | 1.083 | 1.040 | 1.088 | 1.041 | 1.059 | 2.14  |
| 16) 2,2'-oxybis(1-...      |                | 2.144 | 2.104 | 2.053 | 1.954 | 1.805 | 1.868 | 1.703 | 1.947 | 8.40  |
| 17) 2-Methylphenol         |                | 1.052 | 1.042 | 1.056 | 1.011 | 0.975 | 1.011 | 0.960 | 1.015 | 3.67  |
| 18) Hexachloroethane       |                | 0.495 | 0.498 | 0.502 | 0.491 | 0.462 | 0.489 | 0.456 | 0.485 | 3.79  |
| 19) P n-Nitroso-di-n...    | 0.922          | 0.943 | 0.912 | 0.879 | 0.844 | 0.800 | 0.832 | 0.781 | 0.864 | 6.86  |
| 20) 3+4-Methylphenols      |                |       | 1.337 | 1.320 | 1.218 | 1.143 | 1.178 | 1.076 | 1.212 | 8.40  |
|                            |                |       |       |       |       |       |       |       |       |       |
| 21) I Naphthalene-d8       | -----ISTD----- |       |       |       |       |       |       |       |       |       |
| 22) Acetophenone           |                | 0.489 | 0.464 | 0.459 | 0.418 | 0.402 | 0.410 | 0.382 | 0.432 | 8.99  |
| 23) S Nitrobenzene-d5      |                | 0.354 | 0.345 | 0.357 | 0.344 | 0.335 | 0.340 | 0.325 | 0.343 | 3.18  |
| 24) Nitrobenzene           |                | 0.353 | 0.351 | 0.366 | 0.356 | 0.348 | 0.358 | 0.337 | 0.353 | 2.55  |
| 25) Isophorone             |                | 0.666 | 0.630 | 0.644 | 0.621 | 0.608 | 0.625 | 0.603 | 0.628 | 3.41  |
| 26) C 2-Nitrophenol        |                | 0.162 | 0.165 | 0.177 | 0.177 | 0.175 | 0.180 | 0.173 | 0.173 | 3.76  |
| 27) 2,4-Dimethylph...      |                | 0.283 | 0.272 | 0.281 | 0.270 | 0.262 | 0.265 | 0.251 | 0.269 | 4.11  |
| 28) bis(2-Chloroet...      |                | 0.409 | 0.399 | 0.401 | 0.386 | 0.371 | 0.379 | 0.357 | 0.386 | 4.74  |
| 29) C 2,4-Dichloroph...    |                | 0.293 | 0.299 | 0.303 | 0.288 | 0.281 | 0.283 | 0.266 | 0.288 | 4.30  |
| 30) 1,2,4-Trichlor...      |                | 0.318 | 0.300 | 0.314 | 0.292 | 0.282 | 0.291 | 0.274 | 0.296 | 5.46  |
| 31) Naphthalene            |                | 1.069 | 1.009 | 1.014 | 0.943 | 0.907 | 0.921 | 0.859 | 0.960 | 7.58  |
| 32) Benzoic acid           |                |       | 0.086 | 0.116 | 0.145 | 0.153 | 0.163 | 0.160 | 0.137 | 21.97 |
| 33) 4-Chloroaniline        |                | 0.359 | 0.345 | 0.362 | 0.343 | 0.331 | 0.330 | 0.315 | 0.341 | 4.87  |
| 34) C Hexachlorobuta...    |                | 0.207 | 0.199 | 0.201 | 0.192 | 0.186 | 0.190 | 0.177 | 0.193 | 5.21  |
| 35) Caprolactam            |                |       | 0.063 | 0.068 | 0.071 | 0.071 | 0.070 | 0.072 | 0.069 | 4.96  |
| 36) C 4-Chloro-3-met...    |                | 0.303 | 0.292 | 0.303 | 0.287 | 0.280 | 0.284 | 0.269 | 0.288 | 4.25  |
| 37) 2-Methylnaphth...      |                | 0.716 | 0.684 | 0.684 | 0.633 | 0.600 | 0.610 | 0.562 | 0.641 | 8.62  |
| 38) 1-Methylnaphth...      |                | 0.692 | 0.651 | 0.650 | 0.603 | 0.577 | 0.581 | 0.540 | 0.613 | 8.62  |

Method Path : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\  
 Method File : 8270-BF082025.M

|       |                   |                |       |       |       |       |       |       |       |       |  |
|-------|-------------------|----------------|-------|-------|-------|-------|-------|-------|-------|-------|--|
| 39) I | Acenaphthene-d10  | -----ISTD----- |       |       |       |       |       |       |       |       |  |
| 40)   | 1,2,4,5-Tetrac... | 0.629          | 0.584 | 0.613 | 0.565 | 0.544 | 0.552 | 0.517 | 0.572 | 6.88  |  |
| 41) P | Hexachlorocycl... |                | 0.098 | 0.153 | 0.198 | 0.205 | 0.226 | 0.222 | 0.184 | 26.99 |  |
| 42) S | 2,4,6-Tribromo... | 0.191          | 0.188 | 0.194 | 0.188 | 0.181 | 0.184 | 0.177 | 0.186 | 3.12  |  |
| 43) C | 2,4,6-Trichlor... | 0.338          | 0.347 | 0.364 | 0.364 | 0.351 | 0.360 | 0.339 | 0.352 | 3.18  |  |
| 44)   | 2,4,5-Trichlor... | 0.391          | 0.376 | 0.413 | 0.382 | 0.377 | 0.387 | 0.365 | 0.385 | 3.94  |  |
| 45) S | 2-Fluorobiphenyl  | 1.445          | 1.336 | 1.323 | 1.156 | 1.106 | 1.106 | 0.999 | 1.210 | 13.20 |  |
| 46)   | 1,1'-Biphenyl     | 1.582          | 1.505 | 1.519 | 1.399 | 1.335 | 1.349 | 1.247 | 1.420 | 8.43  |  |
| 47)   | 2-Chloronaphth... | 1.222          | 1.165 | 1.172 | 1.105 | 1.056 | 1.068 | 1.008 | 1.114 | 6.77  |  |
| 48)   | 2-Nitroaniline    | 0.314          | 0.313 | 0.338 | 0.337 | 0.329 | 0.331 | 0.320 | 0.326 | 3.19  |  |
| 49)   | Acenaphthylene    | 1.747          | 1.663 | 1.687 | 1.580 | 1.516 | 1.533 | 1.437 | 1.595 | 6.85  |  |
| 50)   | Dimethylphthalate | 1.293          | 1.219 | 1.252 | 1.201 | 1.159 | 1.175 | 1.133 | 1.205 | 4.58  |  |
| 51)   | 2,6-Dinitrotol... | 0.224          | 0.237 | 0.259 | 0.259 | 0.260 | 0.257 | 0.251 | 0.250 | 5.63  |  |
| 52) C | Acenaphthene      | 1.170          | 1.131 | 1.137 | 1.076 | 1.037 | 1.044 | 0.985 | 1.083 | 6.07  |  |
| 53)   | 3-Nitroaniline    | 0.260          | 0.257 | 0.280 | 0.278 | 0.273 | 0.271 | 0.269 | 0.270 | 3.21  |  |
| 54) P | 2,4-Dinitrophenol |                | 0.072 | 0.099 | 0.120 | 0.132 | 0.137 | 0.129 | 0.115 | 21.92 |  |
| 55)   | Dibenzofuran      | 1.717          | 1.625 | 1.647 | 1.523 | 1.461 | 1.453 | 1.381 | 1.544 | 7.91  |  |
| 56) P | 4-Nitrophenol     |                | 0.115 | 0.147 | 0.171 | 0.177 | 0.172 | 0.181 | 0.161 | 15.78 |  |
| 57)   | 2,4-Dinitrotol... | 0.292          | 0.318 | 0.349 | 0.351 | 0.344 | 0.344 | 0.336 | 0.333 | 6.38  |  |
| 58)   | Fluorene          | 1.367          | 1.286 | 1.270 | 1.158 | 1.111 | 1.102 | 1.025 | 1.188 | 10.25 |  |
| 59)   | 2,3,4,6-Tetrac... | 0.261          | 0.264 | 0.299 | 0.297 | 0.294 | 0.297 | 0.288 | 0.286 | 5.59  |  |
| 60)   | Diethylphthalate  | 1.172          | 1.097 | 1.110 | 1.084 | 1.053 | 1.036 | 1.008 | 1.080 | 5.01  |  |
| 61)   | 4-Chlorophenyl... | 0.671          | 0.624 | 0.628 | 0.585 | 0.558 | 0.567 | 0.527 | 0.594 | 8.30  |  |
| 62)   | 4-Nitroaniline    | 0.230          | 0.237 | 0.258 | 0.257 | 0.255 | 0.248 | 0.250 | 0.248 | 4.28  |  |
| 63)   | Azobenzene        | 1.221          | 1.154 | 1.177 | 1.122 | 1.080 | 1.067 | 1.018 | 1.120 | 6.26  |  |
| 64) I | Phenanthrene-d10  | -----ISTD----- |       |       |       |       |       |       |       |       |  |
| 65)   | 4,6-Dinitro-2-... |                | 0.085 | 0.103 | 0.119 | 0.120 | 0.124 | 0.126 | 0.113 | 13.97 |  |
| 66) c | n-Nitrosodiphe... | 0.684          | 0.661 | 0.679 | 0.641 | 0.615 | 0.636 | 0.590 | 0.644 | 5.29  |  |
| 67)   | 4-Bromophenyl-... | 0.243          | 0.240 | 0.250 | 0.241 | 0.230 | 0.240 | 0.227 | 0.239 | 3.35  |  |
| 68)   | Hexachlorobenzene | 0.266          | 0.262 | 0.267 | 0.257 | 0.248 | 0.256 | 0.243 | 0.257 | 3.53  |  |
| 69)   | Atrazine          | 0.154          | 0.156 | 0.164 | 0.168 | 0.171 | 0.174 | 0.176 | 0.166 | 5.09  |  |
| 70) C | Pentachlorophenol |                | 0.088 | 0.106 | 0.124 | 0.127 | 0.131 | 0.131 | 0.118 | 14.47 |  |
| 71)   | Phenanthrene      | 1.188          | 1.123 | 1.130 | 1.039 | 1.015 | 1.033 | 0.969 | 1.071 | 7.23  |  |
| 72)   | Anthracene        | 1.191          | 1.144 | 1.148 | 1.074 | 1.022 | 1.036 | 0.980 | 1.085 | 7.16  |  |
| 73)   | Carbazole         | 0.949          | 0.925 | 0.932 | 0.882 | 0.849 | 0.861 | 0.825 | 0.889 | 5.30  |  |
| 74)   | Di-n-butylphth... | 0.782          | 0.800 | 0.830 | 0.850 | 0.833 | 0.857 | 0.855 | 0.829 | 3.47  |  |
| 75) C | Fluoranthene      | 1.061          | 1.020 | 1.002 | 0.946 | 0.914 | 0.923 | 0.862 | 0.961 | 7.21  |  |
| 76) I | Chrysene-d12      | -----ISTD----- |       |       |       |       |       |       |       |       |  |
| 77)   | Benzidine         |                | 0.611 | 0.665 | 0.516 | 0.391 | 0.368 | 0.372 | 0.487 | 26.70 |  |
| 78)   | Pyrene            | 1.891          | 1.768 | 1.793 | 1.635 | 1.590 | 1.512 | 1.327 | 1.645 | 11.63 |  |
| 79) S | Terphenyl-d14     | 1.341          | 1.255 | 1.248 | 1.130 | 1.094 | 1.052 | 0.904 | 1.146 | 12.86 |  |
| 80)   | Butylbenzylpht... | 0.380          | 0.402 | 0.443 | 0.467 | 0.461 | 0.509 | 0.485 | 0.449 | 10.15 |  |
| 81)   | Benzo(a)anthra... | 1.302          | 1.274 | 1.300 | 1.318 | 1.266 | 1.316 | 1.237 | 1.288 | 2.30  |  |
| 82)   | 3,3'-Dichlorob... |                | 0.353 | 0.386 | 0.373 | 0.354 | 0.377 | 0.370 | 0.369 | 3.60  |  |
| 83)   | Chrysene          | 1.251          | 1.195 | 1.213 | 1.113 | 1.089 | 1.125 | 1.114 | 1.157 | 5.33  |  |
| 84)   | Bis(2-ethylhex... | 0.600          | 0.626 | 0.699 | 0.710 | 0.690 | 0.767 | 0.700 | 0.685 | 8.12  |  |
| 85) c | Di-n-octyl pht... |                | 1.129 | 1.289 | 1.298 | 1.251 | 1.413 | 1.228 | 1.268 | 7.36  |  |

Method Path : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\  
Method File : 8270-BF082025.M

|       |                   |                |       |       |       |       |       |       |       |      |
|-------|-------------------|----------------|-------|-------|-------|-------|-------|-------|-------|------|
| 86) I | Perylene-d12      | -----ISTD----- |       |       |       |       |       |       |       |      |
| 87)   | Indeno(1,2,3-c... | 1.425          | 1.433 | 1.565 | 1.449 | 1.427 | 1.466 | 1.295 | 1.437 | 5.51 |
| 88)   | Benzo(b)fluora... | 1.135          | 1.147 | 1.182 | 1.074 | 1.058 | 1.068 | 1.119 | 1.112 | 4.20 |
| 89)   | Benzo(k)fluora... | 1.157          | 1.014 | 1.056 | 1.081 | 1.034 | 1.104 | 1.005 | 1.064 | 5.10 |
| 90) C | Benzo(a)pyrene    | 1.033          | 1.021 | 1.074 | 1.032 | 1.010 | 1.043 | 1.011 | 1.032 | 2.15 |
| 91)   | Dibenzo(a,h)an... | 1.139          | 1.164 | 1.260 | 1.162 | 1.142 | 1.168 | 1.023 | 1.151 | 6.04 |
| 92)   | Benzo(g,h,i)pe... | 1.128          | 1.129 | 1.242 | 1.156 | 1.145 | 1.172 | 1.039 | 1.145 | 5.30 |

(#) = Out of Range

Method Path : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\  
 Method File : 8270-BF082625.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 Last Update : Tue Aug 26 16:30:52 2025  
 Response Via : Initial Calibration

## Calibration Files

2.5 =BF143585.D 5 =BF143586.D 10 =BF143587.D 20 =BF143588.D 40 =BF143589.D 50 =BF143590.D 60 =BF143591.D 80 =BF143592.D

|       | Compound              | 2.5            | 5     | 10    | 20    | 40    | 50    | 60    | 80    | Avg    | %RSD |
|-------|-----------------------|----------------|-------|-------|-------|-------|-------|-------|-------|--------|------|
| 1) I  | 1,4-Dichlorobenzen... | -----ISTD----- |       |       |       |       |       |       |       |        |      |
| 2)    | 1,4-Dioxane           | 0.606          | 0.579 | 0.592 | 0.571 | 0.546 | 0.570 | 0.561 | 0.575 | 3.46   |      |
| 3)    | Pyridine              | 1.218          | 1.341 | 1.525 | 1.504 | 1.491 | 1.540 | 1.487 | 1.444 | 8.25   |      |
| 4)    | n-Nitrosodimet...     |                | 0.669 | 0.683 | 0.695 | 0.690 | 0.703 | 0.682 | 0.687 | 1.71   |      |
| 5) S  | 2-Fluorophenol        | 1.245          | 1.223 | 1.243 | 1.152 | 1.114 | 1.151 | 1.098 | 1.175 | 5.20   |      |
| 6)    | Aniline               | 2.116          | 2.036 | 2.151 | 1.980 | 1.970 | 2.048 | 1.921 | 2.032 | 4.04   |      |
| 7) S  | Phenol-d6             | 1.584          | 1.519 | 1.521 | 1.422 | 1.373 | 1.424 | 1.346 | 1.455 | 6.00   |      |
| 8)    | 2-Chlorophenol        | 1.255          | 1.263 | 1.286 | 1.239 | 1.207 | 1.259 | 1.195 | 1.243 | 2.61   |      |
| 9)    | Benzaldehyde          |                | 1.075 | 1.018 | 1.209 | 1.011 | 1.162 |       | 1.095 | 8.00   |      |
| 10) C | Phenol                | 1.726          | 1.685 | 1.720 | 1.584 | 1.558 | 1.617 | 1.536 | 1.632 | 4.78   |      |
| 11)   | bis(2-Chloroet...     | 1.358          | 1.306 | 1.341 | 1.211 | 1.157 | 1.224 | 1.158 | 1.251 | 6.73   |      |
| 12)   | 1,3-Dichlorobe...     | 1.574          | 1.506 | 1.524 | 1.378 | 1.339 | 1.377 | 1.298 | 1.428 | 7.37   |      |
| 13) C | 1,4-Dichlorobe...     | 1.615          | 1.517 | 1.536 | 1.400 | 1.328 | 1.388 | 1.296 | 1.440 | 8.18   |      |
| 14)   | 1,2-Dichlorobe...     | 1.494          | 1.423 | 1.457 | 1.312 | 1.280 | 1.328 | 1.266 | 1.366 | 6.65   |      |
| 15)   | Benzyl Alcohol        |                | 1.074 | 1.135 | 1.088 | 1.074 | 1.117 | 1.057 | 1.091 | 2.71   |      |
| 16)   | 2,2'-oxybis(1-...     | 2.622          | 2.449 | 2.481 | 2.223 | 2.120 | 2.217 | 2.050 | 2.309 | 9.11   |      |
| 17)   | 2-Methylphenol        | 1.099          | 1.076 | 1.105 | 1.031 | 1.017 | 1.059 | 1.000 | 1.055 | 3.86   |      |
| 18)   | Hexachloroethane      | 0.451          | 0.452 | 0.484 | 0.460 | 0.449 | 0.471 | 0.436 | 0.458 | 3.40   |      |
| 19) P | n-Nitroso-di-n...     | 0.959          | 0.979 | 0.951 | 0.968 | 0.876 | 0.860 | 0.921 | 0.863 | 0.922  | 5.35 |
| 20)   | 3+4-Methylphenols     |                | 1.389 | 1.411 | 1.288 | 1.227 | 1.268 | 1.195 | 1.296 | 6.69   |      |
| 21) I | Naphthalene-d8        | -----ISTD----- |       |       |       |       |       |       |       |        |      |
| 22)   | Acetophenone          | 0.513          | 0.484 | 0.479 | 0.437 | 0.418 | 0.423 | 0.399 | 0.450 | 9.28   |      |
| 23) S | Nitrobenzene-d5       |                | 0.207 | 0.260 | 0.293 | 0.300 | 0.315 | 0.307 | 0.280 | 14.57  |      |
| 24)   | Nitrobenzene          | 0.217          | 0.250 | 0.305 | 0.330 | 0.329 | 0.342 | 0.331 | 0.300 | 15.99  |      |
| 25)   | Isophorone            | 0.694          | 0.662 | 0.683 | 0.645 | 0.628 | 0.654 | 0.623 | 0.656 | 4.02   |      |
| 26) C | 2-Nitrophenol         | 0.043          | 0.056 | 0.080 | 0.114 | 0.130 | 0.142 | 0.142 | 0.101 | 40.76# |      |
| 27)   | 2,4-Dimethylph...     | 0.292          | 0.280 | 0.297 | 0.285 | 0.282 | 0.286 | 0.274 | 0.285 | 2.65   |      |
| 28)   | bis(2-Chloroet...     | 0.431          | 0.411 | 0.424 | 0.385 | 0.377 | 0.386 | 0.363 | 0.397 | 6.40   |      |
| 29) C | 2,4-Dichloroph...     | 0.257          | 0.268 | 0.289 | 0.284 | 0.273 | 0.284 | 0.269 | 0.275 | 4.18   |      |
| 30)   | 1,2,4-Trichlor...     | 0.312          | 0.297 | 0.296 | 0.284 | 0.275 | 0.281 | 0.263 | 0.287 | 5.67   |      |
| 31)   | Naphthalene           | 1.098          | 1.029 | 1.031 | 0.938 | 0.900 | 0.909 | 0.860 | 0.966 | 9.00   |      |
| 32)   | Benzoic acid          |                | 0.047 | 0.091 | 0.126 | 0.150 | 0.172 | 0.168 | 0.126 | 38.83  |      |
| 33)   | 4-Chloroaniline       | 0.373          | 0.352 | 0.360 | 0.338 | 0.325 | 0.326 | 0.315 | 0.341 | 6.16   |      |
| 34) C | Hexachlorobuta...     | 0.199          | 0.189 | 0.189 | 0.177 | 0.171 | 0.175 | 0.165 | 0.181 | 6.63   |      |
| 35)   | Caprolactam           |                | 0.071 | 0.078 | 0.078 | 0.079 | 0.083 | 0.078 | 0.078 | 4.87   |      |
| 36) C | 4-Chloro-3-met...     | 0.282          | 0.284 | 0.300 | 0.288 | 0.283 | 0.296 | 0.278 | 0.287 | 2.75   |      |
| 37)   | 2-Methylnaphth...     | 0.727          | 0.685 | 0.684 | 0.615 | 0.586 | 0.605 | 0.565 | 0.638 | 9.46   |      |
| 38)   | 1-Methylnaphth...     | 0.694          | 0.666 | 0.655 | 0.599 | 0.566 | 0.583 | 0.548 | 0.616 | 9.05   |      |

Method Path : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\  
 Method File : 8270-BF082625.M

|       |                   |                |       |       |       |       |       |       |       |       |  |
|-------|-------------------|----------------|-------|-------|-------|-------|-------|-------|-------|-------|--|
| 39) I | Acenaphthene-d10  | -----ISTD----- |       |       |       |       |       |       |       |       |  |
| 40)   | 1,2,4,5-Tetrac... | 0.576          | 0.542 | 0.554 | 0.510 | 0.498 | 0.491 | 0.470 | 0.520 | 7.33  |  |
| 41) P | Hexachlorocycl... | 0.201          | 0.249 | 0.264 | 0.273 | 0.279 | 0.270 | 0.256 | 11.23 |       |  |
| 42) S | 2,4,6-Tribromo... | 0.130          | 0.143 | 0.166 | 0.162 | 0.163 | 0.170 | 0.162 | 0.156 | 9.38  |  |
| 43) C | 2,4,6-Trichlor... | 0.285          | 0.320 | 0.357 | 0.374 | 0.359 | 0.364 | 0.348 | 0.344 | 9.00  |  |
| 44)   | 2,4,5-Trichlor... | 0.317          | 0.321 | 0.360 | 0.353 | 0.356 | 0.368 | 0.358 | 0.347 | 5.74  |  |
| 45) S | 2-Fluorobiphenyl  | 1.404          | 1.290 | 1.276 | 1.120 | 1.041 | 1.056 | 1.001 | 1.170 | 13.11 |  |
| 46)   | 1,1'-Biphenyl     | 1.588          | 1.493 | 1.480 | 1.358 | 1.282 | 1.287 | 1.226 | 1.388 | 9.67  |  |
| 47)   | 2-Chloronaphth... | 1.210          | 1.161 | 1.143 | 1.086 | 1.035 | 1.036 | 1.004 | 1.096 | 6.99  |  |
| 48)   | 2-Nitroaniline    | 0.118          | 0.147 | 0.225 | 0.287 | 0.307 | 0.328 | 0.318 | 0.247 | 34.69 |  |
| 49)   | Acenaphthylene    | 1.772          | 1.673 | 1.711 | 1.557 | 1.489 | 1.511 | 1.451 | 1.595 | 7.74  |  |
| 50)   | Dimethylphthalate | 1.309          | 1.266 | 1.306 | 1.199 | 1.180 | 1.204 | 1.147 | 1.230 | 5.18  |  |
| 51)   | 2,6-Dinitrotol... | 0.065          | 0.090 | 0.145 | 0.189 | 0.208 | 0.226 | 0.219 | 0.163 | 39.69 |  |
| 52) C | Acenaphthene      | 1.169          | 1.103 | 1.115 | 1.014 | 0.969 | 0.992 | 0.953 | 1.045 | 7.97  |  |
| 53)   | 3-Nitroaniline    | 0.092          | 0.127 | 0.194 | 0.233 | 0.246 | 0.262 | 0.251 | 0.201 | 33.23 |  |
| 54) P | 2,4-Dinitrophenol | 0.020          | 0.032 | 0.050 | 0.066 | 0.077 | 0.081 | 0.054 | 45.09 |       |  |
| 55)   | Dibenzofuran      | 1.716          | 1.633 | 1.648 | 1.470 | 1.425 | 1.451 | 1.379 | 1.532 | 8.54  |  |
| 56) P | 4-Nitrophenol     | 0.114          | 0.161 | 0.189 | 0.201 | 0.218 | 0.209 | 0.182 | 21.33 |       |  |
| 57)   | 2,4-Dinitrotol... | 0.081          | 0.120 | 0.186 | 0.246 | 0.276 | 0.294 | 0.289 | 0.213 | 40.38 |  |
| 58)   | Fluorene          | 1.368          | 1.270 | 1.231 | 1.089 | 1.035 | 1.056 | 1.002 | 1.150 | 12.10 |  |
| 59)   | 2,3,4,6-Tetrac... | 0.190          | 0.224 | 0.273 | 0.271 | 0.277 | 0.287 | 0.275 | 0.257 | 13.91 |  |
| 60)   | Diethylphthalate  | 1.219          | 1.198 | 1.229 | 1.110 | 1.097 | 1.147 | 1.068 | 1.153 | 5.54  |  |
| 61)   | 4-Chlorophenyl... | 0.656          | 0.616 | 0.603 | 0.533 | 0.515 | 0.525 | 0.489 | 0.562 | 11.02 |  |
| 62)   | 4-Nitroaniline    | 0.115          | 0.143 | 0.191 | 0.212 | 0.229 | 0.244 | 0.237 | 0.196 | 25.31 |  |
| 63)   | Azobenzene        | 1.269          | 1.227 | 1.262 | 1.128 | 1.111 | 1.126 | 1.058 | 1.169 | 7.11  |  |
| 64) I | Phenanthrene-d10  | -----ISTD----- |       |       |       |       |       |       |       |       |  |
| 65)   | 4,6-Dinitro-2-... | 0.020          | 0.035 | 0.058 | 0.071 | 0.080 | 0.080 | 0.057 | 42.94 |       |  |
| 66) c | n-Nitrosodiphe... | 0.694          | 0.664 | 0.662 | 0.632 | 0.598 | 0.610 | 0.593 | 0.636 | 6.02  |  |
| 67)   | 4-Bromophenyl-... | 0.223          | 0.219 | 0.228 | 0.219 | 0.208 | 0.211 | 0.206 | 0.216 | 3.74  |  |
| 68)   | Hexachlorobenzene | 0.247          | 0.239 | 0.242 | 0.229 | 0.220 | 0.226 | 0.217 | 0.231 | 5.05  |  |
| 69)   | Atrazine          | 0.179          | 0.184 | 0.188 | 0.181 | 0.182 | 0.189 | 0.177 | 0.183 | 2.48  |  |
| 70) C | Pentachlorophenol | 0.101          | 0.126 | 0.139 | 0.140 | 0.145 | 0.141 | 0.132 | 12.49 |       |  |
| 71)   | Phenanthrene      | 1.223          | 1.162 | 1.127 | 1.040 | 0.995 | 0.994 | 0.955 | 1.071 | 9.38  |  |
| 72)   | Anthracene        | 1.223          | 1.168 | 1.151 | 1.041 | 0.998 | 1.004 | 0.969 | 1.079 | 9.22  |  |
| 73)   | Carbazole         | 1.078          | 1.019 | 0.991 | 0.892 | 0.855 | 0.865 | 0.830 | 0.933 | 10.23 |  |
| 74)   | Di-n-butylphth... | 0.961          | 0.968 | 1.013 | 0.923 | 0.925 | 0.968 | 0.893 | 0.950 | 4.15  |  |
| 75) C | Fluoranthene      | 1.176          | 1.101 | 1.065 | 0.897 | 0.862 | 0.901 | 0.842 | 0.978 | 13.62 |  |
| 76) I | Chrysene-d12      | -----ISTD----- |       |       |       |       |       |       |       |       |  |
| 77)   | Benzidine         | 0.598          | 0.654 | 0.538 | 0.357 | 0.222 | 0.394 | 0.460 | 35.65 |       |  |
| 78)   | Pyrene            | 1.722          | 1.768 | 1.810 | 1.609 | 1.687 | 1.613 | 1.507 | 1.674 | 6.25  |  |
| 79) S | Terphenyl-d14     | 1.199          | 1.213 | 1.221 | 1.044 | 1.080 | 1.053 | 0.992 | 1.115 | 8.45  |  |
| 80)   | Butylbenzylpht... | 0.211          | 0.276 | 0.386 | 0.456 | 0.455 | 0.513 | 0.490 | 0.398 | 28.68 |  |
| 81)   | Benzo(a)anthra... | 1.365          | 1.276 | 1.288 | 1.257 | 1.216 | 1.220 | 1.186 | 1.258 | 4.73  |  |
| 82)   | 3,3'-Dichlorob... | 0.271          | 0.339 | 0.373 | 0.340 | 0.334 | 0.367 | 0.337 | 10.77 |       |  |
| 83)   | Chrysene          | 1.239          | 1.174 | 1.244 | 1.125 | 1.129 | 1.168 | 1.129 | 1.173 | 4.35  |  |
| 84)   | Bis(2-ethylhex... | 0.285          | 0.397 | 0.529 | 0.632 | 0.580 | 0.637 | 0.632 | 0.528 | 25.91 |  |
| 85) c | Di-n-octyl pht... | 0.564          | 0.840 | 1.167 | 1.128 | 1.179 | 1.223 | 1.017 | 25.62 |       |  |

Method Path : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\  
Method File : 8270-BF082625.M

|       |                   |                |       |       |       |       |       |       |       |       |
|-------|-------------------|----------------|-------|-------|-------|-------|-------|-------|-------|-------|
| 86) I | Perylene-d12      | -----ISTD----- |       |       |       |       |       |       |       |       |
| 87)   | Indeno(1,2,3-c... | 1.127          | 1.316 | 1.462 | 1.429 | 1.401 | 1.430 | 1.359 | 1.361 | 8.36  |
| 88)   | Benzo(b)fluora... | 1.354          | 1.208 | 1.131 | 1.045 | 1.008 | 1.184 | 1.069 | 1.143 | 10.35 |
| 89)   | Benzo(k)fluora... | 1.169          | 1.128 | 1.165 | 1.061 | 1.028 | 0.982 | 0.947 | 1.068 | 8.28  |
| 90) C | Benzo(a)pyrene    | 1.007          | 0.994 | 1.034 | 1.032 | 0.989 | 1.030 | 0.982 | 1.010 | 2.20  |
| 91)   | Dibenzo(a,h)an... | 0.980          | 1.093 | 1.191 | 1.154 | 1.127 | 1.163 | 1.089 | 1.114 | 6.25  |
| 92)   | Benzo(g,h,i)pe... | 0.935          | 1.071 | 1.154 | 1.148 | 1.148 | 1.157 | 1.106 | 1.103 | 7.31  |

(#) = Out of Range

Method Path : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\  
 Method File : 8270E-BP081425.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 Last Update : Thu Aug 14 18:14:31 2025  
 Response Via : Initial Calibration

## Calibration Files

2.5 =BP025392.D 5 =BP025393.D 10 =BP025394.D 20 =BP025395.D 40 =BP025396.D 50 =BP025397.D 60 =BP025398.D 80 =BP025399.D

| Compound                   | 2.5            | 5     | 10    | 20    | 40    | 50    | 60    | 80    | Avg   | %RSD  |
|----------------------------|----------------|-------|-------|-------|-------|-------|-------|-------|-------|-------|
| -----                      |                |       |       |       |       |       |       |       |       |       |
| 1) I 1,4-Dichlorobenzen... | -----ISTD----- |       |       |       |       |       |       |       |       |       |
| 2) 1,4-Dioxane             |                | 0.520 | 0.464 | 0.475 | 0.445 | 0.479 | 0.462 | 0.458 | 0.472 | 5.07  |
| 3) Pyridine                |                | 1.446 | 1.250 | 1.256 | 1.220 | 1.317 | 1.281 | 1.268 | 1.291 | 5.77  |
| 4) n-Nitrosodimet...       |                |       | 0.515 | 0.512 | 0.511 | 0.559 | 0.555 | 0.541 | 0.532 | 4.13  |
| 5) S 2-Fluorophenol        |                | 1.232 | 1.177 | 1.194 | 1.162 | 1.240 | 1.226 | 1.199 | 1.204 | 2.42  |
| 6) Aniline                 |                | 2.098 | 1.962 | 2.030 | 2.020 | 2.194 | 2.164 | 2.094 | 2.080 | 3.96  |
| 7) S Phenol-d6             |                | 1.483 | 1.467 | 1.533 | 1.513 | 1.641 | 1.609 | 1.563 | 1.544 | 4.16  |
| 8) 2-Chlorophenol          |                | 1.351 | 1.283 | 1.347 | 1.313 | 1.428 | 1.414 | 1.378 | 1.359 | 3.83  |
| 9) Benzaldehyde            |                |       | 0.987 | 0.956 | 1.366 | 1.262 | 1.005 | 0.915 | 1.082 | 17.12 |
| 10) C Phenol               |                | 1.594 | 1.529 | 1.613 | 1.575 | 1.707 | 1.688 | 1.635 | 1.620 | 3.87  |
| 11) bis(2-Chloroet...      |                | 1.283 | 1.183 | 1.273 | 1.222 | 1.321 | 1.303 | 1.255 | 1.263 | 3.78  |
| 12) 1,3-Dichlorobe...      |                | 1.532 | 1.433 | 1.503 | 1.443 | 1.555 | 1.530 | 1.493 | 1.499 | 3.07  |
| 13) C 1,4-Dichlorobe...    |                | 1.587 | 1.490 | 1.527 | 1.467 | 1.596 | 1.544 | 1.510 | 1.532 | 3.12  |
| 14) 1,2-Dichlorobe...      |                | 1.558 | 1.430 | 1.491 | 1.431 | 1.533 | 1.487 | 1.449 | 1.483 | 3.35  |
| 15) Benzyl Alcohol         |                |       | 1.138 | 1.205 | 1.193 | 1.287 | 1.289 | 1.250 | 1.227 | 4.83  |
| 16) 2,2'-oxybis(1-...      |                | 1.762 | 1.592 | 1.712 | 1.620 | 1.751 | 1.740 | 1.665 | 1.692 | 3.96  |
| 17) 2-Methylphenol         |                | 1.075 | 1.030 | 1.123 | 1.107 | 1.198 | 1.191 | 1.154 | 1.125 | 5.41  |
| 18) Hexachloroethane       |                | 0.577 | 0.544 | 0.552 | 0.545 | 0.588 | 0.578 | 0.559 | 0.563 | 3.13  |
| 19) P n-Nitroso-di-n...    | 1.014          | 1.066 | 0.959 | 1.056 | 0.992 | 1.053 | 1.042 | 1.000 | 1.023 | 3.68  |
| 20) 3+4-Methylphenols      |                |       | 1.442 | 1.533 | 1.525 | 1.639 | 1.640 | 1.580 | 1.560 | 4.87  |
|                            |                |       |       |       |       |       |       |       |       |       |
| 21) I Naphthalene-d8       | -----ISTD----- |       |       |       |       |       |       |       |       |       |
| 22) Acetophenone           |                | 0.515 | 0.487 | 0.506 | 0.485 | 0.523 | 0.506 | 0.488 | 0.502 | 2.96  |
| 23) S Nitrobenzene-d5      |                | 0.395 | 0.376 | 0.392 | 0.383 | 0.419 | 0.408 | 0.394 | 0.395 | 3.65  |
| 24) Nitrobenzene           |                | 0.350 | 0.342 | 0.348 | 0.344 | 0.375 | 0.362 | 0.352 | 0.353 | 3.21  |
| 25) Isophorone             |                | 0.734 | 0.672 | 0.700 | 0.674 | 0.730 | 0.709 | 0.680 | 0.700 | 3.70  |
| 26) C 2-Nitrophenol        |                | 0.167 | 0.163 | 0.177 | 0.181 | 0.197 | 0.195 | 0.190 | 0.181 | 7.43  |
| 27) 2,4-Dimethylph...      |                | 0.303 | 0.303 | 0.296 | 0.308 | 0.336 | 0.325 | 0.313 | 0.312 | 4.46  |
| 28) bis(2-Chloroet...      |                | 0.443 | 0.407 | 0.425 | 0.407 | 0.439 | 0.430 | 0.410 | 0.423 | 3.61  |
| 29) C 2,4-Dichloroph...    |                | 0.293 | 0.293 | 0.308 | 0.305 | 0.331 | 0.324 | 0.314 | 0.310 | 4.71  |
| 30) 1,2,4-Trichlor...      |                | 0.357 | 0.329 | 0.335 | 0.327 | 0.354 | 0.344 | 0.332 | 0.340 | 3.61  |
| 31) Naphthalene            |                | 1.069 | 1.020 | 1.037 | 1.000 | 1.087 | 1.052 | 1.011 | 1.040 | 3.06  |
| 32) Benzoic acid           |                |       | 0.188 | 0.207 | 0.227 | 0.249 | 0.255 | 0.255 | 0.230 | 12.20 |
| 33) 4-Chloroaniline        |                | 0.453 | 0.435 | 0.447 | 0.452 | 0.492 | 0.478 | 0.458 | 0.459 | 4.19  |
| 34) C Hexachlorobuta...    |                | 0.222 | 0.205 | 0.211 | 0.205 | 0.218 | 0.213 | 0.207 | 0.211 | 3.05  |
| 35) Caprolactam            |                |       | 0.114 | 0.110 | 0.109 | 0.120 | 0.116 | 0.113 | 0.114 | 3.67  |
| 36) C 4-Chloro-3-met...    |                | 0.366 | 0.335 | 0.347 | 0.346 | 0.373 | 0.368 | 0.353 | 0.355 | 3.85  |
| 37) 2-Methylnaphth...      |                | 0.716 | 0.653 | 0.678 | 0.652 | 0.700 | 0.682 | 0.654 | 0.676 | 3.69  |
| 38) 1-Methylnaphth...      |                | 0.777 | 0.706 | 0.713 | 0.686 | 0.751 | 0.719 | 0.685 | 0.719 | 4.67  |

Method Path : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\  
 Method File : 8270E-BP081425.M

|       |                   |                |       |       |       |       |       |       |       |       |
|-------|-------------------|----------------|-------|-------|-------|-------|-------|-------|-------|-------|
| 39) I | Acenaphthene-d10  | -----ISTD----- |       |       |       |       |       |       |       |       |
| 40)   | 1,2,4,5-Tetrac... | 0.578          | 0.559 | 0.585 | 0.560 | 0.610 | 0.581 | 0.571 | 0.578 | 3.00  |
| 41) P | Hexachlorocycl... | 0.274          | 0.313 | 0.348 | 0.387 | 0.388 | 0.383 | 0.349 |       | 13.46 |
| 42) S | 2,4,6-Tribromo... | 0.266          | 0.259 | 0.267 | 0.263 | 0.286 | 0.275 | 0.268 | 0.269 | 3.39  |
| 43) C | 2,4,6-Trichlor... | 0.368          | 0.364 | 0.381 | 0.382 | 0.421 | 0.408 | 0.405 | 0.390 | 5.56  |
| 44)   | 2,4,5-Trichlor... | 0.417          | 0.385 | 0.424 | 0.424 | 0.457 | 0.448 | 0.437 | 0.427 | 5.52  |
| 45) S | 2-Fluorobiphenyl  | 1.558          | 1.483 | 1.507 | 1.392 | 1.517 | 1.435 | 1.351 | 1.463 | 5.04  |
| 46)   | 1,1'-Biphenyl     | 1.495          | 1.418 | 1.457 | 1.408 | 1.537 | 1.445 | 1.407 | 1.453 | 3.37  |
| 47)   | 2-Chloronaphth... | 1.152          | 1.093 | 1.144 | 1.101 | 1.184 | 1.142 | 1.100 | 1.131 | 2.99  |
| 48)   | 2-Nitroaniline    | 0.301          | 0.306 | 0.317 | 0.330 | 0.360 | 0.351 | 0.341 | 0.329 | 6.85  |
| 49)   | Acenaphthylene    | 1.907          | 1.806 | 1.848 | 1.806 | 1.992 | 1.907 | 1.832 | 1.871 | 3.63  |
| 50)   | Dimethylphthalate | 1.569          | 1.449 | 1.438 | 1.403 | 1.512 | 1.479 | 1.403 | 1.465 | 4.12  |
| 51)   | 2,6-Dinitrotol... | 0.297          | 0.287 | 0.302 | 0.309 | 0.331 | 0.323 | 0.311 | 0.309 | 4.95  |
| 52) C | Acenaphthene      | 1.149          | 1.082 | 1.100 | 1.041 | 1.155 | 1.106 | 1.056 | 1.098 | 3.93  |
| 53)   | 3-Nitroaniline    | 0.318          | 0.329 | 0.337 | 0.342 | 0.384 | 0.367 | 0.354 | 0.347 | 6.56  |
| 54) P | 2,4-Dinitrophenol | 0.114          | 0.140 | 0.163 | 0.186 | 0.192 | 0.194 | 0.165 |       | 19.83 |
| 55)   | Dibenzofuran      | 1.820          | 1.711 | 1.737 | 1.677 | 1.811 | 1.729 | 1.672 | 1.736 | 3.40  |
| 56) P | 4-Nitrophenol     | 0.251          | 0.265 | 0.279 | 0.315 | 0.304 | 0.298 | 0.285 |       | 8.56  |
| 57)   | 2,4-Dinitrotol... | 0.404          | 0.407 | 0.431 | 0.438 | 0.479 | 0.471 | 0.454 | 0.440 | 6.61  |
| 58)   | Fluorene          | 1.506          | 1.391 | 1.390 | 1.306 | 1.415 | 1.330 | 1.253 | 1.370 | 6.02  |
| 59)   | 2,3,4,6-Tetrac... | 0.375          | 0.360 | 0.366 | 0.368 | 0.402 | 0.392 | 0.379 | 0.377 | 4.01  |
| 60)   | Diethylphthalate  | 1.615          | 1.475 | 1.465 | 1.434 | 1.540 | 1.459 | 1.430 | 1.488 | 4.48  |
| 61)   | 4-Chlorophenyl... | 0.761          | 0.684 | 0.679 | 0.651 | 0.690 | 0.671 | 0.634 | 0.681 | 5.91  |
| 62)   | 4-Nitroaniline    | 0.341          | 0.336 | 0.349 | 0.351 | 0.387 | 0.370 | 0.359 | 0.356 | 4.89  |
| 63)   | Azobenzene        | 1.315          | 1.226 | 1.274 | 1.237 | 1.341 | 1.301 | 1.227 | 1.274 | 3.62  |
| 64) I | Phenanthrene-d10  | -----ISTD----- |       |       |       |       |       |       |       |       |
| 65)   | 4,6-Dinitro-2-... | 0.096          | 0.113 | 0.123 | 0.139 | 0.139 | 0.138 | 0.125 |       | 14.21 |
| 66) c | n-Nitrosodiphe... | 0.636          | 0.603 | 0.607 | 0.591 | 0.634 | 0.605 | 0.594 | 0.610 | 2.99  |
| 67)   | 4-Bromophenyl-... | 0.233          | 0.206 | 0.217 | 0.213 | 0.230 | 0.223 | 0.218 | 0.220 | 4.34  |
| 68)   | Hexachlorobenzene | 0.264          | 0.253 | 0.254 | 0.242 | 0.264 | 0.257 | 0.251 | 0.255 | 3.06  |
| 69)   | Atrazine          | 0.229          | 0.226 | 0.228 | 0.218 | 0.240 | 0.229 | 0.229 | 0.228 | 2.74  |
| 70) C | Pentachlorophenol | 0.135          | 0.154 | 0.157 | 0.177 | 0.171 | 0.172 | 0.161 |       | 9.76  |
| 71)   | Phenanthrene      | 1.162          | 1.101 | 1.111 | 1.040 | 1.135 | 1.085 | 1.057 | 1.099 | 3.88  |
| 72)   | Anthracene        | 1.161          | 1.124 | 1.133 | 1.079 | 1.175 | 1.101 | 1.080 | 1.122 | 3.37  |
| 73)   | Carbazole         | 1.079          | 1.052 | 1.074 | 1.007 | 1.111 | 1.053 | 1.022 | 1.057 | 3.32  |
| 74)   | Di-n-butylphth... | 1.345          | 1.255 | 1.332 | 1.244 | 1.374 | 1.285 | 1.265 | 1.300 | 3.86  |
| 75) C | Fluoranthene      | 1.370          | 1.323 | 1.348 | 1.228 | 1.361 | 1.286 | 1.258 | 1.311 | 4.13  |
| 76) I | Chrysene-d12      | -----ISTD----- |       |       |       |       |       |       |       |       |
| 77)   | Benzidine         | 0.642          | 0.642 | 0.967 | 0.664 | 0.608 | 0.573 | 0.683 |       | 20.94 |
| 78)   | Pyrene            | 1.398          | 1.261 | 1.287 | 1.274 | 1.320 | 1.319 | 1.241 | 1.300 | 4.00  |
| 79) S | Terphenyl-d14     | 1.249          | 1.100 | 1.097 | 1.045 | 1.100 | 1.061 | 0.999 | 1.093 | 7.15  |
| 80)   | Butylbenzylpht... | 0.589          | 0.557 | 0.592 | 0.579 | 0.614 | 0.616 | 0.578 | 0.589 | 3.55  |
| 81)   | Benzo(a)anthra... | 1.384          | 1.302 | 1.313 | 1.268 | 1.359 | 1.333 | 1.275 | 1.319 | 3.21  |
| 82)   | 3,3'-Dichlorob... | 0.482          | 0.507 | 0.488 | 0.521 | 0.506 | 0.480 | 0.497 |       | 3.31  |
| 83)   | Chrysene          | 1.292          | 1.214 | 1.249 | 1.182 | 1.280 | 1.243 | 1.186 | 1.235 | 3.49  |
| 84)   | Bis(2-ethylhex... | 0.907          | 0.826 | 0.874 | 0.831 | 0.901 | 0.882 | 0.850 | 0.867 | 3.73  |
| 85) c | Di-n-octyl pht... | 1.381          | 1.498 | 1.444 | 1.572 | 1.563 | 1.491 | 1.492 |       | 4.84  |

Method Path : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\  
Method File : 8270E-BP081425.M

|       |                   |                |       |       |       |       |       |       |       |      |
|-------|-------------------|----------------|-------|-------|-------|-------|-------|-------|-------|------|
| 86) I | Perylene-d12      | -----ISTD----- |       |       |       |       |       |       |       |      |
| 87)   | Indeno(1,2,3-c... | 1.454          | 1.385 | 1.440 | 1.411 | 1.569 | 1.522 | 1.486 | 1.467 | 4.36 |
| 88)   | Benzo(b)fluora... | 1.203          | 1.126 | 1.171 | 1.134 | 1.250 | 1.197 | 1.175 | 1.179 | 3.60 |
| 89)   | Benzo(k)fluora... | 1.202          | 1.157 | 1.200 | 1.130 | 1.226 | 1.206 | 1.141 | 1.180 | 3.16 |
| 90) C | Benzo(a)pyrene    | 1.139          | 1.093 | 1.150 | 1.108 | 1.219 | 1.185 | 1.141 | 1.148 | 3.77 |
| 91)   | Dibenzo(a,h)an... | 1.180          | 1.125 | 1.179 | 1.156 | 1.290 | 1.244 | 1.216 | 1.199 | 4.64 |
| 92)   | Benzo(g,h,i)pe... | 1.162          | 1.118 | 1.169 | 1.138 | 1.248 | 1.220 | 1.192 | 1.178 | 3.86 |

(#) = Out of Range

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>AEC002</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q2924</u>                 |
| Instrument ID:  | <u>BNA_F</u>      | Calibration Date/Time: | <u>08/22/2025 10:29</u>      |
| Lab File ID:    | <u>BF143563.D</u> | Init. Calib. Date(s):  | <u>08/20/2025 08/20/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>10:55 14:20</u>           |
| GC Column:      | <u>DB-UI</u>      | ID:                    | <u>0.18</u> (mm)             |

| COMPOUND                    | RRF   | RRF040 | MIN RRF | %D    | MAX%D |
|-----------------------------|-------|--------|---------|-------|-------|
| 2-Fluorophenol              | 1.145 | 1.105  |         | -3.5  |       |
| Benzaldehyde                | 1.145 | 0.999  |         | -12.8 |       |
| Phenol-d6                   | 1.414 | 1.366  |         | -3.4  |       |
| Phenol                      | 1.629 | 1.599  |         | -1.8  | 20.0  |
| bis(2-Chloroethyl)ether     | 1.217 | 1.157  |         | -4.9  |       |
| 2-Chlorophenol              | 1.265 | 1.220  |         | -3.6  |       |
| 2-Methylphenol              | 1.015 | 0.981  |         | -3.3  |       |
| 2,2-oxybis(1-Chloropropane) | 1.947 | 1.762  |         | -9.5  |       |
| Acetophenone                | 0.432 | 0.414  |         | -4.2  |       |
| 3+4-Methylphenols           | 1.212 | 1.227  |         | 1.2   |       |
| n-Nitroso-di-n-propylamine  | 0.864 | 0.857  | 0.050   | -0.8  |       |
| Nitrobenzene-d5             | 0.343 | 0.333  |         | -2.9  |       |
| Hexachloroethane            | 0.485 | 0.478  |         | -1.4  |       |
| Nitrobenzene                | 0.353 | 0.349  |         | -1.1  |       |
| Isophorone                  | 0.628 | 0.612  |         | -2.5  |       |
| 2-Nitrophenol               | 0.173 | 0.169  |         | -2.3  | 20.0  |
| 2,4-Dimethylphenol          | 0.269 | 0.253  |         | -5.9  |       |
| bis(2-Chloroethoxy)methane  | 0.386 | 0.363  |         | -6.0  |       |
| 2,4-Dichlorophenol          | 0.288 | 0.278  |         | -3.5  | 20.0  |
| Naphthalene                 | 0.960 | 0.899  |         | -6.4  |       |
| 4-Chloroaniline             | 0.341 | 0.325  |         | -4.7  |       |
| Hexachlorobutadiene         | 0.193 | 0.186  |         | -3.6  | 20.0  |
| Caprolactam                 | 0.069 | 0.078  |         | 13.0  |       |
| 4-Chloro-3-methylphenol     | 0.288 | 0.292  |         | 1.4   | 20.0  |
| 2-Methylnaphthalene         | 0.641 | 0.610  |         | -4.8  |       |
| Hexachlorocyclopentadiene   | 0.184 | 0.193  | 0.050   | 4.9   |       |
| 2,4,6-Trichlorophenol       | 0.352 | 0.336  |         | -4.5  | 20.0  |
| 2-Fluorobiphenyl            | 1.210 | 1.143  |         | -5.5  |       |
| 2,4,5-Trichlorophenol       | 0.385 | 0.355  |         | -7.8  |       |
| 1,1-Biphenyl                | 1.420 | 1.319  |         | -7.1  |       |
| 2-Chloronaphthalene         | 1.114 | 1.033  |         | -7.3  |       |
| 2-Nitroaniline              | 0.326 | 0.330  |         | 1.2   |       |
| Dimethylphthalate           | 1.205 | 1.196  |         | -0.7  |       |
| Acenaphthylene              | 1.595 | 1.486  |         | -6.8  |       |
| 2,6-Dinitrotoluene          | 0.250 | 0.249  |         | -0.4  |       |
| 3-Nitroaniline              | 0.270 | 0.267  |         | -1.1  |       |
| Acenaphthene                | 1.083 | 1.038  |         | -4.2  | 20.0  |
| 2,4-Dinitrophenol           | 0.115 | 0.130  | 0.050   | 13.0  |       |
| 4-Nitrophenol               | 0.161 | 0.193  | 0.050   | 19.9  |       |

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>AECO02</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q2924</u>                 |
| Instrument ID:  | <u>BNA_F</u>      | Calibration Date/Time: | <u>08/22/2025 10:29</u>      |
| Lab File ID:    | <u>BF143563.D</u> | Init. Calib. Date(s):  | <u>08/20/2025 08/20/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>10:55 14:20</u>           |
| GC Column:      | <u>DB-UI</u>      | ID:                    | <u>0.18 (mm)</u>             |

| COMPOUND                   | RRF   | RRF040 | MIN<br>RRF | %D    | MAX%D |
|----------------------------|-------|--------|------------|-------|-------|
| Dibenzofuran               | 1.544 | 1.440  |            | -6.7  |       |
| 2,4-Dinitrotoluene         | 0.333 | 0.338  |            | 1.5   |       |
| Diethylphthalate           | 1.080 | 1.129  |            | 4.5   |       |
| 4-Chlorophenyl-phenylether | 0.594 | 0.562  |            | -5.4  |       |
| Fluorene                   | 1.188 | 1.122  |            | -5.6  |       |
| 4-Nitroaniline             | 0.248 | 0.264  |            | 6.5   |       |
| 4,6-Dinitro-2-methylphenol | 0.113 | 0.102  |            | -9.7  |       |
| n-Nitrosodiphenylamine     | 0.644 | 0.599  |            | -7.0  | 20.0  |
| 2,4,6-Tribromophenol       | 0.186 | 0.178  |            | -4.3  |       |
| 4-Bromophenyl-phenylether  | 0.239 | 0.222  |            | -7.1  |       |
| Hexachlorobenzene          | 0.257 | 0.239  |            | -7.0  |       |
| Atrazine                   | 0.166 | 0.199  |            | 19.9  |       |
| Pentachlorophenol          | 0.118 | 0.108  |            | -8.5  | 20.0  |
| Phenanthrene               | 1.071 | 0.996  |            | -7.0  |       |
| Anthracene                 | 1.085 | 1.022  |            | -5.8  |       |
| Carbazole                  | 0.889 | 0.865  |            | -2.7  |       |
| Di-n-butylphthalate        | 0.829 | 0.965  |            | 16.4  |       |
| Fluoranthene               | 0.961 | 0.923  |            | -4.0  | 20.0  |
| Pyrene                     | 1.645 | 1.624  |            | -1.3  |       |
| Terphenyl-d14              | 1.146 | 1.162  |            | 1.4   |       |
| Butylbenzylphthalate       | 0.449 | 0.488  |            | 8.7   |       |
| 3,3-Dichlorobenzidine      | 0.369 | 0.335  |            | -9.2  |       |
| Benzo (a) anthracene       | 1.288 | 1.184  |            | -8.1  |       |
| Chrysene                   | 1.157 | 1.118  |            | -3.4  |       |
| Bis(2-ethylhexyl)phthalate | 0.685 | 0.617  |            | -9.9  |       |
| Di-n-octyl phthalate       | 1.268 | 1.115  |            | -12.1 | 20.0  |
| Benzo (b) fluoranthene     | 1.112 | 1.078  |            | -3.1  |       |
| Benzo (k) fluoranthene     | 1.064 | 1.024  |            | -3.8  |       |
| Benzo (a) pyrene           | 1.032 | 1.000  |            | -3.1  | 20.0  |
| Indeno (1,2,3-cd) pyrene   | 1.437 | 1.391  |            | -3.2  |       |
| Dibenzo (a,h) anthracene   | 1.151 | 1.121  |            | -2.6  |       |
| Benzo (g,h,i) perylene     | 1.145 | 1.103  |            | -3.7  |       |
| 1,2,4,5-Tetrachlorobenzene | 0.572 | 0.527  |            | -7.9  |       |
| 1,4-Dioxane                | 0.531 | 0.503  |            | -5.3  | 20.0  |
| 2,3,4,6-Tetrachlorophenol  | 0.286 | 0.274  |            | -4.2  |       |

All other compounds must meet a minimum RRF of 0.010.

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>AECO02</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q2924</u>                 |
| Instrument ID:  | <u>BNA_F</u>      | Calibration Date/Time: | <u>08/27/2025 09:16</u>      |
| Lab File ID:    | <u>BF143596.D</u> | Init. Calib. Date(s):  | <u>08/26/2025 08/26/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>09:11 12:32</u>           |
| GC Column:      | <u>DB-UI</u>      | ID:                    | <u>0.18</u> (mm)             |

| COMPOUND                    | RRF   | RRF040 | MIN RRF | %D   | MAX%D |
|-----------------------------|-------|--------|---------|------|-------|
| 2-Fluorophenol              | 1.175 | 1.200  |         | 2.1  |       |
| Benzaldehyde                | 1.095 | 1.252  |         | 14.3 |       |
| Phenol-d6                   | 1.455 | 1.474  |         | 1.3  |       |
| Phenol                      | 1.632 | 1.764  |         | 8.1  | 20.0  |
| bis(2-Chloroethyl)ether     | 1.251 | 1.256  |         | 0.4  |       |
| 2-Chlorophenol              | 1.243 | 1.295  |         | 4.2  |       |
| 2-Methylphenol              | 1.055 | 1.066  |         | 1.0  |       |
| 2,2-oxybis(1-Chloropropane) | 2.309 | 2.190  |         | -5.2 |       |
| Acetophenone                | 0.450 | 0.451  |         | 0.2  |       |
| 3+4-Methylphenols           | 1.296 | 1.370  |         | 5.7  |       |
| n-Nitroso-di-n-propylamine  | 0.922 | 0.915  | 0.050   | -0.8 |       |
| Nitrobenzene-d5             | 0.280 | 0.318  |         | 13.6 |       |
| Hexachloroethane            | 0.458 | 0.487  |         | 6.3  |       |
| Nitrobenzene                | 0.300 | 0.351  |         | 17.0 |       |
| Isophorone                  | 0.656 | 0.659  |         | 0.5  |       |
| 2-Nitrophenol               | 0.101 | 0.136  |         | 34.7 | 20.0  |
| 2,4-Dimethylphenol          | 0.285 | 0.296  |         | 3.9  |       |
| bis(2-Chloroethoxy)methane  | 0.397 | 0.401  |         | 1.0  |       |
| 2,4-Dichlorophenol          | 0.275 | 0.294  |         | 6.9  | 20.0  |
| Naphthalene                 | 0.966 | 0.981  |         | 1.6  |       |
| 4-Chloroaniline             | 0.341 | 0.352  |         | 3.2  |       |
| Hexachlorobutadiene         | 0.181 | 0.186  |         | 2.8  | 20.0  |
| Caprolactam                 | 0.078 | 0.086  |         | 10.3 |       |
| 4-Chloro-3-methylphenol     | 0.287 | 0.303  |         | 5.6  | 20.0  |
| 2-Methylnaphthalene         | 0.638 | 0.651  |         | 2.0  |       |
| Hexachlorocyclopentadiene   | 0.256 | 0.275  | 0.050   | 7.4  |       |
| 2,4,6-Trichlorophenol       | 0.344 | 0.359  |         | 4.4  | 20.0  |
| 2-Fluorobiphenyl            | 1.170 | 1.158  |         | -1.0 |       |
| 2,4,5-Trichlorophenol       | 0.347 | 0.387  |         | 11.5 |       |
| 1,1-Biphenyl                | 1.388 | 1.375  |         | -0.9 |       |
| 2-Chloronaphthalene         | 1.096 | 1.113  |         | 1.6  |       |
| 2-Nitroaniline              | 0.247 | 0.328  |         | 32.8 |       |
| Dimethylphthalate           | 1.230 | 1.274  |         | 3.6  |       |
| Acenaphthylene              | 1.595 | 1.622  |         | 1.7  |       |
| 2,6-Dinitrotoluene          | 0.163 | 0.229  |         | 40.5 |       |
| 3-Nitroaniline              | 0.201 | 0.274  |         | 36.3 |       |
| Acenaphthene                | 1.045 | 1.066  |         | 2.0  | 20.0  |
| 2,4-Dinitrophenol           | 0.054 | 0.068  | 0.050   | 25.9 |       |
| 4-Nitrophenol               | 0.182 | 0.219  | 0.050   | 20.3 |       |

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>AECO02</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q2924</u>                 |
| Instrument ID:  | <u>BNA_F</u>      | Calibration Date/Time: | <u>08/27/2025 09:16</u>      |
| Lab File ID:    | <u>BF143596.D</u> | Init. Calib. Date(s):  | <u>08/26/2025 08/26/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>09:11 12:32</u>           |
| GC Column:      | <u>DB-UI</u>      | ID:                    | <u>0.18</u> (mm)             |

| COMPOUND                   | RRF   | RRF040 | MIN<br>RRF | %D   | MAX%D |
|----------------------------|-------|--------|------------|------|-------|
| Dibenzofuran               | 1.532 | 1.562  |            | 2.0  |       |
| 2,4-Dinitrotoluene         | 0.213 | 0.308  |            | 44.6 |       |
| Diethylphthalate           | 1.153 | 1.218  |            | 5.6  |       |
| 4-Chlorophenyl-phenylether | 0.562 | 0.566  |            | 0.7  |       |
| Fluorene                   | 1.150 | 1.146  |            | -0.3 |       |
| 4-Nitroaniline             | 0.196 | 0.264  |            | 34.7 |       |
| 4,6-Dinitro-2-methylphenol | 0.057 | 0.073  |            | 28.1 |       |
| n-Nitrosodiphenylamine     | 0.636 | 0.634  |            | -0.3 | 20.0  |
| 2,4,6-Tribromophenol       | 0.156 | 0.181  |            | 16.0 |       |
| 4-Bromophenyl-phenylether  | 0.216 | 0.220  |            | 1.9  |       |
| Hexachlorobenzene          | 0.231 | 0.236  |            | 2.2  |       |
| Atrazine                   | 0.183 | 0.197  |            | 7.7  |       |
| Pentachlorophenol          | 0.132 | 0.147  |            | 11.4 | 20.0  |
| Phenanthrene               | 1.071 | 1.067  |            | -0.4 |       |
| Anthracene                 | 1.079 | 1.085  |            | 0.6  |       |
| Carbazole                  | 0.933 | 0.966  |            | 3.5  |       |
| Di-n-butylphthalate        | 0.950 | 1.072  |            | 12.8 |       |
| Fluoranthene               | 0.978 | 1.026  |            | 4.9  | 20.0  |
| Pyrene                     | 1.674 | 1.787  |            | 6.8  |       |
| Terphenyl-d14              | 1.115 | 1.214  |            | 8.9  |       |
| Butylbenzylphthalate       | 0.398 | 0.493  |            | 23.9 |       |
| 3,3-Dichlorobenzidine      | 0.337 | 0.364  |            | 8.0  |       |
| Benzo (a) anthracene       | 1.258 | 1.278  |            | 1.6  |       |
| Chrysene                   | 1.173 | 1.229  |            | 4.8  |       |
| Bis(2-ethylhexyl)phthalate | 0.528 | 0.593  |            | 12.3 |       |
| Di-n-octyl phthalate       | 1.017 | 0.974  |            | -4.2 | 20.0  |
| Benzo (b) fluoranthene     | 1.143 | 1.127  |            | -1.4 |       |
| Benzo (k) fluoranthene     | 1.068 | 1.128  |            | 5.6  |       |
| Benzo (a) pyrene           | 1.010 | 1.052  |            | 4.2  | 20.0  |
| Indeno (1,2,3-cd) pyrene   | 1.361 | 1.495  |            | 9.8  |       |
| Dibenzo (a,h) anthracene   | 1.114 | 1.224  |            | 9.9  |       |
| Benzo (g,h,i) perylene     | 1.103 | 1.199  |            | 8.7  |       |
| 1,2,4,5-Tetrachlorobenzene | 0.520 | 0.531  |            | 2.1  |       |
| 1,4-Dioxane                | 0.575 | 0.578  |            | 0.5  | 20.0  |
| 2,3,4,6-Tetrachlorophenol  | 0.257 | 0.297  |            | 15.6 |       |

All other compounds must meet a minimum RRF of 0.010.

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>AECO02</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q2924</u>                 |
| Instrument ID:  | <u>BNA_P</u>      | Calibration Date/Time: | <u>08/22/2025 10:49</u>      |
| Lab File ID:    | <u>BP025544.D</u> | Init. Calib. Date(s):  | <u>08/14/2025 08/14/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>12:33 17:24</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                    | RRF   | RRF040 | MIN RRF | %D   | MAX%D |
|-----------------------------|-------|--------|---------|------|-------|
| 2-Fluorophenol              | 1.204 | 1.201  |         | -0.2 |       |
| Benzaldehyde                | 1.082 | 1.377  |         | 27.3 |       |
| Phenol-d6                   | 1.544 | 1.586  |         | 2.7  |       |
| Phenol                      | 1.620 | 1.656  |         | 2.2  | 20.0  |
| bis(2-Chloroethyl)ether     | 1.263 | 1.266  |         | 0.2  |       |
| 2-Chlorophenol              | 1.359 | 1.376  |         | 1.3  |       |
| 2-Methylphenol              | 1.125 | 1.147  |         | 2.0  |       |
| 2,2-oxybis(1-Chloropropane) | 1.692 | 1.738  |         | 2.7  |       |
| Acetophenone                | 0.502 | 0.494  |         | -1.6 |       |
| 3+4-Methylphenols           | 1.560 | 1.584  |         | 1.5  |       |
| n-Nitroso-di-n-propylamine  | 1.023 | 1.030  | 0.050   | 0.7  |       |
| Nitrobenzene-d5             | 0.395 | 0.396  |         | 0.3  |       |
| Hexachloroethane            | 0.563 | 0.554  |         | -1.6 |       |
| Nitrobenzene                | 0.353 | 0.357  |         | 1.1  |       |
| Isophorone                  | 0.700 | 0.697  |         | -0.4 |       |
| 2-Nitrophenol               | 0.181 | 0.184  |         | 1.7  | 20.0  |
| 2,4-Dimethylphenol          | 0.312 | 0.307  |         | -1.6 |       |
| bis(2-Chloroethoxy)methane  | 0.423 | 0.416  |         | -1.7 |       |
| 2,4-Dichlorophenol          | 0.310 | 0.307  |         | -1.0 | 20.0  |
| Naphthalene                 | 1.040 | 1.006  |         | -3.3 |       |
| 4-Chloroaniline             | 0.459 | 0.460  |         | 0.2  |       |
| Hexachlorobutadiene         | 0.211 | 0.200  |         | -5.2 | 20.0  |
| Caprolactam                 | 0.114 | 0.124  |         | 8.8  |       |
| 4-Chloro-3-methylphenol     | 0.355 | 0.364  |         | 2.5  | 20.0  |
| 2-Methylnaphthalene         | 0.676 | 0.652  |         | -3.5 |       |
| Hexachlorocyclopentadiene   | 0.349 | 0.397  | 0.050   | 13.8 |       |
| 2,4,6-Trichlorophenol       | 0.390 | 0.387  |         | -0.8 | 20.0  |
| 2-Fluorobiphenyl            | 1.463 | 1.364  |         | -6.8 |       |
| 2,4,5-Trichlorophenol       | 0.427 | 0.430  |         | 0.7  |       |
| 1,1-Biphenyl                | 1.453 | 1.392  |         | -4.2 |       |
| 2-Chloronaphthalene         | 1.131 | 1.080  |         | -4.5 |       |
| 2-Nitroaniline              | 0.329 | 0.351  |         | 6.7  |       |
| Dimethylphthalate           | 1.465 | 1.455  |         | -0.7 |       |
| Acenaphthylene              | 1.871 | 1.834  |         | -2.0 |       |
| 2,6-Dinitrotoluene          | 0.309 | 0.324  |         | 4.9  |       |
| 3-Nitroaniline              | 0.347 | 0.370  |         | 6.6  |       |
| Acenaphthene                | 1.098 | 1.042  |         | -5.1 | 20.0  |
| 2,4-Dinitrophenol           | 0.165 | 0.228  | 0.050   | 38.2 |       |
| 4-Nitrophenol               | 0.285 | 0.311  | 0.050   | 9.1  |       |

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>AECO02</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q2924</u>                 |
| Instrument ID:  | <u>BNA_P</u>      | Calibration Date/Time: | <u>08/22/2025 10:49</u>      |
| Lab File ID:    | <u>BP025544.D</u> | Init. Calib. Date(s):  | <u>08/14/2025 08/14/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>12:33 17:24</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                   | RRF   | RRF040 | MIN RRF | %D   | MAX%D |
|----------------------------|-------|--------|---------|------|-------|
| Dibenzofuran               | 1.736 | 1.669  |         | -3.9 |       |
| 2,4-Dinitrotoluene         | 0.440 | 0.471  |         | 7.0  |       |
| Diethylphthalate           | 1.488 | 1.520  |         | 2.2  |       |
| 4-Chlorophenyl-phenylether | 0.681 | 0.653  |         | -4.1 |       |
| Fluorene                   | 1.370 | 1.350  |         | -1.5 |       |
| 4-Nitroaniline             | 0.356 | 0.382  |         | 7.3  |       |
| 4,6-Dinitro-2-methylphenol | 0.125 | 0.146  |         | 16.8 |       |
| n-Nitrosodiphenylamine     | 0.610 | 0.590  |         | -3.3 | 20.0  |
| 2,4,6-Tribromophenol       | 0.269 | 0.273  |         | 1.5  |       |
| 4-Bromophenyl-phenylether  | 0.220 | 0.209  |         | -5.0 |       |
| Hexachlorobenzene          | 0.255 | 0.238  |         | -6.7 |       |
| Atrazine                   | 0.228 | 0.228  |         | 0.0  |       |
| Pentachlorophenol          | 0.161 | 0.169  |         | 5.0  | 20.0  |
| Phenanthrene               | 1.099 | 1.058  |         | -3.7 |       |
| Anthracene                 | 1.122 | 1.113  |         | -0.8 |       |
| Carbazole                  | 1.057 | 1.064  |         | 0.7  |       |
| Di-n-butylphthalate        | 1.300 | 1.314  |         | 1.1  |       |
| Fluoranthene               | 1.311 | 1.296  |         | -1.1 | 20.0  |
| Pyrene                     | 1.300 | 1.255  |         | -3.5 |       |
| Terphenyl-d14              | 1.093 | 1.041  |         | -4.8 |       |
| Butylbenzylphthalate       | 0.589 | 0.607  |         | 3.1  |       |
| 3,3-Dichlorobenzidine      | 0.497 | 0.497  |         | 0.0  |       |
| Benzo (a) anthracene       | 1.319 | 1.273  |         | -3.5 |       |
| Chrysene                   | 1.235 | 1.201  |         | -2.8 |       |
| Bis(2-ethylhexyl)phthalate | 0.867 | 0.866  |         | -0.1 |       |
| Di-n-octyl phthalate       | 1.492 | 1.541  |         | 3.3  | 20.0  |
| Benzo (b) fluoranthene     | 1.179 | 1.115  |         | -5.4 |       |
| Benzo (k) fluoranthene     | 1.180 | 1.139  |         | -3.5 |       |
| Benzo (a) pyrene           | 1.148 | 1.107  |         | -3.6 | 20.0  |
| Indeno (1,2,3-cd) pyrene   | 1.467 | 1.402  |         | -4.4 |       |
| Dibenzo (a,h) anthracene   | 1.199 | 1.138  |         | -5.1 |       |
| Benzo (g,h,i) perylene     | 1.178 | 1.141  |         | -3.1 |       |
| 1,2,4,5-Tetrachlorobenzene | 0.578 | 0.548  |         | -5.2 |       |
| 1,4-Dioxane                | 0.472 | 0.461  |         | -2.3 | 20.0  |
| 2,3,4,6-Tetrachlorophenol  | 0.377 | 0.381  |         | 1.1  |       |

All other compounds must meet a minimum RRF of 0.010.

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>AEC002</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q2924</u>                 |
| Instrument ID:  | <u>BNA_P</u>      | Calibration Date/Time: | <u>08/25/2025 12:13</u>      |
| Lab File ID:    | <u>BP025561.D</u> | Init. Calib. Date(s):  | <u>08/14/2025 08/14/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>12:33 17:24</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                    | RRF   | RRF040 | MIN RRF | %D   | MAX%D |
|-----------------------------|-------|--------|---------|------|-------|
| 2-Fluorophenol              | 1.204 | 1.230  |         | 2.2  |       |
| Benzaldehyde                | 1.082 | 1.338  |         | 23.7 |       |
| Phenol-d6                   | 1.544 | 1.578  |         | 2.2  |       |
| Phenol                      | 1.620 | 1.637  |         | 1.0  | 20.0  |
| bis(2-Chloroethyl)ether     | 1.263 | 1.249  |         | -1.1 |       |
| 2-Chlorophenol              | 1.359 | 1.362  |         | 0.2  |       |
| 2-Methylphenol              | 1.125 | 1.108  |         | -1.5 |       |
| 2,2-oxybis(1-Chloropropane) | 1.692 | 1.716  |         | 1.4  |       |
| Acetophenone                | 0.502 | 0.490  |         | -2.4 |       |
| 3+4-Methylphenols           | 1.560 | 1.501  |         | -3.8 |       |
| n-Nitroso-di-n-propylamine  | 1.023 | 0.970  | 0.050   | -5.2 |       |
| Nitrobenzene-d5             | 0.395 | 0.399  |         | 1.0  |       |
| Hexachloroethane            | 0.563 | 0.547  |         | -2.8 |       |
| Nitrobenzene                | 0.353 | 0.358  |         | 1.4  |       |
| Isophorone                  | 0.700 | 0.668  |         | -4.6 |       |
| 2-Nitrophenol               | 0.181 | 0.186  |         | 2.8  | 20.0  |
| 2,4-Dimethylphenol          | 0.312 | 0.304  |         | -2.6 |       |
| bis(2-Chloroethoxy)methane  | 0.423 | 0.405  |         | -4.3 |       |
| 2,4-Dichlorophenol          | 0.310 | 0.306  |         | -1.3 | 20.0  |
| Naphthalene                 | 1.040 | 1.011  |         | -2.8 |       |
| 4-Chloroaniline             | 0.459 | 0.446  |         | -2.8 |       |
| Hexachlorobutadiene         | 0.211 | 0.197  |         | -6.6 | 20.0  |
| Caprolactam                 | 0.114 | 0.111  |         | -2.6 |       |
| 4-Chloro-3-methylphenol     | 0.355 | 0.342  |         | -3.7 | 20.0  |
| 2-Methylnaphthalene         | 0.676 | 0.636  |         | -5.9 |       |
| Hexachlorocyclopentadiene   | 0.349 | 0.387  | 0.050   | 10.9 |       |
| 2,4,6-Trichlorophenol       | 0.390 | 0.396  |         | 1.5  | 20.0  |
| 2-Fluorobiphenyl            | 1.463 | 1.453  |         | -0.7 |       |
| 2,4,5-Trichlorophenol       | 0.427 | 0.435  |         | 1.9  |       |
| 1,1-Biphenyl                | 1.453 | 1.456  |         | 0.2  |       |
| 2-Chloronaphthalene         | 1.131 | 1.146  |         | 1.3  |       |
| 2-Nitroaniline              | 0.329 | 0.351  |         | 6.7  |       |
| Dimethylphthalate           | 1.465 | 1.428  |         | -2.5 |       |
| Acenaphthylene              | 1.871 | 1.872  |         | 0.1  |       |
| 2,6-Dinitrotoluene          | 0.309 | 0.318  |         | 2.9  |       |
| 3-Nitroaniline              | 0.347 | 0.364  |         | 4.9  |       |
| Acenaphthene                | 1.098 | 1.074  |         | -2.2 | 20.0  |
| 2,4-Dinitrophenol           | 0.165 | 0.220  | 0.050   | 33.3 |       |
| 4-Nitrophenol               | 0.285 | 0.297  | 0.050   | 4.2  |       |

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>AECO02</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q2924</u>                 |
| Instrument ID:  | <u>BNA_P</u>      | Calibration Date/Time: | <u>08/25/2025 12:13</u>      |
| Lab File ID:    | <u>BP025561.D</u> | Init. Calib. Date(s):  | <u>08/14/2025 08/14/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>12:33 17:24</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                   | RRF   | RRF040 | MIN<br>RRF | %D   | MAX%D |
|----------------------------|-------|--------|------------|------|-------|
| Dibenzofuran               | 1.736 | 1.714  |            | -1.3 |       |
| 2,4-Dinitrotoluene         | 0.440 | 0.461  |            | 4.8  |       |
| Diethylphthalate           | 1.488 | 1.451  |            | -2.5 |       |
| 4-Chlorophenyl-phenylether | 0.681 | 0.659  |            | -3.2 |       |
| Fluorene                   | 1.370 | 1.353  |            | -1.2 |       |
| 4-Nitroaniline             | 0.356 | 0.380  |            | 6.7  |       |
| 4,6-Dinitro-2-methylphenol | 0.125 | 0.147  |            | 17.6 |       |
| n-Nitrosodiphenylamine     | 0.610 | 0.601  |            | -1.5 | 20.0  |
| 2,4,6-Tribromophenol       | 0.269 | 0.270  |            | 0.4  |       |
| 4-Bromophenyl-phenylether  | 0.220 | 0.213  |            | -3.2 |       |
| Hexachlorobenzene          | 0.255 | 0.243  |            | -4.7 |       |
| Atrazine                   | 0.228 | 0.221  |            | -3.1 |       |
| Pentachlorophenol          | 0.161 | 0.156  |            | -3.1 | 20.0  |
| Phenanthrene               | 1.099 | 1.083  |            | -1.5 |       |
| Anthracene                 | 1.122 | 1.113  |            | -0.8 |       |
| Carbazole                  | 1.057 | 1.061  |            | 0.4  |       |
| Di-n-butylphthalate        | 1.300 | 1.268  |            | -2.5 |       |
| Fluoranthene               | 1.311 | 1.276  |            | -2.7 | 20.0  |
| Pyrene                     | 1.300 | 1.273  |            | -2.1 |       |
| Terphenyl-d14              | 1.093 | 1.041  |            | -4.8 |       |
| Butylbenzylphthalate       | 0.589 | 0.574  |            | -2.5 |       |
| 3,3-Dichlorobenzidine      | 0.497 | 0.487  |            | -2.0 |       |
| Benzo (a) anthracene       | 1.319 | 1.282  |            | -2.8 |       |
| Chrysene                   | 1.235 | 1.205  |            | -2.4 |       |
| Bis(2-ethylhexyl)phthalate | 0.867 | 0.799  |            | -7.8 |       |
| Di-n-octyl phthalate       | 1.492 | 1.388  |            | -7.0 | 20.0  |
| Benzo (b) fluoranthene     | 1.179 | 1.144  |            | -3.0 |       |
| Benzo (k) fluoranthene     | 1.180 | 1.167  |            | -1.1 |       |
| Benzo (a) pyrene           | 1.148 | 1.140  |            | -0.7 | 20.0  |
| Indeno (1,2,3-cd) pyrene   | 1.467 | 1.451  |            | -1.1 |       |
| Dibenzo (a,h) anthracene   | 1.199 | 1.203  |            | 0.3  |       |
| Benzo (g,h,i) perylene     | 1.178 | 1.174  |            | -0.3 |       |
| 1,2,4,5-Tetrachlorobenzene | 0.578 | 0.586  |            | 1.4  |       |
| 1,4-Dioxane                | 0.472 | 0.466  |            | -1.3 | 20.0  |
| 2,3,4,6-Tetrachlorophenol  | 0.377 | 0.380  |            | 0.8  |       |

All other compounds must meet a minimum RRF of 0.010.

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>AEC002</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q2924</u>                 |
| Instrument ID:  | <u>BNA_P</u>      | Calibration Date/Time: | <u>08/26/2025 10:21</u>      |
| Lab File ID:    | <u>BP025579.D</u> | Init. Calib. Date(s):  | <u>08/14/2025 08/14/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>12:33 17:24</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                    | RRF   | RRF040 | MIN RRF | %D   | MAX%D |
|-----------------------------|-------|--------|---------|------|-------|
| 2-Fluorophenol              | 1.204 | 1.238  |         | 2.8  |       |
| Benzaldehyde                | 1.082 | 1.311  |         | 21.2 |       |
| Phenol-d6                   | 1.544 | 1.536  |         | -0.5 |       |
| Phenol                      | 1.620 | 1.579  |         | -2.5 | 20.0  |
| bis(2-Chloroethyl)ether     | 1.263 | 1.232  |         | -2.5 |       |
| 2-Chlorophenol              | 1.359 | 1.354  |         | -0.4 |       |
| 2-Methylphenol              | 1.125 | 1.095  |         | -2.7 |       |
| 2,2-oxybis(1-Chloropropane) | 1.692 | 1.682  |         | -0.6 |       |
| Acetophenone                | 0.502 | 0.498  |         | -0.8 |       |
| 3+4-Methylphenols           | 1.560 | 1.460  |         | -6.4 |       |
| n-Nitroso-di-n-propylamine  | 1.023 | 0.944  | 0.050   | -7.7 |       |
| Nitrobenzene-d5             | 0.395 | 0.403  |         | 2.0  |       |
| Hexachloroethane            | 0.563 | 0.567  |         | 0.7  |       |
| Nitrobenzene                | 0.353 | 0.356  |         | 0.9  |       |
| Isophorone                  | 0.700 | 0.679  |         | -3.0 |       |
| 2-Nitrophenol               | 0.181 | 0.189  |         | 4.4  | 20.0  |
| 2,4-Dimethylphenol          | 0.312 | 0.308  |         | -1.3 |       |
| bis(2-Chloroethoxy)methane  | 0.423 | 0.409  |         | -3.3 |       |
| 2,4-Dichlorophenol          | 0.310 | 0.308  |         | -0.6 | 20.0  |
| Naphthalene                 | 1.040 | 1.016  |         | -2.3 |       |
| 4-Chloroaniline             | 0.459 | 0.440  |         | -4.1 |       |
| Hexachlorobutadiene         | 0.211 | 0.212  |         | 0.5  | 20.0  |
| Caprolactam                 | 0.114 | 0.106  |         | -7.0 |       |
| 4-Chloro-3-methylphenol     | 0.355 | 0.344  |         | -3.1 | 20.0  |
| 2-Methylnaphthalene         | 0.676 | 0.650  |         | -3.8 |       |
| Hexachlorocyclopentadiene   | 0.349 | 0.378  | 0.050   | 8.3  |       |
| 2,4,6-Trichlorophenol       | 0.390 | 0.414  |         | 6.2  | 20.0  |
| 2-Fluorobiphenyl            | 1.463 | 1.454  |         | -0.6 |       |
| 2,4,5-Trichlorophenol       | 0.427 | 0.455  |         | 6.6  |       |
| 1,1-Biphenyl                | 1.453 | 1.465  |         | 0.8  |       |
| 2-Chloronaphthalene         | 1.131 | 1.145  |         | 1.2  |       |
| 2-Nitroaniline              | 0.329 | 0.352  |         | 7.0  |       |
| Dimethylphthalate           | 1.465 | 1.457  |         | -0.5 |       |
| Acenaphthylene              | 1.871 | 1.885  |         | 0.7  |       |
| 2,6-Dinitrotoluene          | 0.309 | 0.317  |         | 2.6  |       |
| 3-Nitroaniline              | 0.347 | 0.352  |         | 1.4  |       |
| Acenaphthene                | 1.098 | 1.071  |         | -2.5 | 20.0  |
| 2,4-Dinitrophenol           | 0.165 | 0.206  | 0.050   | 24.8 |       |
| 4-Nitrophenol               | 0.285 | 0.275  | 0.050   | -3.5 |       |

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>AECO02</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q2924</u>                 |
| Instrument ID:  | <u>BNA_P</u>      | Calibration Date/Time: | <u>08/26/2025 10:21</u>      |
| Lab File ID:    | <u>BP025579.D</u> | Init. Calib. Date(s):  | <u>08/14/2025 08/14/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>12:33 17:24</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                   | RRF   | RRF040 | MIN RRF | %D   | MAX%D |
|----------------------------|-------|--------|---------|------|-------|
| Dibenzofuran               | 1.736 | 1.708  |         | -1.6 |       |
| 2,4-Dinitrotoluene         | 0.440 | 0.449  |         | 2.0  |       |
| Diethylphthalate           | 1.488 | 1.476  |         | -0.8 |       |
| 4-Chlorophenyl-phenylether | 0.681 | 0.673  |         | -1.2 |       |
| Fluorene                   | 1.370 | 1.330  |         | -2.9 |       |
| 4-Nitroaniline             | 0.356 | 0.360  |         | 1.1  |       |
| 4,6-Dinitro-2-methylphenol | 0.125 | 0.142  |         | 13.6 |       |
| n-Nitrosodiphenylamine     | 0.610 | 0.600  |         | -1.6 | 20.0  |
| 2,4,6-Tribromophenol       | 0.269 | 0.279  |         | 3.7  |       |
| 4-Bromophenyl-phenylether  | 0.220 | 0.223  |         | 1.4  |       |
| Hexachlorobenzene          | 0.255 | 0.259  |         | 1.6  |       |
| Atrazine                   | 0.228 | 0.224  |         | -1.8 |       |
| Pentachlorophenol          | 0.161 | 0.160  |         | -0.6 | 20.0  |
| Phenanthrene               | 1.099 | 1.075  |         | -2.2 |       |
| Anthracene                 | 1.122 | 1.109  |         | -1.2 |       |
| Carbazole                  | 1.057 | 1.018  |         | -3.7 |       |
| Di-n-butylphthalate        | 1.300 | 1.288  |         | -0.9 |       |
| Fluoranthene               | 1.311 | 1.234  |         | -5.9 | 20.0  |
| Pyrene                     | 1.300 | 1.297  |         | -0.2 |       |
| Terphenyl-d14              | 1.093 | 1.098  |         | 0.5  |       |
| Butylbenzylphthalate       | 0.589 | 0.609  |         | 3.4  |       |
| 3,3-Dichlorobenzidine      | 0.497 | 0.499  |         | 0.4  |       |
| Benzo (a) anthracene       | 1.319 | 1.294  |         | -1.9 |       |
| Chrysene                   | 1.235 | 1.223  |         | -1.0 |       |
| Bis(2-ethylhexyl)phthalate | 0.867 | 0.871  |         | 0.5  |       |
| Di-n-octyl phthalate       | 1.492 | 1.545  |         | 3.6  | 20.0  |
| Benzo (b) fluoranthene     | 1.179 | 1.158  |         | -1.8 |       |
| Benzo (k) fluoranthene     | 1.180 | 1.141  |         | -3.3 |       |
| Benzo (a) pyrene           | 1.148 | 1.118  |         | -2.6 | 20.0  |
| Indeno (1,2,3-cd) pyrene   | 1.467 | 1.442  |         | -1.7 |       |
| Dibenzo (a,h) anthracene   | 1.199 | 1.179  |         | -1.7 |       |
| Benzo (g,h,i) perylene     | 1.178 | 1.147  |         | -2.6 |       |
| 1,2,4,5-Tetrachlorobenzene | 0.578 | 0.600  |         | 3.8  |       |
| 1,4-Dioxane                | 0.472 | 0.478  |         | 1.3  | 20.0  |
| 2,3,4,6-Tetrachlorophenol  | 0.377 | 0.382  |         | 1.3  |       |

All other compounds must meet a minimum RRF of 0.010.

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>AEC002</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q2924</u>                 |
| Instrument ID:  | <u>BNA_P</u>      | Calibration Date/Time: | <u>08/27/2025 09:39</u>      |
| Lab File ID:    | <u>BP025596.D</u> | Init. Calib. Date(s):  | <u>08/14/2025 08/14/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>12:33 17:24</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                    | RRF   | RRF040 | MIN<br>RRF | %D    | MAX%D |
|-----------------------------|-------|--------|------------|-------|-------|
| 2-Fluorophenol              | 1.204 | 1.204  |            | 0.0   |       |
| Benzaldehyde                | 1.082 | 1.320  |            | 22.0  |       |
| Phenol-d6                   | 1.544 | 1.563  |            | 1.2   |       |
| Phenol                      | 1.620 | 1.636  |            | 1.0   | 20.0  |
| bis(2-Chloroethyl)ether     | 1.263 | 1.242  |            | -1.7  |       |
| 2-Chlorophenol              | 1.359 | 1.356  |            | -0.2  |       |
| 2-Methylphenol              | 1.125 | 1.127  |            | 0.2   |       |
| 2,2-oxybis(1-Chloropropane) | 1.692 | 1.723  |            | 1.8   |       |
| Acetophenone                | 0.502 | 0.496  |            | -1.2  |       |
| 3+4-Methylphenols           | 1.560 | 1.513  |            | -3.0  |       |
| n-Nitroso-di-n-propylamine  | 1.023 | 1.032  | 0.050      | 0.9   |       |
| Nitrobenzene-d5             | 0.395 | 0.395  |            | 0.0   |       |
| Hexachloroethane            | 0.563 | 0.563  |            | 0.0   |       |
| Nitrobenzene                | 0.353 | 0.352  |            | -0.3  |       |
| Isophorone                  | 0.700 | 0.708  |            | 1.1   |       |
| 2-Nitrophenol               | 0.181 | 0.187  |            | 3.3   | 20.0  |
| 2,4-Dimethylphenol          | 0.312 | 0.310  |            | -0.6  |       |
| bis(2-Chloroethoxy)methane  | 0.423 | 0.417  |            | -1.4  |       |
| 2,4-Dichlorophenol          | 0.310 | 0.310  |            | 0.0   | 20.0  |
| Naphthalene                 | 1.040 | 1.018  |            | -2.1  |       |
| 4-Chloroaniline             | 0.459 | 0.458  |            | -0.2  |       |
| Hexachlorobutadiene         | 0.211 | 0.206  |            | -2.4  | 20.0  |
| Caprolactam                 | 0.114 | 0.121  |            | 6.1   |       |
| 4-Chloro-3-methylphenol     | 0.355 | 0.363  |            | 2.3   | 20.0  |
| 2-Methylnaphthalene         | 0.676 | 0.657  |            | -2.8  |       |
| Hexachlorocyclopentadiene   | 0.349 | 0.280  | 0.050      | -19.8 |       |
| 2,4,6-Trichlorophenol       | 0.390 | 0.385  |            | -1.3  | 20.0  |
| 2-Fluorobiphenyl            | 1.463 | 1.415  |            | -3.3  |       |
| 2,4,5-Trichlorophenol       | 0.427 | 0.424  |            | -0.7  |       |
| 1,1-Biphenyl                | 1.453 | 1.400  |            | -3.6  |       |
| 2-Chloronaphthalene         | 1.131 | 1.094  |            | -3.3  |       |
| 2-Nitroaniline              | 0.329 | 0.360  |            | 9.4   |       |
| Dimethylphthalate           | 1.465 | 1.499  |            | 2.3   |       |
| Acenaphthylene              | 1.871 | 1.827  |            | -2.4  |       |
| 2,6-Dinitrotoluene          | 0.309 | 0.331  |            | 7.1   |       |
| 3-Nitroaniline              | 0.347 | 0.364  |            | 4.9   |       |
| Acenaphthene                | 1.098 | 1.046  |            | -4.7  | 20.0  |
| 2,4-Dinitrophenol           | 0.165 | 0.204  | 0.050      | 23.6  |       |
| 4-Nitrophenol               | 0.285 | 0.275  | 0.050      | -3.5  |       |

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                   |                        |                              |
|-----------------|-------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>   | Contract:              | <u>AECO02</u>                |
| Lab Code:       | <u>ACE</u>        | SDG No.:               | <u>Q2924</u>                 |
| Instrument ID:  | <u>BNA_P</u>      | Calibration Date/Time: | <u>08/27/2025 09:39</u>      |
| Lab File ID:    | <u>BP025596.D</u> | Init. Calib. Date(s):  | <u>08/14/2025 08/14/2025</u> |
| EPA Sample No.: | <u>SSTDCCC040</u> | Init. Calib. Time(s):  | <u>12:33 17:24</u>           |
| GC Column:      | <u>ZB-GR</u>      | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                   | RRF   | RRF040 | MIN RRF | %D   | MAX%D |
|----------------------------|-------|--------|---------|------|-------|
| Dibenzofuran               | 1.736 | 1.718  |         | -1.0 |       |
| 2,4-Dinitrotoluene         | 0.440 | 0.481  |         | 9.3  |       |
| Diethylphthalate           | 1.488 | 1.550  |         | 4.2  |       |
| 4-Chlorophenyl-phenylether | 0.681 | 0.690  |         | 1.3  |       |
| Fluorene                   | 1.370 | 1.386  |         | 1.2  |       |
| 4-Nitroaniline             | 0.356 | 0.391  |         | 9.8  |       |
| 4,6-Dinitro-2-methylphenol | 0.125 | 0.142  |         | 13.6 |       |
| n-Nitrosodiphenylamine     | 0.610 | 0.593  |         | -2.8 | 20.0  |
| 2,4,6-Tribromophenol       | 0.269 | 0.296  |         | 10.0 |       |
| 4-Bromophenyl-phenylether  | 0.220 | 0.217  |         | -1.4 |       |
| Hexachlorobenzene          | 0.255 | 0.256  |         | 0.4  |       |
| Atrazine                   | 0.228 | 0.234  |         | 2.6  |       |
| Pentachlorophenol          | 0.161 | 0.155  |         | -3.7 | 20.0  |
| Phenanthrene               | 1.099 | 1.066  |         | -3.0 |       |
| Anthracene                 | 1.122 | 1.113  |         | -0.8 |       |
| Carbazole                  | 1.057 | 1.047  |         | -0.9 |       |
| Di-n-butylphthalate        | 1.300 | 1.367  |         | 5.2  |       |
| Fluoranthene               | 1.311 | 1.324  |         | 1.0  | 20.0  |
| Pyrene                     | 1.300 | 1.296  |         | -0.3 |       |
| Terphenyl-d14              | 1.093 | 1.122  |         | 2.7  |       |
| Butylbenzylphthalate       | 0.589 | 0.630  |         | 7.0  |       |
| 3,3-Dichlorobenzidine      | 0.497 | 0.494  |         | -0.6 |       |
| Benzo (a) anthracene       | 1.319 | 1.299  |         | -1.5 |       |
| Chrysene                   | 1.235 | 1.231  |         | -0.3 |       |
| Bis(2-ethylhexyl)phthalate | 0.867 | 0.884  |         | 2.0  |       |
| Di-n-octyl phthalate       | 1.492 | 1.545  |         | 3.6  | 20.0  |
| Benzo (b) fluoranthene     | 1.179 | 1.165  |         | -1.2 |       |
| Benzo (k) fluoranthene     | 1.180 | 1.169  |         | -0.9 |       |
| Benzo (a) pyrene           | 1.148 | 1.139  |         | -0.8 | 20.0  |
| Indeno (1,2,3-cd) pyrene   | 1.467 | 1.427  |         | -2.7 |       |
| Dibenzo (a,h) anthracene   | 1.199 | 1.176  |         | -1.9 |       |
| Benzo (g,h,i) perylene     | 1.178 | 1.148  |         | -2.5 |       |
| 1,2,4,5-Tetrachlorobenzene | 0.578 | 0.565  |         | -2.2 |       |
| 1,4-Dioxane                | 0.472 | 0.463  |         | -1.9 | 20.0  |
| 2,3,4,6-Tetrachlorophenol  | 0.377 | 0.394  |         | 4.5  |       |

All other compounds must meet a minimum RRF of 0.010.