

## LAB CHRONICLE

|                 |                      |                   |                               |
|-----------------|----------------------|-------------------|-------------------------------|
| <b>OrderID:</b> | Q2881                | <b>OrderDate:</b> | 8/15/2025 10:55:00 AM         |
| <b>Client:</b>  | Tetra Tech NUS, Inc. | <b>Project:</b>   | NWIRP Bethpage 112G08005-WE13 |
| <b>Contact:</b> | Ernie Wu             | <b>Location:</b>  | J32                           |

| LabID           | ClientID                             | Matrix       | Test           | Method        | Sample Date     | Prep Date | Anal Date | Received        |
|-----------------|--------------------------------------|--------------|----------------|---------------|-----------------|-----------|-----------|-----------------|
| <b>Q2881-01</b> | <b>RW8-SP100-2025081</b><br><b>4</b> | <b>Water</b> |                |               | <b>08/14/25</b> |           |           | <b>08/15/25</b> |
|                 |                                      |              | SVOC-SIMGroup1 | 8270-Modified |                 | 08/18/25  | 08/20/25  |                 |
| <b>Q2881-02</b> | <b>RW8-SP303-2025081</b><br><b>4</b> | <b>Water</b> |                |               | <b>08/14/25</b> |           |           | <b>08/15/25</b> |
|                 |                                      |              | SVOC-SIMGroup1 | 8270-Modified |                 | 08/18/25  | 08/20/25  |                 |



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### Hit Summary Sheet SW-846

**SDG No.:** Q2881  
**Client:** Tetra Tech NUS, Inc.

| Sample ID            | Client ID          | Parameter | Concentration | C     | MDL | LOD  | RDL | Units    |
|----------------------|--------------------|-----------|---------------|-------|-----|------|-----|----------|
| Client ID :          | RW8-SP100-20250814 |           |               |       |     |      |     |          |
| Q2881-01             | RW8-SP100-20250814 | WATER     | 1,4-Dioxane   | 0.070 | J   | 0.07 | 0.2 | 0.2 ug/L |
| Total Svoc :         |                    |           |               | 0.07  |     |      |     |          |
| Total Concentration: |                    |           |               | 0.07  |     |      |     |          |



# SAMPLE DATA

## Report of Analysis

|                    |                               |                 |                |
|--------------------|-------------------------------|-----------------|----------------|
| Client:            | Tetra Tech NUS, Inc.          | Date Collected: | 08/14/25       |
| Project:           | NWIRP Bethpage 112G08005-WE13 | Date Received:  | 08/15/25       |
| Client Sample ID:  | RW8-SP100-20250814            | SDG No.:        | Q2881          |
| Lab Sample ID:     | Q2881-01                      | Matrix:         | Water          |
| Analytical Method: | SW8270ESIM                    | % Solid:        | 0              |
| Sample Wt/Vol:     | 1000 Units: mL                | Final Vol:      | 1000 uL        |
| Soil Aliquot Vol:  | uL                            | Test:           | SVOC-SIMGroup1 |
| 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 |
| BN037623.D        | 1         | 08/18/25 08:41 | 08/20/25 07:13 | PB169286      |

| CAS Number                | Parameter               | Conc. | Qualifier | MDL      | LOD  | LOQ / CRQL | Units    |
|---------------------------|-------------------------|-------|-----------|----------|------|------------|----------|
| <b>TARGETS</b>            |                         |       |           |          |      |            |          |
| 123-91-1                  | 1,4-Dioxane             | 0.070 | J         | 0.070    | 0.20 | 0.20       | ug/L     |
| <b>SURROGATES</b>         |                         |       |           |          |      |            |          |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.26  |           | 30 - 150 |      | 65%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.37  |           | 30 - 150 |      | 92%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.32  |           | 55 - 111 |      | 79%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.31  |           | 53 - 106 |      | 77%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.42  |           | 58 - 132 |      | 105%       | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 2130  | 7.717     |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 4990  | 10.498    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 2390  | 14.345    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 5110  | 17.086    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 4520  | 21.268    |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 4090  | 23.499    |          |      |            |          |

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:            | Tetra Tech NUS, Inc.          | Date Collected: | 08/14/25       |
| Project:           | NWIRP Bethpage 112G08005-WE13 | Date Received:  | 08/15/25       |
| Client Sample ID:  | RW8-SP303-20250814            | SDG No.:        | Q2881          |
| Lab Sample ID:     | Q2881-02                      | Matrix:         | Water          |
| Analytical Method: | SW8270ESIM                    | % Solid:        | 0              |
| Sample Wt/Vol:     | 1000 Units: mL                | Final Vol:      | 1000 uL        |
| Soil Aliquot Vol:  | uL                            | Test:           | SVOC-SIMGroup1 |
| 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 |
| BN037624.D        | 1         | 08/18/25 08:41 | 08/20/25 07:49 | PB169286      |

| CAS Number                | Parameter               | Conc. | Qualifier | MDL      | LOD  | LOQ / CRQL | Units    |
|---------------------------|-------------------------|-------|-----------|----------|------|------------|----------|
| <b>TARGETS</b>            |                         |       |           |          |      |            |          |
| 123-91-1                  | 1,4-Dioxane             | 0.20  | U         | 0.070    | 0.20 | 0.20       | ug/L     |
| <b>SURROGATES</b>         |                         |       |           |          |      |            |          |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.27  |           | 30 - 150 |      | 68%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.37  |           | 30 - 150 |      | 93%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.33  |           | 55 - 111 |      | 81%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.30  |           | 53 - 106 |      | 75%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.43  |           | 58 - 132 |      | 107%       | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 2200  | 7.71      |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 5180  | 10.498    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 2520  | 14.345    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 5410  | 17.086    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 4670  | 21.268    |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 4150  | 23.502    |          |      |            |          |

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.: Q2881

Client: Tetra Tech NUS, Inc.

Analytical Method: 8270-Modified

| Lab Sample ID | Client ID               | Parameter               | Spike (PPM) | Result (PPM) | Recovery (%) | Qual | Limits (%) |      |
|---------------|-------------------------|-------------------------|-------------|--------------|--------------|------|------------|------|
|               |                         |                         |             |              |              |      | Low        | High |
| PB169286BL    | PB169286BL              | 2-Methylnaphthalene-d10 | 0.4         | 0.32         | 81           |      | 30         | 150  |
|               |                         | Fluoranthene-d10        | 0.4         | 0.33         | 83           |      | 30         | 150  |
|               |                         | Nitrobenzene-d5         | 0.4         | 0.36         | 90           |      | 55         | 111  |
|               |                         | 2-Fluorobiphenyl        | 0.4         | 0.35         | 86           |      | 53         | 106  |
|               |                         | Terphenyl-d14           | 0.4         | 0.37         | 93           |      | 58         | 132  |
| PB169286BS    | PB169286BS              | 2-Methylnaphthalene-d10 | 0.4         | 0.33         | 82           |      | 30         | 150  |
|               |                         | Fluoranthene-d10        | 0.4         | 0.31         | 76           |      | 30         | 150  |
|               |                         | Nitrobenzene-d5         | 0.4         | 0.35         | 87           |      | 55         | 111  |
|               |                         | 2-Fluorobiphenyl        | 0.4         | 0.36         | 90           |      | 53         | 106  |
|               |                         | Terphenyl-d14           | 0.4         | 0.35         | 88           |      | 58         | 132  |
| Q2842-04MS    | MW-17B-55.5-08122025MS  | 2-Methylnaphthalene-d10 | 0.4         | 0.31         | 77           |      | 30         | 150  |
|               |                         | Fluoranthene-d10        | 0.4         | 0.38         | 94           |      | 30         | 150  |
|               |                         | Nitrobenzene-d5         | 0.4         | 0.33         | 83           |      | 55         | 111  |
|               |                         | 2-Fluorobiphenyl        | 0.4         | 0.32         | 81           |      | 53         | 106  |
|               |                         | Terphenyl-d14           | 0.4         | 0.47         | 117          |      | 58         | 132  |
| Q2842-05MSD   | MW-17B-55.5-08122025MSD | 2-Methylnaphthalene-d10 | 0.4         | 0.30         | 76           |      | 30         | 150  |
|               |                         | Fluoranthene-d10        | 0.4         | 0.37         | 93           |      | 30         | 150  |
|               |                         | Nitrobenzene-d5         | 0.4         | 0.32         | 79           |      | 55         | 111  |
|               |                         | 2-Fluorobiphenyl        | 0.4         | 0.32         | 81           |      | 53         | 106  |
|               |                         | Terphenyl-d14           | 0.4         | 0.47         | 117          |      | 58         | 132  |
| Q2881-01      | RW8-SP100-20250814      | 2-Methylnaphthalene-d10 | 0.4         | 0.26         | 65           |      | 30         | 150  |
|               |                         | Fluoranthene-d10        | 0.4         | 0.37         | 92           |      | 30         | 150  |
|               |                         | Nitrobenzene-d5         | 0.4         | 0.32         | 79           |      | 55         | 111  |
|               |                         | 2-Fluorobiphenyl        | 0.4         | 0.31         | 77           |      | 53         | 106  |
|               |                         | Terphenyl-d14           | 0.4         | 0.42         | 105          |      | 58         | 132  |
| Q2881-02      | RW8-SP303-20250814      | 2-Methylnaphthalene-d10 | 0.4         | 0.27         | 68           |      | 30         | 150  |
|               |                         | Fluoranthene-d10        | 0.4         | 0.37         | 93           |      | 30         | 150  |
|               |                         | Nitrobenzene-d5         | 0.4         | 0.33         | 81           |      | 55         | 111  |
|               |                         | 2-Fluorobiphenyl        | 0.4         | 0.30         | 75           |      | 53         | 106  |
|               |                         | Terphenyl-d14           | 0.4         | 0.43         | 107          |      | 58         | 132  |

Matrix Spike/Matrix Spike Duplicate Summary  
SW-846

|          |                      |                    |                 |
|----------|----------------------|--------------------|-----------------|
| SDG No.: | Q2881                | Analytical Method: | SW8270-Modified |
| Client:  | Tetra Tech NUS, Inc. | DataFile:          | BN037605.D      |

| Parameter      | Spike      | Sample Result     | Result                 | Units | Rec | Rec Qual | RPD | RPD Qual | Low | Limits High | RPD |
|----------------|------------|-------------------|------------------------|-------|-----|----------|-----|----------|-----|-------------|-----|
| Lab Sample ID: | Q2842-04MS | Client Sample ID: | MW-17B-55.5-08122025MS |       |     |          |     |          |     |             |     |
| 1,4-Dioxane    | 0.43       | 2.90              | 3.00                   | ug/L  | 23  | *        |     |          | 70  | 130         |     |

Matrix Spike/Matrix Spike Duplicate Summary  
SW-846

SDG No.:

Q2881

Analytical Method:

SW8270-Modified

Client:

Tetra Tech NUS, Inc.

DataFile:

BN037606.D

| Parameter      | Spike       | Sample Result     | Result                  | Units | Rec | Rec Qual | RPD | RPD Qual | Low | Limits High | RPD |
|----------------|-------------|-------------------|-------------------------|-------|-----|----------|-----|----------|-----|-------------|-----|
| Lab Sample ID: | Q2842-05MSD | Client Sample ID: | MW-17B-55.5-08122025MSD |       |     |          |     |          |     |             |     |
| 1,4-Dioxane    | 0.42        | 2.90              | 3.00                    | ug/L  | 24  | *        | 4   |          | 70  | 130         | 20  |

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

SW-846

SDG No.: Q2881

Analytical Method: 8270-Modified

Client: Tetra Tech NUS, Inc.

DataFile: BN037630.D

| Lab Sample ID | Parameter   | Spike | Result | Unit | Rec | RPD | Qual | RPD  | Limits |      | RPD |
|---------------|-------------|-------|--------|------|-----|-----|------|------|--------|------|-----|
|               |             |       |        |      |     |     |      | Qual | Low    | High |     |
| PB169286BS    | 1,4-Dioxane | 0.4   | 0.30   | ug/L | 75  |     |      |      | 70     | 130  |     |

4B

SEMIVOLATILE METHOD BLANK SUMMARY

Client ID

PB169286BL

Lab Name: Alliance

Contract: TETR06

Lab Code: ACE

SDG NO.: Q2881

Lab File ID: BN037619.D

Lab Sample ID: PB169286BL

Instrument ID: BNA\_N

Date Extracted: 08/18/2025

Matrix: (soil/water) Water

Date Analyzed: 08/20/2025

Level: (low/med) LOW

Time Analyzed: 04:50

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 |
|-------------------------|------------------|----------------|------------------|
| PB169286BS              | PB169286BS       | BN037630.D     | 08/20/2025       |
| MW-17B-55.5-08122025MS  | Q2842-04MS       | BN037605.D     | 08/19/2025       |
| MW-17B-55.5-08122025MSD | Q2842-05MSD      | BN037606.D     | 08/19/2025       |
| RW8-SP100-20250814      | Q2881-01         | BN037623.D     | 08/20/2025       |
| RW8-SP303-20250814      | Q2881-02         | BN037624.D     | 08/20/2025       |

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

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BN037576.D  
Instrument ID: BNA\_N

Contract: TETR06  
SDG NO.: Q2881  
DFTPP Injection Date: 08/12/2025  
DFTPP Injection Time: 15:05

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.0 ( 0.0 ) 1        |
| 69  | Mass 69 relative abundance         | 39.9                 |
| 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            | 7.1                  |
| 365 | Greater than 1% of mass 198        | 3.9                  |
| 441 | Present, but less than mass 443    | 14.2                 |
| 442 | Greater than 50% of mass 198       | 84.5                 |
| 443 | 15.0 - 24.0% of mass 442           | 16.2 ( 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 |
|-------------------|------------------|----------------|------------------|------------------|
| SSTDICC0.1        | SSTDICC0.1       | BN037578.D     | 08/12/2025       | 16:26            |
| SSTDICC0.2        | SSTDICC0.2       | BN037579.D     | 08/12/2025       | 17:03            |
| SSTDICCC0.4       | SSTDICCC0.4      | BN037580.D     | 08/12/2025       | 17:39            |
| SSTDICC0.8        | SSTDICC0.8       | BN037581.D     | 08/12/2025       | 18:16            |
| SSTDICC1.6        | SSTDICC1.6       | BN037582.D     | 08/12/2025       | 18:52            |
| SSTDICC3.2        | SSTDICC3.2       | BN037583.D     | 08/12/2025       | 19:29            |
| SSTDICC5.0        | SSTDICC5.0       | BN037584.D     | 08/12/2025       | 20:05            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BN037599.D  
Instrument ID: BNA\_N

Contract: TETR06  
SDG NO.: Q2881  
DFTPP Injection Date: 08/19/2025  
DFTPP Injection Time: 10:00

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.0 ( 0.0 ) 1        |
| 69  | Mass 69 relative abundance         | 35.9                 |
| 70  | Less than 2.0% of mass 69          | 0.2 ( 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            | 6.9                  |
| 365 | Greater than 1% of mass 198        | 4                    |
| 441 | Present, but less than mass 443    | 14.4                 |
| 442 | Greater than 50% of mass 198       | 97.7                 |
| 443 | 15.0 - 24.0% of mass 442           | 17.9 ( 18.3 ) 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 |
|-------------------------|------------------|----------------|------------------|------------------|
| SSTDCCC0.4              | SSTDCCC0.4       | BN037600.D     | 08/19/2025       | 10:39            |
| MW-17B-55.5-08122025MS  | Q2842-04MS       | BN037605.D     | 08/19/2025       | 13:40            |
| MW-17B-55.5-08122025MSD | Q2842-05MSD      | BN037606.D     | 08/19/2025       | 14:17            |
| SSTDCCC0.4EC            | SSTDCCC0.4       | BN037616.D     | 08/19/2025       | 20:19            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BN037617.D  
Instrument ID: BNA\_N

Contract: TETR06  
SDG NO.: Q2881  
DFTPP Injection Date: 08/20/2025  
DFTPP Injection Time: 03:35

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.4 ( 0.9 ) 1        |
| 69  | Mass 69 relative abundance         | 39.2                 |
| 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            | 6.8                  |
| 365 | Greater than 1% of mass 198        | 3.6                  |
| 441 | Present, but less than mass 443    | 11.8                 |
| 442 | Greater than 50% of mass 198       | 78.3                 |
| 443 | 15.0 - 24.0% of mass 442           | 14.9 ( 19 ) 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 |
|--------------------|------------------|----------------|------------------|------------------|
| SSTDCCC0.4         | SSTDCCC0.4       | BN037618.D     | 08/20/2025       | 04:14            |
| PB169286BL         | PB169286BL       | BN037619.D     | 08/20/2025       | 04:50            |
| RW8-SP100-20250814 | Q2881-01         | BN037623.D     | 08/20/2025       | 07:13            |
| RW8-SP303-20250814 | Q2881-02         | BN037624.D     | 08/20/2025       | 07:49            |
| PB169286BS         | PB169286BS       | BN037630.D     | 08/20/2025       | 11:25            |
| SSTDCCC0.4EC       | SSTDCCC0.4       | BN037631.D     | 08/20/2025       | 14:56            |

8B

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q2881

Client ID : SSTDCCC0.4

Date Analyzed: 08/19/2025

Lab File ID: BN037600.D

Time Analyzed: 10:39

Instrument ID: BNA\_N

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                | 2628                | 7.717 | 6457                | 10.50  | 3216                | 14.34  |
| UPPER LIMIT                | 5256                | 8.217 | 12914               | 10.998 | 6432                | 14.844 |
| LOWER LIMIT                | 1314                | 7.217 | 3228.5              | 9.998  | 1608                | 13.844 |
| EPA SAMPLE NO.             |                     |       |                     |        |                     |        |
| 01 MW-17B-55.5-08122025MS  | 2023                | 7.72  | 4765                | 10.50  | 2380                | 14.35  |
| 02 MW-17B-55.5-08122025MSD | 2022                | 7.72  | 4839                | 10.50  | 2406                | 14.34  |

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.: Q2881

Client ID: SSTDCCC0.4

Date Analyzed: 08/19/2025

Lab File ID: BN037600.D

Time Analyzed: 10:39

Instrument ID: BNA\_N

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                | 6293                | 17.086 | 5356                | 21.267 | 5379                | 23.504 |
| UPPER LIMIT                | 12586               | 17.586 | 10712               | 21.767 | 10758               | 24.004 |
| LOWER LIMIT                | 3146.5              | 16.586 | 2678                | 20.767 | 2689.5              | 23.004 |
| EPA SAMPLE NO.             |                     |        |                     |        |                     |        |
| 01 MW-17B-55.5-08122025MS  | 4922                | 17.09  | 4507                | 21.27  | 3909                | 23.51  |
| 02 MW-17B-55.5-08122025MSD | 5009                | 17.09  | 4561                | 21.27  | 3935                | 23.50  |

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.: Q2881

Client ID : SSTDCCC0.4

Date Analyzed: 08/20/2025

Lab File ID: BN037618.D

Time Analyzed: 04:14

Instrument ID: BNA\_N

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           | 2622                | 7.71 | 6166                | 10.50  | 3094                | 14.35  |
| UPPER LIMIT           | 5244                | 8.21 | 12332               | 10.998 | 6188                | 14.845 |
| LOWER LIMIT           | 1311                | 7.21 | 3083                | 9.998  | 1547                | 13.845 |
| EPA SAMPLE NO.        |                     |      |                     |        |                     |        |
| 01 RW8-SP100-20250814 | 2130                | 7.72 | 4994                | 10.50  | 2388                | 14.35  |
| 02 RW8-SP303-20250814 | 2203                | 7.71 | 5181                | 10.50  | 2518                | 14.35  |
| 03 PB169286BL         | 2323                | 7.71 | 5229                | 10.50  | 2344                | 14.35  |
| 04 PB169286BS         | 2389                | 7.71 | 5616                | 10.49  | 2629                | 14.35  |

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.: Q2881

Client ID: SSTDCCC0.4

Date Analyzed: 08/20/2025

Lab File ID: BN037618.D

Time Analyzed: 04:14

Instrument ID: BNA\_N

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        | 6113                | 17.086 | 5494                | 21.268 | 5322                | 23.499 |
|    | UPPER LIMIT        | 12226               | 17.586 | 10988               | 21.768 | 10644               | 23.999 |
|    | LOWER LIMIT        | 3056.5              | 16.586 | 2747                | 20.768 | 2661                | 22.999 |
|    | EPA SAMPLE NO.     |                     |        |                     |        |                     |        |
| 01 | RW8-SP100-20250814 | 5113                | 17.09  | 4521                | 21.27  | 4086                | 23.50  |
| 02 | RW8-SP303-20250814 | 5410                | 17.09  | 4670                | 21.27  | 4150                | 23.50  |
| 03 | PB169286BL         | 4466                | 17.09  | 3733                | 21.27  | 3552                | 23.50  |
| 04 | PB169286BS         | 5040                | 17.09  | 4175                | 21.27  | 3780                | 23.50  |

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:            | Tetra Tech NUS, Inc.          |                  | Date Collected: |                |
| Project:           | NWIRP Bethpage 112G08005-WE13 |                  | Date Received:  |                |
| Client Sample ID:  | PB169286BL                    |                  | SDG No.:        | Q2881          |
| Lab Sample ID:     | PB169286BL                    |                  | Matrix:         | Water          |
| Analytical Method: | SW8270ESIM                    |                  | % Solid:        | 0              |
| Sample Wt/Vol:     | 1000                          | Units: mL        | Final Vol:      | 1000 uL        |
| Soil Aliquot Vol:  |                               | uL               | Test:           | SVOC-SIMGroup1 |
| 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 |
| BN037619.D        | 1         | 08/18/25 08:41 | 08/20/25 04:50 | PB169286      |

| CAS Number                | Parameter               | Conc. | Qualifier | MDL      | LOD  | LOQ / CRQL | Units    |
|---------------------------|-------------------------|-------|-----------|----------|------|------------|----------|
| <b>TARGETS</b>            |                         |       |           |          |      |            |          |
| 123-91-1                  | 1,4-Dioxane             | 0.20  | U         | 0.070    | 0.20 | 0.20       | ug/L     |
| <b>SURROGATES</b>         |                         |       |           |          |      |            |          |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.32  |           | 30 - 150 |      | 81%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.33  |           | 30 - 150 |      | 83%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.36  |           | 55 - 111 |      | 90%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.35  |           | 53 - 106 |      | 86%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.37  |           | 58 - 132 |      | 93%        | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 2320  | 7.71      |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 5230  | 10.498    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 2340  | 14.345    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 4470  | 17.087    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 3730  | 21.268    |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 3550  | 23.499    |          |      |            |          |

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:            | Tetra Tech NUS, Inc.          |                  | Date Collected: |                |
| Project:           | NWIRP Bethpage 112G08005-WE13 |                  | Date Received:  |                |
| Client Sample ID:  | PB169286BS                    |                  | SDG No.:        | Q2881          |
| Lab Sample ID:     | PB169286BS                    |                  | Matrix:         | Water          |
| Analytical Method: | SW8270ESIM                    |                  | % Solid:        | 0              |
| Sample Wt/Vol:     | 1000                          | Units: mL        | Final Vol:      | 1000 uL        |
| Soil Aliquot Vol:  |                               | uL               | Test:           | SVOC-SIMGroup1 |
| 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 |
| BN037630.D        | 1         | 08/18/25 08:41 | 08/20/25 11:25 | PB169286      |

| CAS Number                | Parameter               | Conc. | Qualifier | MDL      | LOD  | LOQ / CRQL | Units    |
|---------------------------|-------------------------|-------|-----------|----------|------|------------|----------|
| <b>TARGETS</b>            |                         |       |           |          |      |            |          |
| 123-91-1                  | 1,4-Dioxane             | 0.30  |           | 0.070    | 0.20 | 0.20       | ug/L     |
| <b>SURROGATES</b>         |                         |       |           |          |      |            |          |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.33  |           | 30 - 150 |      | 82%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.31  |           | 30 - 150 |      | 76%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.35  |           | 55 - 111 |      | 87%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.36  |           | 53 - 106 |      | 90%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.35  |           | 58 - 132 |      | 88%        | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 2390  | 7.71      |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 5620  | 10.487    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 2630  | 14.345    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 5040  | 17.086    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 4180  | 21.268    |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 3780  | 23.502    |          |      |            |          |

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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:            | Tetra Tech NUS, Inc.          | Date Collected: | 08/12/25       |
| Project:           | NWIRP Bethpage 112G08005-WE13 | Date Received:  | 08/12/25       |
| Client Sample ID:  | MW-17B-55.5-08122025MS        | SDG No.:        | Q2881          |
| Lab Sample ID:     | Q2842-04MS                    | Matrix:         | Water          |
| Analytical Method: | SW8270ESIM                    | % Solid:        | 0              |
| Sample Wt/Vol:     | 940 Units: mL                 | Final Vol:      | 1000 uL        |
| Soil Aliquot Vol:  | uL                            | Test:           | SVOC-SIMGroup1 |
| 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 |
| BN037605.D        | 1         | 08/18/25 08:41 | 08/19/25 13:40 | PB169286      |

| CAS Number                | Parameter               | Conc. | Qualifier | MDL      | LOD  | LOQ / CRQL | Units    |
|---------------------------|-------------------------|-------|-----------|----------|------|------------|----------|
| <b>TARGETS</b>            |                         |       |           |          |      |            |          |
| 123-91-1                  | 1,4-Dioxane             | 3.00  |           | 0.070    | 0.21 | 0.21       | ug/L     |
| <b>SURROGATES</b>         |                         |       |           |          |      |            |          |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.31  |           | 30 - 150 |      | 77%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.38  |           | 30 - 150 |      | 94%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.33  |           | 55 - 111 |      | 83%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.32  |           | 53 - 106 |      | 81%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.47  |           | 58 - 132 |      | 117%       | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 2020  | 7.717     |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 4770  | 10.498    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 2380  | 14.345    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 4920  | 17.087    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 4510  | 21.268    |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 3910  | 23.505    |          |      |            |          |

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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:            | Tetra Tech NUS, Inc.          | Date Collected: | 08/12/25       |
| Project:           | NWIRP Bethpage 112G08005-WE13 | Date Received:  | 08/12/25       |
| Client Sample ID:  | MW-17B-55.5-08122025MSD       | SDG No.:        | Q2881          |
| Lab Sample ID:     | Q2842-05MSD                   | Matrix:         | Water          |
| Analytical Method: | SW8270ESIM                    | % Solid:        | 0              |
| Sample Wt/Vol:     | 960 Units: mL                 | Final Vol:      | 1000 uL        |
| Soil Aliquot Vol:  | uL                            | Test:           | SVOC-SIMGroup1 |
| 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 |
| BN037606.D        | 1         | 08/18/25 08:41 | 08/19/25 14:17 | PB169286      |

| CAS Number                | Parameter               | Conc. | Qualifier | MDL      | LOD  | LOQ / CRQL | Units    |
|---------------------------|-------------------------|-------|-----------|----------|------|------------|----------|
| <b>TARGETS</b>            |                         |       |           |          |      |            |          |
| 123-91-1                  | 1,4-Dioxane             | 3.00  |           | 0.070    | 0.21 | 0.21       | ug/L     |
| <b>SURROGATES</b>         |                         |       |           |          |      |            |          |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.30  |           | 30 - 150 |      | 76%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.37  |           | 30 - 150 |      | 93%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.32  |           | 55 - 111 |      | 79%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.32  |           | 53 - 106 |      | 81%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.47  |           | 58 - 132 |      | 117%       | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 2020  | 7.717     |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 4840  | 10.498    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 2410  | 14.344    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 5010  | 17.086    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 4560  | 21.267    |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 3940  | 23.504    |          |      |            |          |

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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\_N\Methods\  
 Method File : 8270-SIM-BN081225.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 Last Update : Wed Aug 13 05:00:58 2025  
 Response Via : Initial Calibration

## Calibration Files

0.1 =BN037578.D 0.2 =BN037579.D 0.4 =BN037580.D 0.8 =BN037581.D 1.6 =BN037582.D 3.2 =BN037583.D 5 =BN037584.D

|           | Compound              | 0.1            | 0.2   | 0.4   | 0.8   | 1.6   | 3.2   | 5     | Avg   | %RSD  |
|-----------|-----------------------|----------------|-------|-------|-------|-------|-------|-------|-------|-------|
| -----     |                       |                |       |       |       |       |       |       |       |       |
| 1) I      | 1,4-Dichlorobenzen... | -----ISTD----- |       |       |       |       |       |       |       |       |
| 2)        | 1,4-Dioxane           | 0.413          | 0.395 | 0.368 | 0.397 | 0.375 | 0.348 | 0.383 |       | 6.08  |
| 3)        | n-Nitrosodimet...     | 0.471          | 0.488 | 0.471 | 0.493 | 0.503 | 0.508 | 0.489 |       | 3.19  |
| 4) S      | 2-Fluorophenol        | 0.922          | 0.945 | 0.888 | 0.823 | 0.886 | 0.906 | 0.977 | 0.907 | 5.41  |
| 5) S      | Phenol-d6             | 1.045          | 1.051 | 1.044 | 0.983 | 1.198 | 1.117 | 1.201 | 1.091 | 7.66  |
| 6)        | bis(2-Chloroet...     | 0.935          | 0.987 | 0.976 | 0.936 | 1.020 | 1.008 | 1.022 | 0.983 | 3.75  |
|           |                       |                |       |       |       |       |       |       |       |       |
| 7) I      | Naphthalene-d8        | -----ISTD----- |       |       |       |       |       |       |       |       |
| 8) S      | Nitrobenzene-d5       | 0.270          | 0.257 | 0.262 | 0.264 | 0.292 | 0.301 | 0.326 | 0.282 | 9.03  |
| 9)        | Naphthalene           | 1.044          | 1.050 | 1.037 | 1.012 | 1.089 | 1.096 | 1.127 | 1.065 | 3.78  |
| 10)       | Hexachlorobuta...     | 0.252          | 0.258 | 0.262 | 0.253 | 0.267 | 0.264 | 0.265 | 0.260 | 2.22  |
| 11) SURR2 | Methylnaphth...       | 0.488          | 0.492 | 0.508 | 0.493 | 0.543 | 0.578 | 0.705 | 0.544 | 14.39 |
| 12)       | 2-Methylnaphth...     | 0.589          | 0.631 | 0.635 | 0.629 | 0.701 | 0.731 | 0.760 | 0.668 | 9.39  |
|           |                       |                |       |       |       |       |       |       |       |       |
| 13) I     | Acenaphthene-d10      | -----ISTD----- |       |       |       |       |       |       |       |       |
| 14) S     | 2,4,6-Tribromo...     | 0.139          | 0.149 | 0.156 | 0.160 | 0.185 | 0.203 | 0.234 | 0.175 | 19.39 |
| 15) S     | 2-Fluorobiphenyl      | 2.226          | 2.243 | 2.300 | 2.251 | 2.341 | 2.391 | 2.440 | 2.313 | 3.50  |
| 16)       | Acenaphthylene        | 1.740          | 1.760 | 1.710 | 1.653 | 1.850 | 1.858 | 1.980 | 1.793 | 6.15  |
| 17)       | Acenaphthene          | 1.144          | 1.150 | 1.172 | 1.154 | 1.283 | 1.286 | 1.348 | 1.220 | 6.86  |
| 18)       | Fluorene              | 1.466          | 1.510 | 1.524 | 1.521 | 1.686 | 1.693 | 1.770 | 1.596 | 7.38  |
|           |                       |                |       |       |       |       |       |       |       |       |
| 19) I     | Phenanthrene-d10      | -----ISTD----- |       |       |       |       |       |       |       |       |
| 20)       | 4,6-Dinitro-2-...     | 0.042          | 0.046 | 0.047 | 0.061 | 0.070 |       | 0.053 |       | 22.37 |
| 21)       | 4-Bromophenyl-...     | 0.226          | 0.236 | 0.234 | 0.237 | 0.265 | 0.271 | 0.292 | 0.251 | 9.77  |
| 22)       | Hexachlorobenzene     | 0.360          | 0.360 | 0.356 | 0.332 | 0.364 | 0.354 | 0.370 | 0.356 | 3.37  |
| 23)       | Atrazine              | 0.127          | 0.126 | 0.130 | 0.134 | 0.160 | 0.176 |       | 0.142 | 14.76 |
| 24)       | Pentachlorophenol     |                | 0.108 | 0.113 | 0.113 | 0.139 | 0.152 |       | 0.125 | 15.36 |
| 25)       | Phenanthrene          | 1.146          | 1.162 | 1.170 | 1.138 | 1.281 | 1.269 | 1.347 | 1.216 | 6.72  |
| 26)       | Anthracene            | 0.950          | 0.985 | 0.995 | 0.988 | 1.145 | 1.186 | 1.286 | 1.077 | 11.93 |
| 27) SURR  | Fluoranthene-d10      | 0.932          | 0.966 | 0.954 | 0.935 | 1.050 | 1.110 | 1.419 | 1.052 | 16.61 |
| 28)       | Fluoranthene          | 1.241          | 1.264 | 1.299 | 1.291 | 1.503 | 1.533 | 1.648 | 1.397 | 11.53 |
|           |                       |                |       |       |       |       |       |       |       |       |
| 29) I     | Chrysene-d12          | -----ISTD----- |       |       |       |       |       |       |       |       |
| 30)       | Pyrene                | 1.539          | 1.486 | 1.464 | 1.403 | 1.502 | 1.520 | 1.552 | 1.495 | 3.38  |
| 31) S     | Terphenyl-d14         | 0.814          | 0.801 | 0.796 | 0.763 | 0.833 | 0.853 | 0.901 | 0.823 | 5.45  |
| 32)       | Benzo(a)anthra...     | 1.266          | 1.275 | 1.263 | 1.253 | 1.351 | 1.458 | 1.487 | 1.336 | 7.40  |
| 33)       | Chrysene              | 1.547          | 1.515 | 1.433 | 1.379 | 1.511 | 1.489 | 1.551 | 1.489 | 4.21  |
| 34)       | Bis(2-ethylhex...     |                | 0.518 | 0.522 | 0.483 | 0.522 | 0.593 | 0.671 | 0.551 | 12.47 |
|           |                       |                |       |       |       |       |       |       |       |       |
| 35) I     | Perylene-d12          | -----ISTD----- |       |       |       |       |       |       |       |       |

Method Path : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\

Method File : 8270-SIM-BN081225.M

|       |                   |       |       |       |       |       |       |       |       |       |
|-------|-------------------|-------|-------|-------|-------|-------|-------|-------|-------|-------|
| 36)   | Indeno(1,2,3-c... | 1.446 | 1.462 | 1.612 | 1.568 | 1.816 | 1.883 | 2.011 | 1.686 | 13.01 |
| 37)   | Benzo(b)fluora... | 1.351 | 1.375 | 1.338 | 1.432 | 1.615 | 1.674 | 1.825 | 1.516 | 12.52 |
| 38)   | Benzo(k)fluora... | 1.597 | 1.580 | 1.651 | 1.628 | 1.821 | 1.797 | 1.898 | 1.710 | 7.36  |
| 39) C | Benzo(a)pyrene    | 1.146 | 1.136 | 1.149 | 1.157 | 1.322 | 1.382 | 1.508 | 1.257 | 11.77 |
| 40)   | Dibenzo(a,h)an... | 1.023 | 1.118 | 1.222 | 1.215 | 1.439 | 1.516 | 1.621 | 1.308 | 16.85 |
| 41)   | Benzo(g,h,i)pe... | 1.207 | 1.274 | 1.262 | 1.261 | 1.467 | 1.532 | 1.643 | 1.378 | 12.17 |

-----  
(#) = Out of Range

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: Alliance Contract: TETR06  
Lab Code: ACE SDG No.: Q2881  
Instrument ID: BNA\_N Calibration Date/Time: 08/19/2025 10:39  
Lab File ID: BN037600.D Init. Calib. Date(s): 08/12/2025 08/12/2025  
EPA Sample No.: SSTDCCC0.4 Init. Calib. Time(s): 16:26 20:05  
GC Column: ZB-GR ID: 0.25 (mm)

| COMPOUND                | RRF   | RRF0.4 | MIN<br>RRF | %D   | MAX%D |
|-------------------------|-------|--------|------------|------|-------|
| 2-Methylnaphthalene-d10 | 0.544 | 0.498  |            | -8.5 | 20.0  |
| Fluoranthene-d10        | 1.052 | 0.957  |            | -9.0 | 20.0  |
| 2-Fluorophenol          | 0.907 | 0.907  |            | 0.0  | 20.0  |
| Phenol-d6               | 1.091 | 1.078  |            | -1.2 | 20.0  |
| Nitrobenzene-d5         | 0.282 | 0.275  |            | -2.5 | 20.0  |
| 2-Fluorobiphenyl        | 2.313 | 2.186  |            | -5.5 | 20.0  |
| 2,4,6-Tribromophenol    | 0.175 | 0.169  |            | -3.4 | 20.0  |
| Terphenyl-d14           | 0.823 | 0.796  |            | -3.3 | 20.0  |
| 1,4-Dioxane             | 0.383 | 0.406  |            | 6.0  | 20.0  |

All other compounds must meet a minimum RRF of 0.010.

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: Alliance Contract: TETR06  
Lab Code: ACE SDG No.: Q2881  
Instrument ID: BNA\_N Calibration Date/Time: 08/19/2025 20:19  
Lab File ID: BN037616.D Init. Calib. Date(s): 08/12/2025 08/12/2025  
EPA Sample No.: SSTDCCC0.4EC Init. Calib. Time(s): 16:26 20:05  
GC Column: ZB-GR ID: 0.25 (mm)

| COMPOUND                | RRF   | RRF0.4 | MIN<br>RRF | %D   | MAX%D |
|-------------------------|-------|--------|------------|------|-------|
| 2-Methylnaphthalene-d10 | 0.544 | 0.508  |            | -6.6 | 50.0  |
| Fluoranthene-d10        | 1.052 | 0.969  |            | -7.9 | 50.0  |
| 2-Fluorophenol          | 0.907 | 0.927  |            | 2.2  | 50.0  |
| Phenol-d6               | 1.091 | 1.101  |            | 0.9  | 50.0  |
| Nitrobenzene-d5         | 0.282 | 0.270  |            | -4.3 | 50.0  |
| 2-Fluorobiphenyl        | 2.313 | 2.275  |            | -1.6 | 50.0  |
| 2,4,6-Tribromophenol    | 0.175 | 0.174  |            | -0.6 | 50.0  |
| Terphenyl-d14           | 0.823 | 0.818  |            | -0.6 | 50.0  |
| 1,4-Dioxane             | 0.383 | 0.398  |            | 3.9  | 50.0  |

All other compounds must meet a minimum RRF of 0.010.

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: Alliance Contract: TETR06  
Lab Code: ACE SDG No.: Q2881  
Instrument ID: BNA\_N Calibration Date/Time: 08/20/2025 04:14  
Lab File ID: BN037618.D Init. Calib. Date(s): 08/12/2025 08/12/2025  
EPA Sample No.: SSTDCCC0.4 Init. Calib. Time(s): 16:26 20:05  
GC Column: ZB-GR ID: 0.25 (mm)

| COMPOUND                | RRF   | RRF0.4 | MIN<br>RRF | %D   | MAX%D |
|-------------------------|-------|--------|------------|------|-------|
| 2-Methylnaphthalene-d10 | 0.544 | 0.498  |            | -8.5 | 20.0  |
| Fluoranthene-d10        | 1.052 | 0.951  |            | -9.6 | 20.0  |
| 2-Fluorophenol          | 0.907 | 0.875  |            | -3.5 | 20.0  |
| Phenol-d6               | 1.091 | 1.045  |            | -4.2 | 20.0  |
| Nitrobenzene-d5         | 0.282 | 0.281  |            | -0.4 | 20.0  |
| 2-Fluorobiphenyl        | 2.313 | 2.246  |            | -2.9 | 20.0  |
| 2,4,6-Tribromophenol    | 0.175 | 0.175  |            | 0.0  | 20.0  |
| Terphenyl-d14           | 0.823 | 0.793  |            | -3.6 | 20.0  |
| 1,4-Dioxane             | 0.383 | 0.380  |            | -0.8 | 20.0  |

All other compounds must meet a minimum RRF of 0.010.

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

|                 |                     |                        |                              |
|-----------------|---------------------|------------------------|------------------------------|
| Lab Name:       | <u>Alliance</u>     | Contract:              | <u>TETR06</u>                |
| Lab Code:       | <u>ACE</u>          | SDG No.:               | <u>Q2881</u>                 |
| Instrument ID:  | <u>BNA_N</u>        | Calibration Date/Time: | <u>08/20/2025 14:56</u>      |
| Lab File ID:    | <u>BN037631.D</u>   | Init. Calib. Date(s):  | <u>08/12/2025 08/12/2025</u> |
| EPA Sample No.: | <u>SSTDCCC0.4EC</u> | Init. Calib. Time(s):  | <u>16:26 20:05</u>           |
| GC Column:      | <u>ZB-GR</u>        | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                | RRF   | RRF0.4 | MIN<br>RRF | %D    | MAX%D |
|-------------------------|-------|--------|------------|-------|-------|
| 2-Methylnaphthalene-d10 | 0.544 | 0.491  |            | -9.7  | 50.0  |
| Fluoranthene-d10        | 1.052 | 0.944  |            | -10.3 | 50.0  |
| 2-Fluorophenol          | 0.907 | 0.853  |            | -6.0  | 50.0  |
| Phenol-d6               | 1.091 | 0.982  |            | -10.0 | 50.0  |
| Nitrobenzene-d5         | 0.282 | 0.275  |            | -2.5  | 50.0  |
| 2-Fluorobiphenyl        | 2.313 | 2.157  |            | -6.7  | 50.0  |
| 2,4,6-Tribromophenol    | 0.175 | 0.172  |            | -1.7  | 50.0  |
| Terphenyl-d14           | 0.823 | 0.717  |            | -12.9 | 50.0  |
| 1,4-Dioxane             | 0.383 | 0.377  |            | -1.6  | 50.0  |

All other compounds must meet a minimum RRF of 0.010.