



LAB CHRONICLE

|          |                      |            |                               |
|----------|----------------------|------------|-------------------------------|
| OrderID: | Q3086                | OrderDate: | 9/12/2025 10:59:00 AM         |
| Client:  | Tetra Tech NUS, Inc. | Project:   | NWIRP Bethpage 112G08005-WE13 |
| Contact: | Ernie Wu             | Location:  | D31                           |

| LabID    | ClientID               | Matrix | Test           | Method        | Sample Date | Prep Date | Anal Date | Received |
|----------|------------------------|--------|----------------|---------------|-------------|-----------|-----------|----------|
| Q3086-01 | RW8-SP100-2025091<br>1 | Water  | SVOC-SIMGroup1 | 8270-Modified | 09/11/25    | 09/15/25  | 09/17/25  | 09/12/25 |
| Q3086-02 | RW8-SP303-2025091<br>1 | Water  | SVOC-SIMGroup1 | 8270-Modified | 09/11/25    | 09/15/25  | 09/17/25  | 09/12/25 |



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

### Hit Summary Sheet SW-846

**SDG No.:** Q3086  
**Client:** Tetra Tech NUS, Inc.

| Sample ID            | Client ID          | Parameter | Concentration | C     | MDL | LOD  | RDL | Units    |
|----------------------|--------------------|-----------|---------------|-------|-----|------|-----|----------|
| Client ID :          | RW8-SP100-20250911 |           |               |       |     |      |     |          |
| Q3086-01             | RW8-SP100-20250911 | 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: | 09/11/25       |
| Project:           | NWIRP Bethpage 112G08005-WE13 | Date Received:  | 09/12/25       |
| Client Sample ID:  | RW8-SP100-20250911            | SDG No.:        | Q3086          |
| Lab Sample ID:     | Q3086-01                      | 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 |
| BN037745.D        | 1         | 09/15/25 10:15 | 09/17/25 14:35 | PB169692      |

| 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 | 0.36  |           | 30 - 150 |      | 89%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.39  |           | 30 - 150 |      | 97%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.39  |           | 55 - 111 |      | 97%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.40  |           | 53 - 106 |      | 101%       | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.55  | *         | 58 - 132 |      | 138%       | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 6570  | 7.84      |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 18000 | 10.637    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 8330  | 14.495    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 15400 | 17.235    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 14200 | 21.411    |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 14200 | 23.75     |          |      |            |          |

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: | 09/11/25       |
| Project:           | NWIRP Bethpage 112G08005-WE13 | Date Received:  | 09/12/25       |
| Client Sample ID:  | RW8-SP303-20250911            | SDG No.:        | Q3086          |
| Lab Sample ID:     | Q3086-02                      | 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 |
| BN037746.D        | 1         | 09/15/25 10:15 | 09/17/25 15:11 | PB169692      |

| 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 | 0.36  |           | 30 - 150 |      | 91%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.35  |           | 30 - 150 |      | 86%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.39  |           | 55 - 111 |      | 97%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.41  |           | 53 - 106 |      | 103%       | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.52  |           | 58 - 132 |      | 130%       | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 6530  | 7.84      |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 17800 | 10.637    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 8070  | 14.495    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 13900 | 17.235    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 10600 | 21.411    |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 10300 | 23.75     |          |      |            |          |

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.: **Q3086**

Client: **Tetra Tech NUS, Inc.**

Analytical Method: **8270-Modified**

| Lab Sample ID | Client ID           | Parameter               | Spike (PPM) | Result (PPM) | Recovery (%) | Qual | Limits (%) |      |
|---------------|---------------------|-------------------------|-------------|--------------|--------------|------|------------|------|
|               |                     |                         |             |              |              |      | Low        | High |
| PB169692BL    | PB169692BL          | 2-Methylnaphthalene-d10 | 0.4         | 0.31         | 79           |      | 30         | 150  |
|               |                     | Fluoranthene-d10        | 0.4         | 0.30         | 75           |      | 30         | 150  |
|               |                     | Nitrobenzene-d5         | 0.4         | 0.39         | 97           |      | 55         | 111  |
|               |                     | 2-Fluorobiphenyl        | 0.4         | 0.37         | 92           |      | 53         | 106  |
|               |                     | Terphenyl-d14           | 0.4         | 0.34         | 84           |      | 58         | 132  |
| PB169692BS    | PB169692BS          | 2-Methylnaphthalene-d10 | 0.4         | 0.30         | 76           |      | 30         | 150  |
|               |                     | Fluoranthene-d10        | 0.4         | 0.30         | 75           |      | 30         | 150  |
|               |                     | Nitrobenzene-d5         | 0.4         | 0.35         | 88           |      | 55         | 111  |
|               |                     | 2-Fluorobiphenyl        | 0.4         | 0.36         | 90           |      | 53         | 106  |
|               |                     | Terphenyl-d14           | 0.4         | 0.31         | 78           |      | 58         | 132  |
| Q3086-01      | RW8-SP100-20250911  | 2-Methylnaphthalene-d10 | 0.4         | 0.36         | 89           |      | 30         | 150  |
|               |                     | Fluoranthene-d10        | 0.4         | 0.39         | 97           |      | 30         | 150  |
|               |                     | Nitrobenzene-d5         | 0.4         | 0.39         | 97           |      | 55         | 111  |
|               |                     | 2-Fluorobiphenyl        | 0.4         | 0.40         | 101          |      | 53         | 106  |
|               |                     | Terphenyl-d14           | 0.4         | 0.55         | 138          | *    | 58         | 132  |
| Q3086-02      | RW8-SP303-20250911  | 2-Methylnaphthalene-d10 | 0.4         | 0.36         | 91           |      | 30         | 150  |
|               |                     | Fluoranthene-d10        | 0.4         | 0.35         | 86           |      | 30         | 150  |
|               |                     | Nitrobenzene-d5         | 0.4         | 0.39         | 97           |      | 55         | 111  |
|               |                     | 2-Fluorobiphenyl        | 0.4         | 0.41         | 103          |      | 53         | 106  |
|               |                     | Terphenyl-d14           | 0.4         | 0.52         | 130          |      | 58         | 132  |
| Q3095-06MS    | MW203D2-20250909MS  | 2-Methylnaphthalene-d10 | 0.4         | 0.29         | 72           |      | 30         | 150  |
|               |                     | Fluoranthene-d10        | 0.4         | 0.33         | 83           |      | 30         | 150  |
|               |                     | Nitrobenzene-d5         | 0.4         | 0.30         | 75           |      | 55         | 111  |
|               |                     | 2-Fluorobiphenyl        | 0.4         | 0.32         | 79           |      | 53         | 106  |
|               |                     | Terphenyl-d14           | 0.4         | 0.40         | 99           |      | 58         | 132  |
| Q3095-07MSD   | MW203D2-20250909MSD | 2-Methylnaphthalene-d10 | 0.4         | 0.30         | 74           |      | 30         | 150  |
|               |                     | Fluoranthene-d10        | 0.4         | 0.33         | 83           |      | 30         | 150  |
|               |                     | Nitrobenzene-d5         | 0.4         | 0.30         | 75           |      | 55         | 111  |
|               |                     | 2-Fluorobiphenyl        | 0.4         | 0.32         | 79           |      | 53         | 106  |
|               |                     | Terphenyl-d14           | 0.4         | 0.40         | 99           |      | 58         | 132  |



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Fax : 908 789 8922

Matrix Spike/Matrix Spike Duplicate Summary  
SW-846

SDG No.:

Q3086

Analytical Method:

SW8270-Modified

Client:

Tetra Tech NUS, Inc.

DataFile:

BN037743.D

| Parameter      | Spike      | Sample Result     | Result             | Units | Rec | Rec Qual | RPD | RPD Qual | Low | Limits High | RPD |
|----------------|------------|-------------------|--------------------|-------|-----|----------|-----|----------|-----|-------------|-----|
| Lab Sample ID: | Q3095-06MS | Client Sample ID: | MW203D2-20250909MS |       |     |          |     |          |     |             |     |
| 1,4-Dioxane    | 0.4        | 0.85              | 0.97               | ug/L  | 30  | *        |     |          | 70  | 130         |     |

Matrix Spike/Matrix Spike Duplicate Summary  
SW-846

|          |                      |                    |                 |
|----------|----------------------|--------------------|-----------------|
| SDG No.: | Q3086                | Analytical Method: | SW8270-Modified |
| Client:  | Tetra Tech NUS, Inc. | DataFile:          | BN037744.D      |

| Parameter      | Spike       | Sample Result     | Result              | Units | Rec | Rec Qual | RPD | RPD Qual | Low | Limits High | RPD |
|----------------|-------------|-------------------|---------------------|-------|-----|----------|-----|----------|-----|-------------|-----|
| Lab Sample ID: | Q3095-07MSD | Client Sample ID: | MW203D2-20250909MSD |       |     |          |     |          |     |             |     |
| 1,4-Dioxane    | 0.4         | 0.85              | 0.96                | ug/L  | 27  | *        | 11  |          | 70  | 130         | 20  |

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q3086

Analytical Method: 8270-Modified

Client: Tetra Tech NUS, Inc.

DataFile: BN037791.D

| Lab Sample ID | Parameter   | Spike | Result | Unit | Rec | RPD | Qual | RPD  | Limits |      | RPD |
|---------------|-------------|-------|--------|------|-----|-----|------|------|--------|------|-----|
|               |             |       |        |      |     |     |      | Qual | Low    | High |     |
| PB169692BS    | 1,4-Dioxane | 0.4   | 0.30   | ug/L | 75  |     |      |      | 70     | 130  |     |

4B

SEMIVOLATILE METHOD BLANK SUMMARY

Client ID

PB169692BL

Lab Name: Alliance

Contract: TETR06

Lab Code: ACE

SDG NO.: Q3086

Lab File ID: BN037796.D

Lab Sample ID: PB169692BL

Instrument ID: BNA\_N

Date Extracted: 09/15/2025

Matrix: (soil/water) Water

Date Analyzed: 09/19/2025

Level: (low/med) LOW

Time Analyzed: 03:24

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 |
|---------------------|------------------|----------------|------------------|
| RW8-SP100-20250911  | Q3086-01         | BN037745.D     | 09/17/2025       |
| RW8-SP303-20250911  | Q3086-02         | BN037746.D     | 09/17/2025       |
| MW203D2-20250909MS  | Q3095-06MS       | BN037743.D     | 09/17/2025       |
| MW203D2-20250909MSD | Q3095-07MSD      | BN037744.D     | 09/17/2025       |
| PB169692BS          | PB169692BS       | BN037791.D     | 09/18/2025       |

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

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BN037689.D  
Instrument ID: BNA\_N

Contract: TETR06  
SDG NO.: Q3086  
DFTPP Injection Date: 09/11/2025  
DFTPP Injection Time: 07:54

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.4 ( 1 ) 1          |
| 69  | Mass 69 relative abundance         | 100                  |
| 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            | 6.2                  |
| 365 | Greater than 1% of mass 198        | 3.5                  |
| 441 | Present, but less than mass 443    | 79.2                 |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 14.5 (20.4) 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       | BN037691.D     | 09/11/2025       | 09:56            |
| SSTDICC0.2        | SSTDICC0.2       | BN037692.D     | 09/11/2025       | 10:32            |
| SSTDICCC0.4       | SSTDICCC0.4      | BN037693.D     | 09/11/2025       | 11:08            |
| SSTDICC0.8        | SSTDICC0.8       | BN037694.D     | 09/11/2025       | 11:44            |
| SSTDICC1.6        | SSTDICC1.6       | BN037695.D     | 09/11/2025       | 12:20            |
| SSTDICC3.2        | SSTDICC3.2       | BN037696.D     | 09/11/2025       | 12:56            |
| SSTDICC5.0        | SSTDICC5.0       | BN037697.D     | 09/11/2025       | 13:33            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BN037739.D  
Instrument ID: BNA\_N

Contract: TETR06  
SDG NO.: Q3086  
DFTPP Injection Date: 09/17/2025  
DFTPP Injection Time: 10:46

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.5 ( 1.3 ) 1        |
| 69  | Mass 69 relative abundance         | 100                  |
| 70  | Less than 2.0% of mass 69          | 0.2 ( 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            | 6.9                  |
| 365 | Greater than 1% of mass 198        | 3.7                  |
| 441 | Present, but less than mass 443    | 82.6                 |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 17.7 (21.3) 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       | BN037740.D     | 09/17/2025       | 11:25            |
| MW203D2-20250909MS  | Q3095-06MS       | BN037743.D     | 09/17/2025       | 13:14            |
| MW203D2-20250909MSD | Q3095-07MSD      | BN037744.D     | 09/17/2025       | 13:59            |
| RW8-SP100-20250911  | Q3086-01         | BN037745.D     | 09/17/2025       | 14:35            |
| RW8-SP303-20250911  | Q3086-02         | BN037746.D     | 09/17/2025       | 15:11            |
| SSTDCCC0.4EC        | SSTDCCC0.4       | BN037757.D     | 09/17/2025       | 21:50            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BN037775.D  
Instrument ID: BNA\_N

Contract: TETR06  
SDG NO.: Q3086  
DFTPP Injection Date: 09/18/2025  
DFTPP Injection Time: 13:53

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.7 ( 1.7 ) 1        |
| 69  | Mass 69 relative abundance         | 100                  |
| 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            | 7                    |
| 365 | Greater than 1% of mass 198        | 3.4                  |
| 441 | Present, but less than mass 443    | 85.6                 |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 16.2 (18.9) 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 |
|-------------------|------------------|----------------|------------------|------------------|
| SSTDCCC0.4        | SSTDCCC0.4       | BN037776.D     | 09/18/2025       | 14:32            |
| PB169692BS        | PB169692BS       | BN037791.D     | 09/18/2025       | 23:40            |
| SSTDCCC0.4EC      | SSTDCCC0.4       | BN037793.D     | 09/19/2025       | 00:52            |

5B

SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: Alliance  
Lab Code: ACE  
Lab File ID: BN037794.D  
Instrument ID: BNA\_N

Contract: TETR06  
SDG NO.: Q3086  
DFTPP Injection Date: 09/19/2025  
DFTPP Injection Time: 02:08

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 68  | Less than 2.0% of mass 69          | 0.5 ( 1.4 ) 1        |
| 69  | Mass 69 relative abundance         | 100                  |
| 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        | 3.8                  |
| 441 | Present, but less than mass 443    | 86.4                 |
| 442 | Greater than 50% of mass 198       | 100                  |
| 443 | 15.0 - 24.0% of mass 442           | 16.9 (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       | BN037795.D     | 09/19/2025       | 02:47            |
| PB169692BL        | PB169692BL       | BN037796.D     | 09/19/2025       | 03:24            |
| SSTDCCC0.4EC      | SSTDCCC0.4       | BN037811.D     | 09/19/2025       | 13:03            |

8B

SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance

Lab Code: ACE

SDG NO.: Q3086

Client ID : SSTDCCC0.4

Date Analyzed: 09/17/2025

Lab File ID: BN037740.D

Time Analyzed: 11:25

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            | 6660                | 7.84 | 18396               | 10.65  | 8930                | 14.50  |
| UPPER LIMIT            | 13320               | 8.34 | 36792               | 11.148 | 17860               | 14.995 |
| LOWER LIMIT            | 3330                | 7.34 | 9198                | 10.148 | 4465                | 13.995 |
| EPA SAMPLE NO.         |                     |      |                     |        |                     |        |
| 01 RW8-SP100-20250911  | 6573                | 7.84 | 18040               | 10.64  | 8326                | 14.50  |
| 02 RW8-SP303-20250911  | 6528                | 7.84 | 17791               | 10.64  | 8068                | 14.50  |
| 03 MW203D2-20250909MS  | 6853                | 7.84 | 18950               | 10.64  | 8815                | 14.49  |
| 04 MW203D2-20250909MSD | 7569                | 7.85 | 20880               | 10.65  | 9823                | 14.49  |

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.: Q3086

Client ID: SSTDCCC0.4

Date Analyzed: 09/17/2025

Lab File ID: BN037740.D

Time Analyzed: 11:25

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            | 15814               | 17.235 | 13841               | 21.42 | 14132               | 23.753 |
| UPPER LIMIT            | 31628               | 17.735 | 27682               | 21.92 | 28264               | 24.253 |
| LOWER LIMIT            | 7907                | 16.735 | 6920.5              | 20.92 | 7066                | 23.253 |
| EPA SAMPLE NO.         |                     |        |                     |       |                     |        |
| 01 RW8-SP100-20250911  | 15420               | 17.24  | 14189               | 21.41 | 14177               | 23.75  |
| 02 RW8-SP303-20250911  | 13903               | 17.24  | 10625               | 21.41 | 10300               | 23.75  |
| 03 MW203D2-20250909MS  | 15289               | 17.24  | 13059               | 21.42 | 12757               | 23.75  |
| 04 MW203D2-20250909MSD | 16954               | 17.24  | 14253               | 21.42 | 13967               | 23.76  |

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.: Q3086

Client ID : SSTDCCC0.4

Date Analyzed: 09/18/2025

Lab File ID: BN037776.D

Time Analyzed: 14:32

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    | 7127                | 7.833 | 18984               | 10.64  | 8698                | 14.48  |
| UPPER LIMIT    | 14254               | 8.333 | 37968               | 11.137 | 17396               | 14.984 |
| LOWER LIMIT    | 3563.5              | 7.333 | 9492                | 10.137 | 4349                | 13.984 |
| EPA SAMPLE NO. |                     |       |                     |        |                     |        |
| 01 PB169692BS  | 6983                | 7.83  | 18103               | 10.64  | 7729                | 14.48  |

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.: Q3086

Client ID: SSTDCCC0.4

Date Analyzed: 09/18/2025

Lab File ID: BN037776.D

Time Analyzed: 14:32

Instrument ID: BNA\_N

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

01

|                | IS4 (PHN)<br>AREA # | RT #   | IS5 (CRY)<br>AREA # | RT #   | IS6 (PRY)<br>AREA # | RT #   |
|----------------|---------------------|--------|---------------------|--------|---------------------|--------|
| 12 HOUR STD    | 14526               | 17.235 | 10438               | 21.411 | 10600               | 23.744 |
| UPPER LIMIT    | 29052               | 17.735 | 20876               | 21.911 | 21200               | 24.244 |
| LOWER LIMIT    | 7263                | 16.735 | 5219                | 20.911 | 5300                | 23.244 |
| EPA SAMPLE NO. |                     |        |                     |        |                     |        |
| PB169692BS     | 12277               | 17.24  | 10644               | 21.41  | 10808               | 23.74  |

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.: Q3086

Client ID : SSTDCCC0.4

Date Analyzed: 09/19/2025

Lab File ID: BN037795.D

Time Analyzed: 02:47

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    | 7724                | 7.833 | 19361               | 10.64  | 8621                | 14.48  |
| UPPER LIMIT    | 15448               | 8.333 | 38722               | 11.137 | 17242               | 14.984 |
| LOWER LIMIT    | 3862                | 7.333 | 9680.5              | 10.137 | 4310.5              | 13.984 |
| EPA SAMPLE NO. |                     |       |                     |        |                     |        |
| 01 PB169692BL  | 7701                | 7.83  | 18665               | 10.64  | 7806                | 14.48  |

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.: Q3086

Client ID: SSTDCCC0.4

Date Analyzed: 09/19/2025

Lab File ID: BN037795.D

Time Analyzed: 02:47

Instrument ID: BNA\_N

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

01

|                | IS4 (PHN)<br>AREA # | RT #   | IS5 (CRY)<br>AREA # | RT #   | IS6 (PRY)<br>AREA # | RT #   |
|----------------|---------------------|--------|---------------------|--------|---------------------|--------|
| 12 HOUR STD    | 14380               | 17.235 | 11169               | 21.411 | 11840               | 23.742 |
| UPPER LIMIT    | 28760               | 17.735 | 22338               | 21.911 | 23680               | 24.242 |
| LOWER LIMIT    | 7190                | 16.735 | 5584.5              | 20.911 | 5920                | 23.242 |
| EPA SAMPLE NO. |                     |        |                     |        |                     |        |
| PB169692BL     | 12511               | 17.24  | 10022               | 21.41  | 10278               | 23.74  |

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:  | PB169692BL                    |                  | SDG No.:        | Q3086          |      |
| Lab Sample ID:     | PB169692BL                    |                  | 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 |
| BN037796.D        | 1         | 09/15/25 10:15 | 09/19/25 03:24 | PB169692      |

| 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 | 0.31  |           | 30 - 150 |      | 79%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.30  |           | 30 - 150 |      | 75%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.39  |           | 55 - 111 |      | 97%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.37  |           | 53 - 106 |      | 92%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.34  |           | 58 - 132 |      | 84%        | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 7700  | 7.833     |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 18700 | 10.637    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 7810  | 14.484    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 12500 | 17.236    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 10000 | 21.411    |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 10300 | 23.739    |          |      |            |          |

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:  | PB169692BS                    |                  | SDG No.:        | Q3086          |
| Lab Sample ID:     | PB169692BS                    |                  | 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 |
| BN037791.D        | 1         | 09/15/25 10:15 | 09/18/25 23:40 | PB169692      |

| 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 | 0.30  |           | 30 - 150 |      | 76%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.30  |           | 30 - 150 |      | 75%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.35  |           | 55 - 111 |      | 88%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.36  |           | 53 - 106 |      | 90%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.31  |           | 58 - 132 |      | 78%        | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 6980  | 7.833     |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 18100 | 10.637    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 7730  | 14.484    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 12300 | 17.235    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 10600 | 21.411    |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 10800 | 23.739    |          |      |            |          |

U = Not Detected

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MDL = Method Detection Limit

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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: | 09/09/25       |
| Project:           | NWIRP Bethpage 112G08005-WE13 | Date Received:  | 09/12/25       |
| Client Sample ID:  | MW203D2-20250909MS            | SDG No.:        | Q3086          |
| Lab Sample ID:     | Q3095-06MS                    | 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 |
| BN037743.D        | 1         | 09/15/25 10:15 | 09/17/25 13:14 | PB169692      |

| CAS Number                | Parameter               | Conc. | Qualifier | MDL      | LOD  | LOQ / CRQL | Units    |
|---------------------------|-------------------------|-------|-----------|----------|------|------------|----------|
| <b>TARGETS</b>            |                         |       |           |          |      |            |          |
| 123-91-1                  | 1,4-Dioxane             | 0.97  |           | 0.070    | 0.20 | 0.20       | ug/L     |
| <b>SURROGATES</b>         |                         |       |           |          |      |            |          |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.29  |           | 30 - 150 |      | 72%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.33  |           | 30 - 150 |      | 83%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.30  |           | 55 - 111 |      | 75%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.32  |           | 53 - 106 |      | 79%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.40  |           | 58 - 132 |      | 99%        | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 6850  | 7.84      |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 19000 | 10.637    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 8820  | 14.494    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 15300 | 17.235    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 13100 | 21.42     |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 12800 | 23.75     |          |      |            |          |

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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: | 09/09/25       |
| Project:           | NWIRP Bethpage 112G08005-WE13 | Date Received:  | 09/12/25       |
| Client Sample ID:  | MW203D2-20250909MSD           | SDG No.:        | Q3086          |
| Lab Sample ID:     | Q3095-07MSD                   | 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 |
| BN037744.D        | 1         | 09/15/25 10:15 | 09/17/25 13:59 | PB169692      |

| CAS Number                | Parameter               | Conc. | Qualifier | MDL      | LOD  | LOQ / CRQL | Units    |
|---------------------------|-------------------------|-------|-----------|----------|------|------------|----------|
| <b>TARGETS</b>            |                         |       |           |          |      |            |          |
| 123-91-1                  | 1,4-Dioxane             | 0.96  |           | 0.070    | 0.20 | 0.20       | ug/L     |
| <b>SURROGATES</b>         |                         |       |           |          |      |            |          |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.30  |           | 30 - 150 |      | 74%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.33  |           | 30 - 150 |      | 83%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.30  |           | 55 - 111 |      | 75%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.32  |           | 53 - 106 |      | 79%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.40  |           | 58 - 132 |      | 99%        | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |      |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 7570  | 7.847     |          |      |            |          |
| 1146-65-2                 | Naphthalene-d8          | 20900 | 10.648    |          |      |            |          |
| 15067-26-2                | Acenaphthene-d10        | 9820  | 14.494    |          |      |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 17000 | 17.235    |          |      |            |          |
| 1719-03-5                 | Chrysene-d12            | 14300 | 21.42     |          |      |            |          |
| 1520-96-3                 | Perylene-d12            | 14000 | 23.756    |          |      |            |          |

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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\_N\Methods\  
 Method File : 8270-SIM-BN091125.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 Last Update : Thu Sep 11 14:20:01 2025  
 Response Via : Initial Calibration

## Calibration Files

0.1 =BN037691.D 0.2 =BN037692.D 0.4 =BN037693.D 0.8 =BN037694.D 1.6 =BN037695.D 3.2 =BN037696.D 5 =BN037697.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           | 0.429          | 0.422 | 0.411 | 0.432 | 0.418 | 0.388 | 0.417 | 3.80  |       |
| 3)       | n-Nitrosodimet...     | 0.530          | 0.564 | 0.515 | 0.556 | 0.550 | 0.551 | 0.544 | 3.34  |       |
| 4) S     | 2-Fluorophenol        | 0.929          | 0.913 | 0.892 | 0.840 | 0.912 | 0.937 | 1.004 | 0.918 | 5.37  |
| 5) S     | Phenol-d6             | 1.191          | 1.193 | 1.156 | 1.088 | 1.193 | 1.225 | 1.289 | 1.191 | 5.15  |
| 6)       | bis(2-Chloroet...     | 1.076          | 1.075 | 1.070 | 1.019 | 1.078 | 1.066 | 1.063 | 1.064 | 1.94  |
|          |                       |                |       |       |       |       |       |       |       |       |
| 7) I     | Naphthalene-d8        | -----ISTD----- |       |       |       |       |       |       |       |       |
| 8) S     | Nitrobenzene-d5       | 0.287          | 0.279 | 0.285 | 0.272 | 0.296 | 0.303 | 0.325 | 0.293 | 6.05  |
| 9)       | Naphthalene           | 1.091          | 1.072 | 1.074 | 1.036 | 1.117 | 1.115 | 1.149 | 1.093 | 3.38  |
| 10)      | Hexachlorobuta...     | 0.199          | 0.196 | 0.197 | 0.189 | 0.203 | 0.199 | 0.202 | 0.198 | 2.29  |
| 11) SURR | 2-Methylnaphth...     | 0.497          | 0.497 | 0.498 | 0.478 | 0.522 | 0.550 | 0.668 | 0.530 | 12.31 |
| 12)      | 2-Methylnaphth...     | 0.643          | 0.651 | 0.659 | 0.633 | 0.697 | 0.712 | 0.740 | 0.676 | 5.91  |
|          |                       |                |       |       |       |       |       |       |       |       |
| 13) I    | Acenaphthene-d10      | -----ISTD----- |       |       |       |       |       |       |       |       |
| 14) S    | 2,4,6-Tribromo...     | 0.088          | 0.098 | 0.101 | 0.096 | 0.113 | 0.131 | 0.104 | 14.74 |       |
| 15) S    | 2-Fluorobiphenyl      | 1.663          | 1.666 | 1.649 | 1.586 | 1.689 | 1.674 | 1.785 | 1.673 | 3.54  |
| 16)      | Acenaphthylene        | 1.754          | 1.756 | 1.783 | 1.727 | 1.885 | 1.916 | 2.020 | 1.834 | 5.90  |
| 17)      | Acenaphthene          | 1.165          | 1.199 | 1.215 | 1.187 | 1.272 | 1.296 | 1.348 | 1.240 | 5.37  |
| 18)      | Fluorene              | 1.404          | 1.402 | 1.441 | 1.419 | 1.545 | 1.633 | 1.671 | 1.502 | 7.59  |
|          |                       |                |       |       |       |       |       |       |       |       |
| 19) I    | Phenanthrene-d10      | -----ISTD----- |       |       |       |       |       |       |       |       |
| 20)      | 4,6-Dinitro-2-...     | 0.032          | 0.033 | 0.033 | 0.040 | 0.046 | 0.037 | 16.34 |       |       |
| 21)      | 4-Bromophenyl-...     | 0.250          | 0.252 | 0.251 | 0.248 | 0.272 | 0.269 | 0.284 | 0.261 | 5.43  |
| 22)      | Hexachlorobenzene     | 0.301          | 0.308 | 0.294 | 0.280 | 0.305 | 0.302 | 0.310 | 0.300 | 3.39  |
| 23)      | Atrazine              | 0.145          | 0.151 | 0.152 | 0.146 | 0.169 | 0.190 | 0.209 | 0.166 | 15.00 |
| 24)      | Pentachlorophenol     | 0.074          | 0.074 | 0.073 | 0.086 | 0.100 | 0.115 | 0.087 | 19.46 |       |
| 25)      | Phenanthrene          | 1.228          | 1.215 | 1.223 | 1.173 | 1.285 | 1.321 | 1.371 | 1.259 | 5.47  |
| 26)      | Anthracene            | 1.038          | 1.029 | 1.070 | 1.030 | 1.140 | 1.221 | 1.281 | 1.116 | 9.09  |
| 27) SURR | Fluoranthene-d10      | 0.912          | 0.887 | 0.935 | 0.883 | 0.954 | 1.108 | 1.350 | 1.004 | 16.98 |
| 28)      | Fluoranthene          | 1.257          | 1.224 | 1.310 | 1.269 | 1.393 | 1.578 | 1.637 | 1.381 | 11.91 |
|          |                       |                |       |       |       |       |       |       |       |       |
| 29) I    | Chrysene-d12          | -----ISTD----- |       |       |       |       |       |       |       |       |
| 30)      | Pyrene                | 1.578          | 1.618 | 1.557 | 1.473 | 1.603 | 1.542 | 1.590 | 1.566 | 3.09  |
| 31) S    | Terphenyl-d14         | 0.806          | 0.833 | 0.801 | 0.754 | 0.812 | 0.784 | 0.892 | 0.812 | 5.30  |
| 32)      | Benzo(a)anthra...     | 1.386          | 1.345 | 1.347 | 1.297 | 1.410 | 1.459 | 1.522 | 1.395 | 5.46  |
| 33)      | Chrysene              | 1.523          | 1.491 | 1.496 | 1.402 | 1.510 | 1.506 | 1.543 | 1.496 | 3.01  |
| 34)      | Bis(2-ethylhex...     | 0.440          | 0.376 | 0.345 | 0.380 | 0.438 | 0.504 | 0.414 | 13.96 |       |
|          |                       |                |       |       |       |       |       |       |       |       |
| 35) I    | Perylene-d12          | -----ISTD----- |       |       |       |       |       |       |       |       |

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

Method File : 8270-SIM-BN091125.M

|       |                   |       |       |       |       |       |       |       |       |       |
|-------|-------------------|-------|-------|-------|-------|-------|-------|-------|-------|-------|
| 36)   | Indeno(1,2,3-c... | 1.517 | 1.491 | 1.543 | 1.532 | 1.791 | 1.865 | 2.025 | 1.681 | 12.58 |
| 37)   | Benzo(b)fluora... | 1.463 | 1.475 | 1.554 | 1.444 | 1.636 | 1.674 | 1.778 | 1.575 | 7.99  |
| 38)   | Benzo(k)fluora... | 1.516 | 1.543 | 1.524 | 1.501 | 1.705 | 1.696 | 1.756 | 1.606 | 6.73  |
| 39) C | Benzo(a)pyrene    | 1.155 | 1.172 | 1.201 | 1.151 | 1.314 | 1.367 | 1.467 | 1.261 | 9.78  |
| 40)   | Dibenzo(a,h)an... | 1.157 | 1.175 | 1.210 | 1.232 | 1.463 | 1.519 | 1.636 | 1.342 | 14.39 |
| 41)   | Benzo(g,h,i)pe... | 1.413 | 1.402 | 1.451 | 1.375 | 1.568 | 1.607 | 1.723 | 1.506 | 8.59  |

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

7C

SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: Alliance Contract: TETR06  
Lab Code: ACE SDG No.: Q3086  
Instrument ID: BNA\_N Calibration Date/Time: 09/17/2025 11:25  
Lab File ID: BN037740.D Init. Calib. Date(s): 09/11/2025 09/11/2025  
EPA Sample No.: SSTDCCC0.4 Init. Calib. Time(s): 09:56 13:33  
GC Column: ZB-GR ID: 0.25 (mm)

| COMPOUND                | RRF   | RRF0.4 | MIN<br>RRF | %D   | MAX%D |
|-------------------------|-------|--------|------------|------|-------|
| 2-Methylnaphthalene-d10 | 0.530 | 0.517  |            | -2.5 | 20.0  |
| Fluoranthene-d10        | 1.004 | 0.936  |            | -6.8 | 20.0  |
| 2-Fluorophenol          | 0.918 | 0.926  |            | 0.9  | 20.0  |
| Phenol-d6               | 1.191 | 1.176  |            | -1.3 | 20.0  |
| Nitrobenzene-d5         | 0.293 | 0.284  |            | -3.1 | 20.0  |
| 2-Fluorobiphenyl        | 1.673 | 1.620  |            | -3.2 | 20.0  |
| 2,4,6-Tribromophenol    | 0.104 | 0.123  |            | 18.3 | 20.0  |
| Terphenyl-d14           | 0.812 | 0.757  |            | -6.8 | 20.0  |
| 1,4-Dioxane             | 0.417 | 0.449  |            | 7.7  | 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>Q3086</u>                 |
| Instrument ID:  | <u>BNA_N</u>        | Calibration Date/Time: | <u>09/17/2025 21:50</u>      |
| Lab File ID:    | <u>BN037757.D</u>   | Init. Calib. Date(s):  | <u>09/11/2025 09/11/2025</u> |
| EPA Sample No.: | <u>SSTDCCC0.4EC</u> | Init. Calib. Time(s):  | <u>09:56 13:33</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 | 0.530 | 0.503  |            | -5.1 | 50.0  |
| Fluoranthene-d10        | 1.004 | 0.936  |            | -6.8 | 50.0  |
| 2-Fluorophenol          | 0.918 | 0.971  |            | 5.8  | 50.0  |
| Phenol-d6               | 1.191 | 1.159  |            | -2.7 | 50.0  |
| Nitrobenzene-d5         | 0.293 | 0.283  |            | -3.4 | 50.0  |
| 2-Fluorobiphenyl        | 1.673 | 1.646  |            | -1.6 | 50.0  |
| 2,4,6-Tribromophenol    | 0.104 | 0.122  |            | 17.3 | 50.0  |
| Terphenyl-d14           | 0.812 | 0.761  |            | -6.3 | 50.0  |
| 1,4-Dioxane             | 0.417 | 0.432  |            | 3.6  | 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.: Q3086  
 Instrument ID: BNA\_N Calibration Date/Time: 09/18/2025 14:32  
 Lab File ID: BN037776.D Init. Calib. Date(s): 09/11/2025 09/11/2025  
 EPA Sample No.: SSTDCCC0.4 Init. Calib. Time(s): 09:56 13:33  
 GC Column: ZB-GR ID: 0.25 (mm)

| COMPOUND                | RRF   | RRF0.4 | MIN<br>RRF | %D    | MAX%D |
|-------------------------|-------|--------|------------|-------|-------|
| 2-Methylnaphthalene-d10 | 0.530 | 0.497  |            | -6.2  | 20.0  |
| Fluoranthene-d10        | 1.004 | 0.839  |            | -16.4 | 20.0  |
| 2-Fluorophenol          | 0.918 | 0.890  |            | -3.0  | 20.0  |
| Phenol-d6               | 1.191 | 1.130  |            | -5.1  | 20.0  |
| Nitrobenzene-d5         | 0.293 | 0.298  |            | 1.7   | 20.0  |
| 2-Fluorobiphenyl        | 1.673 | 1.679  |            | 0.4   | 20.0  |
| 2,4,6-Tribromophenol    | 0.104 | 0.114  |            | 9.6   | 20.0  |
| Terphenyl-d14           | 0.812 | 0.801  |            | -1.4  | 20.0  |
| 1,4-Dioxane             | 0.417 | 0.428  |            | 2.6   | 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>Q3086</u>                 |
| Instrument ID:  | <u>BNA_N</u>        | Calibration Date/Time: | <u>09/19/2025 00:52</u>      |
| Lab File ID:    | <u>BN037793.D</u>   | Init. Calib. Date(s):  | <u>09/11/2025 09/11/2025</u> |
| EPA Sample No.: | <u>SSTDCCC0.4EC</u> | Init. Calib. Time(s):  | <u>09:56 13:33</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 | 0.530 | 0.496  |            | -6.4 | 50.0  |
| Fluoranthene-d10        | 1.004 | 0.954  |            | -5.0 | 50.0  |
| 2-Fluorophenol          | 0.918 | 0.923  |            | 0.5  | 50.0  |
| Phenol-d6               | 1.191 | 1.148  |            | -3.6 | 50.0  |
| Nitrobenzene-d5         | 0.293 | 0.295  |            | 0.7  | 50.0  |
| 2-Fluorobiphenyl        | 1.673 | 1.632  |            | -2.5 | 50.0  |
| 2,4,6-Tribromophenol    | 0.104 | 0.127  |            | 22.1 | 50.0  |
| Terphenyl-d14           | 0.812 | 0.785  |            | -3.3 | 50.0  |
| 1,4-Dioxane             | 0.417 | 0.436  |            | 4.6  | 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.: Q3086  
Instrument ID: BNA\_N Calibration Date/Time: 09/19/2025 02:47  
Lab File ID: BN037795.D Init. Calib. Date(s): 09/11/2025 09/11/2025  
EPA Sample No.: SSTDCCC0.4 Init. Calib. Time(s): 09:56 13:33  
GC Column: ZB-GR ID: 0.25 (mm)

| COMPOUND                | RRF   | RRF0.4 | MIN<br>RRF | %D    | MAX%D |
|-------------------------|-------|--------|------------|-------|-------|
| 2-Methylnaphthalene-d10 | 0.530 | 0.490  |            | -7.5  | 20.0  |
| Fluoranthene-d10        | 1.004 | 0.858  |            | -14.5 | 20.0  |
| 2-Fluorophenol          | 0.918 | 0.830  |            | -9.6  | 20.0  |
| Phenol-d6               | 1.191 | 1.152  |            | -3.3  | 20.0  |
| Nitrobenzene-d5         | 0.293 | 0.308  |            | 5.1   | 20.0  |
| 2-Fluorobiphenyl        | 1.673 | 1.696  |            | 1.4   | 20.0  |
| 2,4,6-Tribromophenol    | 0.104 | 0.120  |            | 15.4  | 20.0  |
| Terphenyl-d14           | 0.812 | 0.771  |            | -5.0  | 20.0  |
| 1,4-Dioxane             | 0.417 | 0.391  |            | -6.2  | 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>Q3086</u>                 |
| Instrument ID:  | <u>BNA_N</u>        | Calibration Date/Time: | <u>09/19/2025 13:03</u>      |
| Lab File ID:    | <u>BN037811.D</u>   | Init. Calib. Date(s):  | <u>09/11/2025 09/11/2025</u> |
| EPA Sample No.: | <u>SSTDCCC0.4EC</u> | Init. Calib. Time(s):  | <u>09:56 13:33</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 | 0.530 | 0.497  |            | -6.2 | 50.0  |
| Fluoranthene-d10        | 1.004 | 0.957  |            | -4.7 | 50.0  |
| 2-Fluorophenol          | 0.918 | 1.044  |            | 13.7 | 50.0  |
| Phenol-d6               | 1.191 | 1.176  |            | -1.3 | 50.0  |
| Nitrobenzene-d5         | 0.293 | 0.292  |            | -0.3 | 50.0  |
| 2-Fluorobiphenyl        | 1.673 | 1.615  |            | -3.5 | 50.0  |
| 2,4,6-Tribromophenol    | 0.104 | 0.131  |            | 26.0 | 50.0  |
| Terphenyl-d14           | 0.812 | 0.765  |            | -5.8 | 50.0  |
| 1,4-Dioxane             | 0.417 | 0.424  |            | 1.7  | 50.0  |

All other compounds must meet a minimum RRF of 0.010.