

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061522\  
 Data File : BN020170.D  
 Acq On : 15 Jun 2022 12:15  
 Operator : CG/JU  
 Sample : N3317-05  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :

Quant Time: Jun 16 01:11:01 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN060622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jun 16 00:50:14 2022  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.799  | 152  | 97271    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.587 | 136  | 458984   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.428 | 164  | 337243   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.169 | 188  | 760331   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.357 | 240  | 745703   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 23.668 | 264  | 699478   | 20.000 | ng/u1 | -0.01    |

|                               |        |     |        |        |       |       |
|-------------------------------|--------|-----|--------|--------|-------|-------|
| System Monitoring Compounds   |        |     |        |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.264  | 96  | 54     | 0.024  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.681  | 84  | 201    | 0.033  | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 0.000  | 99  | 0      | 0.000  | ng/u1 |       |
| 9) Bis-(2-Chloroethyl)eth...  | 0.000  | 67  | 0      | 0.000  | ng/u1 |       |
| 11) 2-Chlorophenol-d4         | 0.000  | 132 | 0      | 0.000  | ng/u1 |       |
| 15) 4-Methylphenol-d8         | 0.000  | 113 | 0      | 0.000  | ng/u1 |       |
| 21) Nitrobenzene-d5           | 0.000  | 128 | 0      | 0.000  | ng/u1 |       |
| 24) 2-Nitrophenol-d4          | 0.000  | 143 | 0      | 0.000  | ng/u1 |       |
| 28) 2,4-Dichlorophenol-d3     | 0.000  | 165 | 0      | 0.000  | ng/u1 |       |
| 31) 4-Chloroaniline-d4        | 10.593 | 131 | 539    | 0.051  | ng/u1 | -0.13 |
| 46) Dimethylphthalate-d6      | 0.000  | 166 | 0      | 0.000  | ng/u1 |       |
| 49) Acenaphthylene-d8         | 0.000  | 160 | 0      | 0.000  | ng/u1 |       |
| 54) 4-Nitrophenol-d4          | 0.000  | 143 | 0      | 0.000  | ng/u1 |       |
| 60) Fluorene-d10              | 0.000  | 176 | 0      | 0.000  | ng/u1 |       |
| 65) 4,6-Dinitro-2-methylph... | 0.000  | 200 | 0      | 0.000  | ng/u1 |       |
| 73) Anthracene-d10            | 17.169 | 188 | 760331 | 22.913 | ng/u1 | -0.09 |
| 81) Pyrene-d10                | 0.000  | 212 | 0      | 0.000  | ng/u1 |       |
| 92) Benzo(a)pyrene-d12        | 23.668 | 264 | 699478 | 20.438 | ng/u1 | 0.14  |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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