

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN101724\  
 Data File : BN034633.D  
 Acq On : 17 Oct 2024 22:33  
 Operator : RC/JU  
 Sample : P4346-10  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 EOAX0

Quant Time: Oct 17 23:49:45 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN101624.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 16 15:50:29 2024  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.664  | 152  | 300750   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.440 | 136  | 1120002  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.298 | 164  | 560188   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.039 | 188  | 974091   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.263 | 240  | 874665   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 24.151 | 264  | 1126913  | 20.000 | ng/u1 | 0.00     |

|                               |        |     |        |        |       |       |
|-------------------------------|--------|-----|--------|--------|-------|-------|
| System Monitoring Compounds   |        |     |        |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.217  | 96  | 14117  | 1.878  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.605  | 84  | 197478 | 8.999  | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 6.834  | 99  | 250022 | 10.130 | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.005  | 67  | 157930 | 10.373 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.199  | 132 | 205640 | 10.785 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.363  | 113 | 189314 | 9.842  | ng/u1 | 0.00  |
| 21) Nitrobenzene-d5           | 8.805  | 128 | 89425  | 11.473 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.528  | 143 | 68766  | 11.165 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.057 | 165 | 150597 | 10.608 | ng/u1 | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.569 | 131 | 253502 | 10.409 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 13.716 | 166 | 404603 | 11.417 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 13.987 | 160 | 514349 | 11.684 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.487 | 143 | 43224  | 6.547  | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.293 | 176 | 347167 | 11.326 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.404 | 200 | 6513   | 1.995  | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.139 | 188 | 491230 | 11.797 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.433 | 212 | 536306 | 11.942 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.957 | 264 | 616691 | 11.877 | ng/u1 | -0.01 |

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

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

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