

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN052024\  
 Data File : BN031579.D  
 Acq On : 20 May 2024 19:55  
 Operator : MA/JU  
 Sample : P2484-06  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 BH9T3

Quant Time: May 20 23:09:19 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN051524.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed May 15 02:37:21 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.699  | 152  | 542489   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.481 | 136  | 2391686  | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.334 | 164  | 1459896  | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.075 | 188  | 2821461  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.298 | 240  | 1721889  | 20.000 | ng/ul | -0.01    |
| 88) Perylene-d12              | 24.239 | 264  | 1713211  | 20.000 | ng/ul | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.205  | 96   | 68644    | 5.212  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.605  | 84   | 759148   | 18.852 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 6.857  | 99   | 1485114  | 32.256 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.034  | 67   | 929573   | 30.841 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.228  | 132  | 1159126  | 36.728 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.393  | 113  | 1228477  | 34.748 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 8.851  | 128  | 609834   | 38.688 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.563  | 143  | 588505   | 49.472 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.098 | 165  | 1108006  | 37.491 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.616 | 131  | 1754623  | 33.380 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.751 | 166  | 3578189  | 38.751 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.028 | 160  | 4252696  | 37.695 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.528 | 143  | 526634   | 29.737 | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.328 | 176  | 3051481  | 36.134 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.439 | 200  | 260798   | 25.816 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.175 | 188  | 4532612  | 37.570 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.469 | 212  | 4388433  | 51.334 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.039 | 264  | 2994145  | 39.376 | ng/ul | -0.01    |

Target Compounds Qvalue

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

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