

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111324\  
 Data File : BP023094.D  
 Acq On : 13 Nov 2024 23:36  
 Operator : RC/JU  
 Sample : P4724-08  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :

Quant Time: Nov 14 01:37:45 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP110924.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 13 04:49:01 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.581  | 152  | 81162    | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.328 | 136  | 304958   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.198 | 164  | 246190   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.010 | 188  | 639045   | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.439 | 240  | 752713   | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 24.657 | 264  | 904950   | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.140  | 96   | 4897     | 2.900  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.558  | 84   | 54597    | 10.883 | ng/ul | 0.01     |
| 7) Phenol-d5                  | 6.787  | 99   | 95887    | 15.980 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.928  | 67   | 58185    | 16.504 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.128  | 132  | 73517    | 16.729 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.293  | 113  | 79084    | 16.016 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 8.722  | 128  | 36058    | 16.984 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.440  | 143  | 43214    | 17.074 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.975  | 165  | 87576    | 16.522 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.481 | 131  | 105804   | 16.732 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.598 | 166  | 339172   | 19.115 | ng/ul | -0.01    |
| 49) Acenaphthylene-d8         | 13.881 | 160  | 353870   | 19.113 | ng/ul | -0.02    |
| 54) 4-Nitrophenol-d4          | 14.481 | 143  | 38062    | 14.274 | ng/ul | 0.03     |
| 60) Fluorene-d10              | 15.222 | 176  | 306072   | 19.675 | ng/ul | 0.01     |
| 65) 4,6-Dinitro-2-methylph... | 15.363 | 200  | 43060    | 11.248 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.098 | 188  | 544190   | 20.112 | ng/ul | -0.02    |
| 81) Pyrene-d10                | 19.492 | 212  | 714525   | 20.609 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.433 | 264  | 877960   | 20.367 | ng/ul | -0.01    |

Target Compounds Qvalue

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

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