

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111324\  
 Data File : BP023095.D  
 Acq On : 14 Nov 2024 00:17  
 Operator : RC/JU  
 Sample : P4724-05  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BH9C0

Quant Time: Nov 14 01:38:04 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  | 82391    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.328 | 136  | 314108   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.192 | 164  | 238586   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 16.998 | 188  | 589930   | 20.000 | ng/u1 | #-0.02   |
| 79) Chrysene-d12              | 21.433 | 240  | 683500   | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 24.651 | 264  | 812615   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.140  | 96   | 6229     | 3.634  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.552  | 84   | 88396    | 17.358 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.787  | 99   | 132244   | 21.710 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.928  | 67   | 79869    | 22.317 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.128  | 132  | 97887    | 21.942 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.299  | 113  | 114275   | 22.797 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.728  | 128  | 51030    | 23.336 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.440  | 143  | 61582    | 23.622 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.975  | 165  | 124753   | 22.851 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.481 | 131  | 146313   | 22.464 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.610 | 166  | 441678   | 25.685 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.881 | 160  | 468874   | 26.132 | ng/u1 | -0.02    |
| 54) 4-Nitrophenol-d4          | 14.475 | 143  | 50974    | 19.725 | ng/u1 | 0.02     |
| 60) Fluorene-d10              | 15.210 | 176  | 400888   | 26.591 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.369 | 200  | 58188    | 16.465 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.104 | 188  | 676674   | 27.091 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.492 | 212  | 861915   | 27.378 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.421 | 264  | 1094273  | 28.270 | ng/u1 | -0.02    |

Target Compounds Qvalue

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

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