

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP120423\  
 Data File : BP018619.D  
 Acq On : 06 Dec 2023 00:11  
 Operator : MA/JU  
 Sample : 05424-04  
 Misc :  
 ALS Vial : 55 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 YCGB9

Quant Time: Dec 06 00:57:41 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Dec 04 21:39:40 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.852  | 152  | 294747   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.646 | 136  | 1316641  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.481 | 164  | 932570   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.228 | 188  | 2151362  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.310 | 240  | 1824513  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.680 | 264  | 1735299  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.264  | 96   | 30874    | 3.569  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.687  | 84   | 244022   | 10.533 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.016  | 99   | 159963   | 5.408  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.187  | 67   | 511988   | 29.205 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.381  | 132  | 465670   | 20.176 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.558  | 113  | 334796   | 13.986 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.010  | 128  | 335872   | 28.662 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.734  | 143  | 332873   | 27.398 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.269 | 165  | 647139   | 26.049 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.787 | 131  | 874688   | 26.941 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.898 | 166  | 2701709  | 32.460 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.175 | 160  | 2808938  | 30.648 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.675 | 143  | 62506    | 4.575  | ng/u1 | -0.01    |
| 60) Fluorene-d10              | 15.475 | 176  | 2346340  | 33.584 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.587 | 200  | 327202   | 25.365 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.328 | 188  | 3911259  | 34.582 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.563 | 212  | 4571912  | 32.191 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.527 | 264  | 3547808  | 35.064 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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