

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP120423\  
 Data File : BP018648.D  
 Acq On : 06 Dec 2023 21:27  
 Operator : MA/JU  
 Sample : 05439-15  
 Misc :  
 ALS Vial : 85 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0AD7

Quant Time: Dec 06 22:44:20 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     | 7.851  | 152  | 366396   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.645 | 136  | 1568672  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.480 | 164  | 1006071  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.227 | 188  | 1941189  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.315 | 240  | 1422579  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.686 | 264  | 1596529  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.263  | 96   | 32560    | 3.028  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.687  | 84   | 93235    | 3.237  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.016  | 99   | 183740   | 4.997  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.187  | 67   | 530446   | 24.341 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.381  | 132  | 521726   | 18.184 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.563  | 113  | 356999   | 11.997 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.010  | 128  | 357926   | 25.636 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.734  | 143  | 368036   | 25.426 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.269 | 165  | 663703   | 22.424 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.786 | 131  | 579755   | 14.988 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.898 | 166  | 2554956  | 28.454 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.174 | 160  | 2629469  | 26.594 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.680 | 143  | 57728    | 3.917  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.474 | 176  | 2105499  | 27.935 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.592 | 200  | 275857   | 23.700 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.327 | 188  | 3033120  | 29.721 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.562 | 212  | 3139043  | 28.347 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.533 | 264  | 2854626  | 30.666 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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