

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP120723\  
 Data File : BP018668.D  
 Acq On : 07 Dec 2023 11:09  
 Operator : MA/JU  
 Sample : PB157420BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK420

Quant Time: Dec 07 22:01:30 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     | 7.846  | 152  | 335121   | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8            | 10.640 | 136  | 1275521  | 20.000 | ng/u1 | -0.01    |
| 38) Acenaphthene-d10          | 14.475 | 164  | 799769   | 20.000 | ng/u1 | -0.01    |
| 64) Phenanthrene-d10          | 17.228 | 188  | 1857883  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.316 | 240  | 2032653  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.686 | 264  | 2426211  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.264  | 96   | 45642    | 4.641  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.681  | 84   | 591131   | 22.441 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.011  | 99   | 697515   | 20.740 | ng/u1 | -0.01    |
| 9) Bis-(2-Chloroethyl)eth...  | 7.181  | 67   | 411117   | 20.625 | ng/u1 | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.375  | 132  | 543633   | 20.716 | ng/u1 | -0.01    |
| 15) 4-Methylphenol-d8         | 8.558  | 113  | 533974   | 19.619 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.005  | 128  | 254305   | 22.401 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.728  | 143  | 274010   | 23.280 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.263 | 165  | 508287   | 21.120 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.781 | 131  | 751986   | 23.908 | ng/u1 | -0.01    |
| 46) Dimethylphthalate-d6      | 13.893 | 166  | 1660543  | 23.263 | ng/u1 | -0.01    |
| 49) Acenaphthylene-d8         | 14.169 | 160  | 1760179  | 22.394 | ng/u1 | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.675 | 143  | 301349   | 25.719 | ng/u1 | -0.01    |
| 60) Fluorene-d10              | 15.469 | 176  | 1178637  | 19.671 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.587 | 200  | 215047   | 19.304 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.322 | 188  | 2356370  | 24.125 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.563 | 212  | 3009880  | 19.023 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.533 | 264  | 3352826  | 23.701 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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