

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP120423\  
 Data File : BP018592.D  
 Acq On : 05 Dec 2023 08:55  
 Operator : MA/JU  
 Sample : PB157362BL  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK362

Quant Time: Dec 05 10:24:47 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  | 416134   | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8            | 10.646 | 136  | 1818523  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.481 | 164  | 1261991  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.233 | 188  | 2880683  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.321 | 240  | 2022233  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.698 | 264  | 2260549  | 20.000 | ng/u1 | 0.01     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.264  | 96   | 64958    | 5.319  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.681  | 84   | 869366   | 26.579 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.010  | 99   | 1110281  | 26.587 | ng/u1 | -0.01    |
| 9) Bis-(2-Chloroethyl)eth...  | 7.181  | 67   | 689591   | 27.861 | ng/u1 | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.375  | 132  | 840085   | 25.781 | ng/u1 | -0.01    |
| 15) 4-Methylphenol-d8         | 8.557  | 113  | 893986   | 26.452 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.004  | 128  | 425465   | 26.287 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.728  | 143  | 430042   | 25.627 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.263 | 165  | 866562   | 25.255 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.787 | 131  | 1024360  | 22.843 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.898 | 166  | 3178282  | 28.218 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.175 | 160  | 3376419  | 27.224 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.681 | 143  | 508196   | 27.487 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.475 | 176  | 2716167  | 28.729 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.592 | 200  | 380082   | 22.005 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.333 | 188  | 4274293  | 28.224 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.569 | 212  | 4790849  | 30.435 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.545 | 264  | 3664538  | 27.802 | ng/u1 | 0.01     |

Target Compounds Qvalue

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

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