

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092923\  
 Data File : BP017451.D  
 Acq On : 29 Sep 2023 18:27  
 Operator : MA/JU  
 Sample : PB155865BL  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK865

Quant Time: Sep 30 00:25:01 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Sep 29 22:31:03 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.952  | 152  | 368731   | 20.000 | ng/ul | -0.01    |
| 20) Naphthalene-d8            | 10.793 | 136  | 1423419  | 20.000 | ng/ul | -0.01    |
| 38) Acenaphthene-d10          | 14.645 | 164  | 806583   | 20.000 | ng/ul | -0.01    |
| 64) Phenanthrene-d10          | 17.434 | 188  | 1715905  | 20.000 | ng/ul | -0.01    |
| 79) Chrysene-d12              | 21.563 | 240  | 1710912  | 20.000 | ng/ul | -0.01    |
| 88) Perylene-d12              | 24.145 | 264  | 1991120  | 20.000 | ng/ul | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.293  | 96   | 40143    | 4.472  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.728  | 84   | 577485   | 23.010 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 7.105  | 99   | 709729   | 23.436 | ng/ul | -0.01    |
| 9) Bis-(2-Chloroethyl)eth...  | 7.293  | 67   | 429354   | 23.493 | ng/ul | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.469  | 132  | 585466   | 24.375 | ng/ul | -0.01    |
| 15) 4-Methylphenol-d8         | 8.675  | 113  | 554900   | 23.568 | ng/ul | -0.01    |
| 21) Nitrobenzene-d5           | 9.163  | 128  | 279585   | 27.169 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.881  | 143  | 316233   | 28.302 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.404 | 165  | 520183   | 24.641 | ng/ul | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.957 | 131  | 743657   | 24.319 | ng/ul | -0.01    |
| 46) Dimethylphthalate-d6      | 14.063 | 166  | 1506966  | 26.081 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.345 | 160  | 1727644  | 25.997 | ng/ul | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.881 | 143  | 225488   | 22.400 | ng/ul | -0.01    |
| 60) Fluorene-d10              | 15.651 | 176  | 1258043  | 25.290 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.792 | 200  | 189094   | 19.552 | ng/ul | -0.01    |
| 73) Anthracene-d10            | 17.534 | 188  | 2044934  | 26.871 | ng/ul | -0.01    |
| 81) Pyrene-d10                | 19.786 | 212  | 2463232  | 26.623 | ng/ul | -0.01    |
| 92) Benzo(a)pyrene-d12        | 23.974 | 264  | 2501788  | 26.097 | ng/ul | -0.01    |

Target Compounds Qvalue

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

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