

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP120123\  
 Data File : BP018537.D  
 Acq On : 03 Dec 2023 03:56  
 Operator : MA/JU  
 Sample : PB157193BL  
 Misc :  
 ALS Vial : 68 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK193

Quant Time: Dec 03 23:48:56 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Dec 01 21:56:43 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.857  | 152  | 405471   | 20.000 | ng/u1 | -0.01    |
| 20) Naphthalene-d8            | 10.657 | 136  | 1798858  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.487 | 164  | 1224310  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.233 | 188  | 2631748  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.316 | 240  | 1786344  | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 23.692 | 264  | 1733290  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.269  | 96   | 74198    | 6.236  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.687  | 84   | 1053572  | 33.057 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.028  | 99   | 1269845  | 31.207 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.193  | 67   | 795055   | 32.967 | ng/u1 | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.387  | 132  | 964128   | 30.366 | ng/u1 | -0.01    |
| 15) 4-Methylphenol-d8         | 8.569  | 113  | 1028647  | 31.237 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.022  | 128  | 491655   | 30.708 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.740  | 143  | 512644   | 30.884 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.275 | 165  | 992521   | 29.242 | ng/u1 | -0.01    |
| 31) 4-Chloroaniline-d4        | 10.793 | 131  | 1526905  | 34.423 | ng/u1 | -0.01    |
| 46) Dimethylphthalate-d6      | 13.904 | 166  | 3557171  | 32.554 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.181 | 160  | 3798480  | 31.569 | ng/u1 | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.687 | 143  | 543906   | 30.324 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.481 | 176  | 2934559  | 31.994 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.592 | 200  | 401778   | 25.461 | ng/u1 | -0.01    |
| 73) Anthracene-d10            | 17.333 | 188  | 4479151  | 32.374 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.569 | 212  | 4700657  | 33.805 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.533 | 264  | 3395043  | 33.593 | ng/u1 | -0.02    |

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

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