

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM120823\  
 Data File : BM043191.D  
 Acq On : 08 Dec 2023 11:48  
 Operator : MA/JU  
 Sample : PB157621BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SBLK621

Quant Time: Dec 08 12:20:10 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM120723.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Dec 08 00:42:52 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.010  | 152  | 493311   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.816 | 136  | 2169580  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.633 | 164  | 1523319  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.368 | 188  | 3357414  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.533 | 240  | 2972396  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.932 | 264  | 3176057  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.398  | 96   | 62379    | 4.963  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.816  | 84   | 842795   | 22.598 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.169  | 99   | 1152580  | 23.534 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.345  | 67   | 775298   | 25.572 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.539  | 132  | 921967   | 24.694 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.716  | 113  | 936904   | 24.181 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.180  | 128  | 483870   | 26.166 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.898  | 143  | 479604   | 26.896 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.433 | 165  | 944526   | 23.764 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.957 | 131  | 1125294  | 19.677 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.051 | 166  | 3574480  | 26.769 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.327 | 160  | 3931135  | 25.446 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.821 | 143  | 540391   | 22.584 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.621 | 176  | 3032993  | 26.056 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.739 | 200  | 338995   | 18.345 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.468 | 188  | 4795225  | 26.400 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.750 | 212  | 5350322  | 28.131 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.779 | 264  | 4820823  | 26.012 | ng/u1 | 0.00     |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

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

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