

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM031323\  
 Data File : BM039009.D  
 Acq On : 14 Mar 2023 12:06  
 Operator : CG/JU  
 Sample : PB151312BL  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SBLK312

Quant Time: Mar 15 01:05:22 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM031323.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Mar 13 16:25:27 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.987  | 152  | 410931   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.798 | 136  | 1874377  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.621 | 164  | 1206035  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.368 | 188  | 2515387  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.550 | 240  | 1946113  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.997 | 264  | 1815227  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.369  | 96   | 82709    | 7.019  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.793  | 84   | 986313   | 29.591 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.140  | 99   | 1342147  | 33.929 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.310  | 67   | 903833   | 35.355 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.510  | 132  | 1041536  | 35.369 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.692  | 113  | 1052059  | 33.994 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.151  | 128  | 550454   | 35.128 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.875  | 143  | 581805   | 36.851 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.410 | 165  | 959714   | 33.207 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.933 | 131  | 993308   | 21.250 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.033 | 166  | 3185436  | 34.872 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.321 | 160  | 3661236  | 34.510 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.810 | 143  | 576521   | 30.292 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.615 | 176  | 2806470  | 35.212 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.727 | 200  | 445829   | 27.984 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.468 | 188  | 4314084  | 35.572 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.762 | 212  | 4615691  | 40.510 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.844 | 264  | 3410137  | 36.594 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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