

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM030223\  
 Data File : BM038843.D  
 Acq On : 02 Mar 2023 19:21  
 Operator : CG/JU  
 Sample : 01710-14  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DBAB1

Quant Time: Mar 03 00:16:53 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM022823.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Mar 01 02:09:21 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc Units   | Dev(Min) |
|-------------------------------|--------|------|----------|--------------|----------|
| -----                         |        |      |          |              |          |
| Internal Standards            |        |      |          |              |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.010  | 152  | 247860   | 20.000 ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.822 | 136  | 1085967  | 20.000 ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.645 | 164  | 674222   | 20.000 ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.386 | 188  | 1486668  | 20.000 ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.562 | 240  | 1480606  | 20.000 ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.015 | 264  | 1805195  | 20.000 ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |              |          |
| 3) 1,4-Dioxane-d8             | 3.381  | 96   | 18452    | 2.780 ng/uL  | 0.00     |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000 ng/u1  |          |
| 7) Phenol-d5                  | 7.157  | 99   | 319143   | 13.734 ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.333  | 67   | 218957   | 14.747 ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.534  | 132  | 263372   | 15.210 ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.710  | 113  | 251084   | 13.716 ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.175  | 128  | 120779   | 15.842 ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.898  | 143  | 112395   | 15.733 ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.439 | 165  | 246045   | 14.822 ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.957 | 131  | 235379   | 8.507 ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.057 | 166  | 829269   | 15.880 ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.339 | 160  | 944239   | 15.470 ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.821 | 143  | 77315    | 7.396 ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.633 | 176  | 735388   | 15.967 ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.739 | 200  | 70645    | 9.140 ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.486 | 188  | 1145261  | 15.623 ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.768 | 212  | 1436632  | 17.642 ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.850 | 264  | 1497280  | 16.459 ng/u1 | -0.02    |

Target Compounds Qvalue

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

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