

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM030223\  
 Data File : BM038837.D  
 Acq On : 02 Mar 2023 15:36  
 Operator : CG/JU  
 Sample : 01710-01  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DBAA0

Quant Time: Mar 03 00:16:01 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  | 217036   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.828 | 136  | 937225   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.645 | 164  | 572332   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.386 | 188  | 1237465  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.562 | 240  | 1206167  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.021 | 264  | 1451127  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.381  | 96   | 23630    | 4.065  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.805  | 84   | 281015   | 17.585 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.157  | 99   | 383905   | 18.867 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.334  | 67   | 256101   | 19.698 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.534  | 132  | 300024   | 19.788 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.710  | 113  | 297277   | 18.545 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.175  | 128  | 135078   | 20.530 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.904  | 143  | 126354   | 20.495 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.439 | 165  | 280104   | 19.552 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.957 | 131  | 456589   | 19.120 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.057 | 166  | 1040248  | 23.467 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.339 | 160  | 1146560  | 22.130 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.822 | 143  | 132821   | 14.968 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.633 | 176  | 916969   | 23.455 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.739 | 200  | 81964    | 12.740 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.486 | 188  | 1447549  | 23.723 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.774 | 212  | 1718980  | 25.912 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.862 | 264  | 1777654  | 24.309 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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