

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM112822\  
 Data File : BM037761.D  
 Acq On : 28 Nov 2022 16:01  
 Operator : CG/JU  
 Sample : N5710-16  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C0L02

Quant Time: Nov 28 22:29:11 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM112322.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Nov 28 22:24:47 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.110  | 152  | 221069   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.928 | 136  | 1040796  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.733 | 164  | 693803   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.468 | 188  | 1429807  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.633 | 240  | 1292059  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.127 | 264  | 1111763  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.434  | 96   | 30422    | 5.707  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.875  | 84   | 184490   | 10.633 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.257  | 99   | 684732   | 31.047 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.428  | 67   | 461186   | 31.347 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.634  | 132  | 520735   | 33.979 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.816  | 113  | 571764   | 32.578 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.281  | 128  | 292766   | 35.881 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.004 | 143  | 315134   | 39.197 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.545 | 165  | 532825   | 36.215 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.063 | 131  | 518141   | 21.223 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.139 | 166  | 1755486  | 35.722 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.433 | 160  | 2090585  | 36.041 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.916 | 143  | 322682   | 29.809 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.722 | 176  | 1582357  | 36.765 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.827 | 200  | 218268   | 28.378 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.568 | 188  | 2467211  | 36.900 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.851 | 212  | 2714741  | 41.346 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.968 | 264  | 2120077  | 37.048 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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