

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM112822\  
 Data File : BM037764.D  
 Acq On : 28 Nov 2022 17:52  
 Operator : CG/JU  
 Sample : N5710-14  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C0KZ7

Quant Time: Nov 28 22:29:35 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  | 216482   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.927 | 136  | 1013149  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.733 | 164  | 645239   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.468 | 188  | 1343904  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.633 | 240  | 1273011  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.127 | 264  | 1099074  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.434  | 96   | 31274    | 5.991  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.875  | 84   | 152079   | 8.951  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.257  | 99   | 675998   | 31.300 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.428  | 67   | 454946   | 31.578 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.634  | 132  | 511282   | 34.069 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.816  | 113  | 542842   | 31.586 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.281  | 128  | 280410   | 35.305 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.004 | 143  | 292332   | 37.353 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.545 | 165  | 493090   | 34.429 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.063 | 131  | 463319   | 19.495 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.139 | 166  | 1572696  | 34.411 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.427 | 160  | 1877096  | 34.796 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.915 | 143  | 285388   | 28.349 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.721 | 176  | 1424714  | 35.593 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.827 | 200  | 193361   | 26.746 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.568 | 188  | 2227014  | 35.436 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.850 | 212  | 2502625  | 38.686 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.968 | 264  | 2010017  | 35.530 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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