

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM112522\  
 Data File : BM037712.D  
 Acq On : 26 Nov 2022 01:23  
 Operator : CG/JU  
 Sample : N5591-10  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C9822

Quant Time: Nov 26 04:18:57 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM112322.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Nov 24 01:27:15 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.110  | 152  | 198129   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.934 | 136  | 981568   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.739 | 164  | 618690   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.474 | 188  | 1249580  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.633 | 240  | 1205679  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 24.127 | 264  | 1167234  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.434  | 96   | 32233    | 6.747  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.869  | 84   | 528204   | 33.967 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.257  | 99   | 717807   | 36.315 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.434  | 67   | 502621   | 38.119 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.634  | 132  | 513552   | 37.390 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.816  | 113  | 583302   | 37.084 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.287  | 128  | 280783   | 36.489 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.010 | 143  | 270558   | 35.683 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.545 | 165  | 479146   | 34.531 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.069 | 131  | 795665   | 34.556 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 14.145 | 166  | 1662020  | 37.926 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.433 | 160  | 1946811  | 37.637 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.916 | 143  | 331643   | 34.357 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.727 | 176  | 1460999  | 38.066 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.833 | 200  | 191527   | 28.493 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.574 | 188  | 2277160  | 38.969 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.851 | 212  | 2537147  | 41.410 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.968 | 264  | 2320281  | 38.619 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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