

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM030624\  
 Data File : BM044533.D  
 Acq On : 07 Mar 2024 01:32  
 Operator : MA/JU  
 Sample : P1676-10  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 E06Z6

Quant Time: Mar 07 03:01:55 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM030524.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Mar 05 23:05:11 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.728  | 152  | 275708   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.510 | 136  | 1219434  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.369 | 164  | 790964   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.115 | 188  | 1633908  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.303 | 240  | 1377966  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.550 | 264  | 1312541  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.287  | 96   | 29297    | 3.604  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.681  | 84   | 426216   | 17.814 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.898  | 99   | 566699   | 19.992 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.081  | 67   | 381096   | 21.766 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.263  | 132  | 433345   | 20.803 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.428  | 113  | 465203   | 20.897 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 8.887  | 128  | 237642   | 23.143 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.604  | 143  | 258635   | 23.364 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.134 | 165  | 427305   | 22.681 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.657 | 131  | 691726   | 22.107 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.786 | 166  | 1525317  | 24.584 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.057 | 160  | 1738918  | 24.506 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.574 | 143  | 252851   | 19.651 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.363 | 176  | 1342029  | 25.739 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.486 | 200  | 215771   | 20.344 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.215 | 188  | 2088047  | 26.736 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.509 | 212  | 2395594  | 28.345 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.409 | 264  | 1884085  | 27.310 | ng/u1 | -0.01    |

Target Compounds Qvalue

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

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