

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM080924\  
 Data File : BM047121.D  
 Acq On : 09 Aug 2024 14:58  
 Operator : RC/JU  
 Sample : P3480-15  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 E0AB9

Quant Time: Aug 09 17:19:46 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM080224.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Aug 09 01:30:34 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.386  | 152  | 233677   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.139 | 136  | 930710   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.039 | 164  | 576996   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 16.803 | 188  | 1179003  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.033 | 240  | 1030024  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.738 | 264  | 1200989  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.016  | 96   | 22872    | 3.878  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.410  | 84   | 114348   | 7.107  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.592  | 99   | 156402   | 8.811  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.745  | 67   | 351398   | 30.086 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 6.928  | 132  | 400814   | 26.656 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.104  | 113  | 270077   | 19.206 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.528  | 128  | 213988   | 28.652 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.239  | 143  | 230724   | 28.289 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.780  | 165  | 418332   | 26.715 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.292 | 131  | 465008   | 22.595 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.468 | 166  | 1196598  | 27.123 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.727 | 160  | 1386396  | 28.153 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.292 | 143  | 51568    | 6.045  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.045 | 176  | 1031229  | 27.325 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.186 | 200  | 164480   | 22.081 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 16.898 | 188  | 1605365  | 28.723 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.215 | 212  | 1789991  | 30.358 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.556 | 264  | 1905689  | 30.134 | ng/u1 | 0.00     |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

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

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