

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM080624\  
 Data File : BM047035.D  
 Acq On : 06 Aug 2024 23:19  
 Operator : RC/JU  
 Sample : P3328-15  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 E20L8

Quant Time: Aug 07 01:24:04 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM080224.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 07 01:21:15 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.398  | 152  | 281971   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.157 | 136  | 1076781  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.057 | 164  | 678705   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 16.815 | 188  | 1387496  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.045 | 240  | 1100858  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.762 | 264  | 1210269  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.022  | 96   | 20078    | 2.821  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.410  | 84   | 304142   | 15.666 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.604  | 99   | 434163   | 20.269 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.757  | 67   | 262406   | 18.619 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 6.940  | 132  | 343370   | 18.924 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.116  | 113  | 353187   | 20.814 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.545  | 128  | 165322   | 19.133 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.257  | 143  | 184143   | 19.515 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.792  | 165  | 358899   | 19.811 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.304 | 131  | 476275   | 20.003 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.480 | 166  | 1091282  | 21.029 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.745 | 160  | 1251547  | 21.606 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.298 | 143  | 170577   | 17.000 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.063 | 176  | 936881   | 21.104 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.198 | 200  | 74072    | 8.450  | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 16.910 | 188  | 1425290  | 21.669 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.227 | 212  | 1601170  | 25.408 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.580 | 264  | 1450729  | 22.764 | ng/u1 | 0.01     |

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

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

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