

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG080824\  
 Data File : BG062332.D  
 Acq On : 8 Aug 2024 16:48  
 Operator : MA/JU  
 Sample : P3437-10  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 E05X4

Quant Time: Aug 08 18:05:46 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG080524.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 07 01:46:39 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.959  | 152  | 70348    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.749 | 136  | 340748   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.573 | 164  | 261167   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.311 | 188  | 550847   | 20.000 | ng/u1 | -0.01    |
| 79) Chrysene-d12              | 21.552 | 240  | 538662   | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12              | 24.636 | 264  | 628310   | 20.000 | ng/u1 | -0.02    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.424  | 96   | 7082     | 4.341  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.829  | 84   | 46834    | 9.125  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.107  | 99   | 62510    | 9.310  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.283  | 67   | 132066   | 31.952 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.489  | 132  | 146204   | 30.391 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.646  | 113  | 120142   | 21.482 | ng/u1 | -0.01    |
| 21) Nitrobenzene-d5           | 9.110  | 128  | 81909    | 38.984 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.827  | 143  | 88882    | 44.969 | ng/u1 | -0.01    |
| 28) 2,4-Dichlorophenol-d3     | 10.367 | 165  | 194058   | 33.804 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.878 | 131  | 155395   | 18.663 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.986 | 166  | 649827   | 32.250 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.268 | 160  | 718309   | 31.441 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.749 | 143  | 26833    | 9.512  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.566 | 176  | 580932   | 32.119 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.672 | 200  | 80159    | 49.136 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.410 | 188  | 855906   | 32.215 | ng/u1 | -0.01    |
| 81) Pyrene-d10                | 19.696 | 212  | 1090482  | 34.060 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.430 | 264  | 1167913  | 34.772 | ng/u1 | -0.01    |

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

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

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