

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062522\  
 Data File : BG054063.D  
 Acq On : 25 Jun 2022 16:02  
 Operator : CG/JU  
 Sample : N3401-11  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 C0AE3

Quant Time: Jun 27 01:45:49 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062022.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 20 15:29:13 2022  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 8.070  | 152  | 23430    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.879 | 136  | 111342   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.698 | 164  | 89042    | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.436 | 188  | 229523   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.713 | 240  | 251538   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 24.968 | 264  | 247525   | 20.000 | ng/u1 | -0.01    |

|                               |        |     |        |        |       |       |
|-------------------------------|--------|-----|--------|--------|-------|-------|
| System Monitoring Compounds   |        |     |        |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.493  | 96  | 2466   | 5.008  | ng/uL | -0.01 |
| 4) Pyridine-d5                | 3.916  | 84  | 11814  | 8.821  | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 7.230  | 99  | 13153  | 7.836  | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.389  | 67  | 35393  | 34.955 | ng/u1 | -0.01 |
| 11) 2-Chlorophenol-d4         | 7.606  | 132 | 34744  | 25.957 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.770  | 113 | 26655  | 18.624 | ng/u1 | 0.00  |
| 21) Nitrobenzene-d5           | 9.228  | 128 | 24252  | 29.345 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.956  | 143 | 25853  | 29.442 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.497 | 165 | 53128  | 27.347 | ng/u1 | 0.00  |
| 31) 4-Chloroaniline-d4        | 11.008 | 131 | 68093  | 27.744 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 14.093 | 166 | 219650 | 31.357 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 14.392 | 160 | 236956 | 31.109 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.886 | 143 | 5115   | 5.333  | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.685 | 176 | 198320 | 31.519 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.797 | 200 | 31188  | 24.909 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.536 | 188 | 347872 | 33.560 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.821 | 212 | 448849 | 34.224 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 24.745 | 264 | 428317 | 34.281 | ng/u1 | -0.01 |

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

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

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