

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG093021\  
 Data File : BG050341.D  
 Acq On : 1 Oct 2021 3:27  
 Operator : CG/JU  
 Sample : M3777-04  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 C0B70

Quant Time: Oct 01 04:11:05 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG093021.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Sep 30 16:27:24 2021  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 8.272  | 152  | 39120    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 11.098 | 136  | 168796   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.888 | 164  | 120468   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.632 | 188  | 290729   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.933 | 240  | 275574   | 20.000 | ng/u1 | -0.02    |
| 88) Perylene-d12          | 25.358 | 264  | 267697   | 20.000 | ng/u1 | -0.02    |

|                               |        |     |        |        |       |      |
|-------------------------------|--------|-----|--------|--------|-------|------|
| System Monitoring Compounds   |        |     |        |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.637  | 96  | 6032   | 5.665  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 4.066  | 84  | 8482   | 2.673  | ng/u1 | 0.00 |
| 7) Phenol-d5                  | 7.403  | 99  | 30194  | 8.834  | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.585  | 67  | 86158  | 40.054 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.796  | 132 | 75372  | 31.473 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 8.960  | 113 | 53563  | 20.705 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5           | 9.442  | 128 | 50008  | 42.350 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 10.164 | 143 | 50163  | 39.582 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.705 | 165 | 85420  | 34.133 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4        | 11.222 | 131 | 24832  | 6.946  | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6      | 14.283 | 166 | 327195 | 39.686 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.588 | 160 | 401261 | 38.755 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 15.047 | 143 | 14136  | 8.902  | ng/u1 | 0.00 |
| 60) Fluorene-d10              | 15.875 | 176 | 299735 | 40.352 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.975 | 200 | 56086  | 38.678 | ng/u1 | 0.00 |
| 73) Anthracene-d10            | 17.732 | 188 | 520963 | 42.416 | ng/u1 | 0.00 |
| 81) Pyrene-d10                | 20.006 | 212 | 625570 | 42.692 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12        | 25.129 | 264 | 534145 | 39.682 | ng/u1 | 0.00 |

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

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

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