

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG031822\  
 Data File : BG052722.D  
 Acq On : 18 Mar 2022 22:54  
 Operator : CG/JU  
 Sample : N1882-19  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 DBRY9

Quant Time: Mar 19 00:02:13 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG031822.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Mar 18 00:55:42 2022  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 8.148  | 152  | 31479    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.962 | 136  | 132191   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.768 | 164  | 92136    | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.512 | 188  | 219536   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.794 | 240  | 245212   | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12          | 25.113 | 264  | 242803   | 20.000 | ng/u1 | -0.01    |

|                               |        |     |        |        |       |      |
|-------------------------------|--------|-----|--------|--------|-------|------|
| System Monitoring Compounds   |        |     |        |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.560  | 96  | 6630   | 7.440  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.983  | 84  | 84253  | 37.869 | ng/u1 | 0.00 |
| 7) Phenol-d5                  | 7.290  | 99  | 127312 | 45.155 | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.467  | 67  | 78262  | 45.449 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.678  | 132 | 85264  | 44.538 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 8.841  | 113 | 93281  | 45.673 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5           | 9.305  | 128 | 41834  | 43.963 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 10.034 | 143 | 45515  | 45.349 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.574 | 165 | 91670  | 41.984 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4        | 11.085 | 131 | 122946 | 42.604 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6      | 14.169 | 166 | 315086 | 44.514 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.463 | 160 | 374918 | 43.589 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.933 | 143 | 51594  | 43.353 | ng/u1 | 0.00 |
| 60) Fluorene-d10              | 15.755 | 176 | 274696 | 44.198 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.861 | 200 | 38767  | 34.700 | ng/u1 | 0.00 |
| 73) Anthracene-d10            | 17.612 | 188 | 456761 | 44.680 | ng/u1 | 0.00 |
| 81) Pyrene-d10                | 19.891 | 212 | 575926 | 41.984 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12        | 24.890 | 264 | 552587 | 44.046 | ng/u1 | 0.00 |

Target Compounds Qvalue

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

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