

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072621\  
 Data File : BG049510.D  
 Acq On : 27 Jul 2021 16:35  
 Operator : CG/JU  
 Sample : PB137903BL  
 Misc :  
 ALS Vial : 40 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SBLK903

Quant Time: Jul 27 17:12:15 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SFAM-EPA-BG072421.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jul 26 11:36:47 2021  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 8.057  | 152  | 23129    | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.871 | 136  | 102477   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.701 | 164  | 75620    | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.450 | 188  | 171469   | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.727 | 240  | 161706   | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12          | 25.010 | 264  | 158287   | 20.000 | ng/ul | # 0.00   |

|                               |        |     |        |        |       |      |
|-------------------------------|--------|-----|--------|--------|-------|------|
| System Monitoring Compounds   |        |     |        |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.434  | 96  | 2491   | 4.927  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.857  | 84  | 43406  | 29.893 | ng/ul | 0.00 |
| 7) Phenol-d5                  | 7.205  | 99  | 65189  | 31.856 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.369  | 67  | 36976  | 31.904 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.581  | 132 | 49365  | 33.654 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8         | 8.756  | 113 | 52565  | 34.079 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5           | 9.220  | 128 | 25746  | 32.678 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.948  | 143 | 30926  | 33.497 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.489 | 165 | 52387  | 31.198 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4        | 11.000 | 131 | 72733  | 31.080 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6      | 14.096 | 166 | 181392 | 31.315 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8         | 14.395 | 160 | 218700 | 32.506 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.895 | 143 | 27676  | 29.013 | ng/ul | 0.00 |
| 60) Fluorene-d10              | 15.694 | 176 | 162621 | 32.587 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.811 | 200 | 26392  | 24.839 | ng/ul | 0.00 |
| 73) Anthracene-d10            | 17.550 | 188 | 274096 | 34.394 | ng/ul | 0.00 |
| 81) Pyrene-d10                | 19.835 | 212 | 304348 | 36.808 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12        | 24.781 | 264 | 282323 | 33.935 | ng/ul | 0.01 |

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

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

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