

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021017\
 Data File : BG025833.D
 Acq On : 10 Feb 2017 18:51
 Operator : SJ/MA
 Sample : I1780-05
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AB6

Quant Time: Feb 10 23:07:39 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG021017MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Feb 10 22:04:56 2017
 Response via : Initial Calibration

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.12 | 152 | 88021 | 20.00 | ng/ul | 0.04 |
| 18) Naphthalene-d8 | 10.91 | 136 | 439713 | 20.00 | ng/ul | 0.02 |
| 35) Acenaphthene-d10 | 14.69 | 164 | 292561 | 20.00 | ng/ul | 0.00 |
| 61) Phenanthrene-d10 | 17.42 | 188 | 616356 | 20.00 | ng/ul | 0.00 |
| 77) Chrysene-d12 | 21.67 | 240 | 635817 | 20.00 | ng/ul | 0.00 |
| 85) Perylene-d12 | 24.88 | 264 | 614546 | 20.00 | ng/ul | 0.00 |

System Monitoring Compounds

| | | | | | | |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8 | 3.52 | 96 | 10144 | 5.15 | ng/uL | 0.00 |
| 5) Phenol-d5 | 7.24 | 99 | 43303 | 5.22 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d | 7.42 | 67 | 165571 | 32.71 | ng/ul | 0.02 |
| 9) 2-Chlorophenol-d4 | 7.63 | 132 | 159038 | 25.43 | ng/ul | 0.01 |
| 13) 4-Methylphenol-d8 | 8.78 | 113 | 89336 | 13.02 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5 | 9.25 | 128 | 104713 | 37.09 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4 | 9.97 | 143 | 113807 | 37.14 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3 | 10.51 | 165 | 184930 | 28.38 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4 | 11.02 | 131 | 3161 | 0.38 | ng/ul | 0.01 |
| 43) Dimethylphthalate-d6 | 14.10 | 166 | 712999 | 33.25 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8 | 14.38 | 160 | 857288 | 34.76 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4 | 14.85 | 143 | 22959 | 5.60 | ng/ul | 0.00 |
| 57) Fluorene-d10 | 15.68 | 176 | 636823 | 33.65 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.78 | 200 | 99312 | 31.58 | ng/ul | 0.00 |
| 70) Anthracene-d10 | 17.52 | 188 | 985168 | 35.23 | ng/ul | 0.00 |
| 78) Pyrene-d10 | 19.79 | 212 | 1111068 | 36.64 | ng/ul | 0.00 |
| 89) Benzo(a)pyrene-d12 | 24.67 | 264 | 995891 | 36.48 | ng/ul | 0.00 |

Target Compounds

| | | | | | Qvalue |
|--------------------------------|-------|-----|--------|--------|-----------|
| 20) Nitrobenzene | 9.29 | 77 | 384600 | 57.67 | ng/ul 97 |
| 23) 2-Nitrophenol | 10.00 | 139 | 12500 | 3.79 | ng/ul# 92 |
| 36) 1,2,4,5-Tetrachlorobenzene | 12.90 | 216 | 246119 | 34.49 | ng/ul 98 |
| 72) 1,2,3,4-Tetrachlorobenzene | 13.50 | 216 | 743403 | 110.26 | ng/uL 98 |
| 73) Pentachlorobenzene | 15.01 | 250 | 24124 | 3.41 | ng/uL 97 |

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

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