

Data Path : Z:\HPCHEM1\BNA G\DATA\BG081316\  
 Data File : BG023697.D  
 Acq On : 13 Aug 2016 20:21  
 Operator : UM/SJ  
 Sample : H4388-22MSD  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PR002-SS001-1218-01MSD

Manual Integrations  
 APPROVED

umangi  
 8/16/2016 6:59:22 PM

Quant Time: Aug 15 14:40:52 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080416.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Aug 13 06:38:14 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.11  | 152  | 102720   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.94 | 136  | 469405   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.79 | 164  | 312997   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.55 | 188  | 716988   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.87 | 240  | 751428   | 20.00 | ng/ul | -0.01    |
| 83) Perylene-d12          | 25.27 | 264  | 752279   | 20.00 | ng/ul | -0.01    |

## System Monitoring Compounds

|                                |       |     |         |       |       |       |
|--------------------------------|-------|-----|---------|-------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.48  | 96  | 11455   | 4.74  | ng/uL | 0.00  |
| 5) Phenol-d5                   | 7.27  | 99  | 308795  | 29.84 | ng/ul | 0.00  |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.43  | 67  | 212494m | 32.05 | ng/ul | 0.00  |
| 9) 2-Chlorophenol-d4           | 7.64  | 132 | 219251  | 31.06 | ng/ul | 0.00  |
| 13) 4-Methylphenol-d8          | 8.82  | 113 | 224551  | 27.80 | ng/ul | 0.00  |
| 19) Nitrobenzene-d5            | 9.29  | 128 | 126972m | 34.32 | ng/ul | 0.00  |
| 22) 2-Nitrophenol-d4           | 10.02 | 143 | 144369  | 33.97 | ng/ul | -0.01 |
| 26) 2,4-Dichlorophenol-d3      | 10.57 | 165 | 250339  | 31.23 | ng/ul | 0.00  |
| 29) 4-Chloroaniline-d4         | 11.09 | 131 | 330259  | 35.12 | ng/ul | 0.00  |
| 43) Dimethylphthalate-d6       | 14.18 | 166 | 869486  | 33.93 | ng/ul | 0.00  |
| 46) Acenaphthylene-d8          | 14.48 | 160 | 1010606 | 35.04 | ng/ul | 0.00  |
| 51) 4-Nitrophenol-d4           | 15.00 | 143 | 146240  | 33.43 | ng/ul | 0.00  |
| 57) Fluorene-d10               | 15.78 | 176 | 762074  | 32.18 | ng/ul | 0.00  |
| 62) 4,6-Dinitro-2-methylphenol | 15.91 | 200 | 134172  | 29.17 | ng/ul | -0.01 |
| 70) Anthracene-d10             | 17.65 | 188 | 1125239 | 34.83 | ng/ul | 0.00  |
| 76) Pyrene-d10                 | 19.94 | 212 | 1240677 | 32.62 | ng/ul | 0.00  |
| 87) Benzo(a)pyrene-d12         | 25.03 | 264 | 1242594 | 36.49 | ng/ul | -0.02 |

## Target Compounds

|                                |       |     |         |       |        | Ovalue |
|--------------------------------|-------|-----|---------|-------|--------|--------|
| 6) Phenol                      | 7.30  | 94  | 339418  | 31.50 | ng/ul# | 77     |
| 10) 2-Chlorophenol             | 7.67  | 128 | 235617  | 32.77 | ng/ul# | 84     |
| 15) N-Nitroso-di-n-propylamine | 8.92  | 70  | 270917  | 34.76 | ng/ul# | 85     |
| 33) 4-Chloro-3-methylphenol    | 12.24 | 107 | 346414  | 33.30 | ng/ul# | 81     |
| 44) Dimethylphthalate          | 14.22 | 163 | 378318  | 14.91 | ng/ul  | 98     |
| 49) Acenaphthene               | 14.85 | 153 | 721309  | 35.09 | ng/ul  | 96     |
| 52) 4-Nitrophenol              | 15.02 | 109 | 141834  | 32.37 | ng/ul# | 75     |
| 54) 2,4-Dinitrotoluene         | 15.15 | 165 | 286050  | 34.31 | ng/ul# | 68     |
| 68) Pentachlorophenol          | 17.20 | 266 | 164027  | 33.94 | ng/ul  | 93     |
| 77) Pyrene                     | 19.97 | 202 | 1548515 | 33.34 | ng/ul# | 91     |

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

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