

Data Path : Z:\HPCHEM1\BNA G\DATA\BG111315\  
 Data File : BG019595.D  
 Acq On : 13 Nov 2015 17:06  
 Operator : UM/NP  
 Sample : G4342-08MSD  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 D9KP7MSD

Manual Integrations  
 APPROVED

mmdadoda  
 11/16/2015 3:26:08 PM

Quant Time: Nov 14 00:29:36 2015  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102815.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Nov 14 00:06:36 2015  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.16  | 152  | 23785    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.99 | 136  | 96999    | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.79 | 164  | 75058    | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.53 | 188  | 221846   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.80 | 240  | 290382   | 20.00 | ng/ul | -0.01    |
| 86) Perylene-d12          | 25.14 | 264  | 275231   | 20.00 | ng/ul | -0.02    |

## System Monitoring Compounds

|                                |       |     |        |       |       |       |
|--------------------------------|-------|-----|--------|-------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.46  | 96  | 2626   | 5.29  | ng/uL | 0.00  |
| 5) Phenol-d5                   | 7.30  | 99  | 55190  | 25.28 | ng/ul | 0.00  |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.47  | 67  | 37597  | 27.16 | ng/ul | 0.00  |
| 9) 2-Chlorophenol-d4           | 7.69  | 132 | 33982  | 24.93 | ng/ul | 0.00  |
| 13) 4-Methylphenol-d8          | 8.87  | 113 | 44709  | 25.71 | ng/ul | 0.00  |
| 19) Nitrobenzene-d5            | 9.34  | 128 | 18600  | 26.09 | ng/ul | 0.00  |
| 22) 2-Nitrophenol-d4           | 10.06 | 143 | 21941  | 25.93 | ng/ul | 0.00  |
| 26) 2,4-Dichlorophenol-d3      | 10.61 | 165 | 41311  | 22.50 | ng/ul | 0.00  |
| 29) 4-Chloroaniline-d4         | 11.12 | 131 | 50608  | 27.25 | ng/ul | 0.00  |
| 44) Dimethylphthalate-d6       | 14.19 | 166 | 186991 | 30.22 | ng/ul | 0.00  |
| 47) Acenaphthylene-d8          | 14.49 | 160 | 176474 | 25.75 | ng/ul | 0.00  |
| 52) 4-Nitrophenol-d4           | 14.98 | 143 | 28898  | 29.39 | ng/ul | 0.00  |
| 58) Fluorene-d10               | 15.78 | 176 | 146933 | 25.66 | ng/ul | 0.00  |
| 63) 4,6-Dinitro-2-methylphenol | 15.89 | 200 | 35529  | 24.74 | ng/ul | 0.00  |
| 71) Anthracene-d10             | 17.63 | 188 | 277422 | 27.43 | ng/ul | 0.00  |
| 79) Pyrene-d10                 | 19.91 | 212 | 348415 | 27.42 | ng/ul | 0.00  |
| 90) Benzo(a)pyrene-d12         | 24.91 | 264 | 380261 | 27.53 | ng/ul | -0.02 |

## Target Compounds

|                                |       |     |        |       | Ovalue    |
|--------------------------------|-------|-----|--------|-------|-----------|
| 6) Phenol                      | 7.33  | 94  | 66968  | 28.79 | ng/ul 90  |
| 10) 2-Chlorophenol             | 7.72  | 128 | 37722  | 26.89 | ng/ul 98  |
| 15) N-Nitroso-di-n-propylamine | 8.97  | 70  | 56866  | 29.72 | ng/ul# 86 |
| 33) 4-Chloro-3-methylphenol    | 12.25 | 107 | 69101  | 29.34 | ng/ul 94  |
| 45) Dimethylphthalate          | 14.24 | 163 | 86560  | 13.68 | ng/ul 97  |
| 50) Acenaphthene               | 14.86 | 153 | 132795 | 28.01 | ng/ul 94  |
| 53) 4-Nitrophenol              | 14.99 | 109 | 40094m | 26.15 | ng/ul     |
| 55) 2,4-Dinitrotoluene         | 15.15 | 165 | 57790  | 28.56 | ng/ul# 78 |
| 69) Pentachlorophenol          | 17.18 | 266 | 44181  | 26.75 | ng/ul 97  |
| 80) Pyrene                     | 19.94 | 202 | 478194 | 29.35 | ng/ul# 90 |
| 84) Bis(2-ethylhexyl)phthalate | 21.67 | 149 | 11940  | 1.59  | ng/ul# 97 |

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

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