

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG121016\  
 Data File : BG025099.D  
 Acq On : 11 Dec 2016 11:37  
 Operator : UM/SJ  
 Sample : H5905-21  
 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 DA820

Manual Integrations  
 APPROVED

Sohil  
 12/12/2016 4:30:08 PM

Quant Time: Dec 12 17:29:51 2016  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Dec 12 06:35:19 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.23  | 152  | 102452   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.07 | 136  | 467434   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.86 | 164  | 333667   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.60 | 188  | 699907   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.89 | 240  | 876115   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.31 | 264  | 915881   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.59  | 96  | 12452  | 6.28  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.38  | 99  | 212123 | 24.00 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.55  | 67  | 136774 | 25.01 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.76  | 132 | 149447 | 24.18 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.94  | 113 | 192300 | 26.03 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.42  | 128 | 92252  | 27.87 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.14 | 143 | 101223 | 24.53 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.68 | 165 | 186906 | 23.55 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.20 | 131 | 518539 | 66.49 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 14.26 | 166 | 561605 | 24.47 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.56 | 160 | 660351 | 26.27 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 15.07 | 143 | 104613 | 26.52 | ng/ul | 0.02 |
| 57) Fluorene-d10               | 15.85 | 176 | 690663 | 31.97 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.97 | 200 | 119205 | 27.55 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.70 | 188 | 785789 | 26.60 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.97 | 212 | 931577 | 25.82 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 25.06 | 264 | 974466 | 26.26 | ng/ul | 0.00 |

## Target Compounds

|                         |       |     |         |        |        | Qvalue |
|-------------------------|-------|-----|---------|--------|--------|--------|
| 11) 2-Methylphenol      | 8.67  | 108 | 19022   | 2.85   | ng/ul  | 87     |
| 14) Acetophenone        | 9.03  | 105 | 103796  | 9.23   | ng/ul# | 30     |
| 16) 4-Methylphenol      | 9.00  | 108 | 54639   | 7.35   | ng/ul  | 94     |
| 24) 2,4-Dimethylphenol  | 10.21 | 107 | 190111m | 19.86  | ng/ul  |        |
| 28) Naphthalene         | 11.12 | 128 | 1237769 | 56.96  | ng/ul  | 96     |
| 34) 2-Methylnaphthalene | 12.71 | 142 | 2146828 | 125.20 | ng/ul  | 93     |
| 40) 1,1'-Biphenyl       | 13.70 | 154 | 478311  | 23.64  | ng/ul  | 98     |
| 44) Dimethylphthalate   | 14.30 | 163 | 405328  | 17.97  | ng/ul# | 98     |
| 49) Acenaphthene        | 14.93 | 153 | 117407  | 7.01   | ng/ul# | 89     |
| 53) Dibenzofuran        | 15.26 | 168 | 307801  | 11.62  | ng/ul  | 94     |
| 58) Fluorene            | 15.90 | 166 | 324870  | 15.07  | ng/ul# | 98     |
| 69) Phenanthrene        | 17.65 | 178 | 140921  | 4.22   | ng/ul# | 91     |
| 71) Anthracene          | 17.74 | 178 | 37200   | 1.08   | ng/ul# | 90     |

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

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