

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122815\  
 Data File : BG020579.D  
 Acq On : 29 Dec 2015 1:04  
 Operator : UM/SJ  
 Sample : G4847-02DL 10X  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleID :  
 D9N68DL

Manual Integrations  
 APPROVED

UMANGI  
 12/29/2015 4:55:08 PM

Quant Time: Dec 29 06:04:20 2015  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG122415.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Dec 29 05:04:12 2015  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.07  | 152  | 12278    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.89 | 136  | 54056    | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.70 | 164  | 43389    | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.45 | 188  | 135861   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.73 | 240  | 186252   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 24.98 | 264  | 175803   | 20.00 | ng/ul | 0.00     |

| System Monitoring Compounds    | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|--------------------------------|-------|------|----------|------|-------|----------|
| 3) 1,4-Dioxane-d8              | 0.00  | 96   | 0        | 0.00 | ng/uL |          |
| 5) Phenol-d5                   | 0.00  | 99   | 0d       | 0.00 | ng/ul |          |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.39  | 67   | 1492     | 2.56 | ng/ul | 0.00     |
| 9) 2-Chlorophenol-d4           | 7.61  | 132  | 1779     | 2.35 | ng/ul | 0.01     |
| 13) 4-Methylphenol-d8          | 8.78  | 113  | 928m     | 1.12 | ng/ul | 0.00     |
| 19) Nitrobenzene-d5            | 9.24  | 128  | 1133     | 2.67 | ng/ul | 0.00     |
| 22) 2-Nitrophenol-d4           | 9.97  | 143  | 1253     | 2.60 | ng/ul | 0.00     |
| 26) 2,4-Dichlorophenol-d3      | 10.52 | 165  | 2220     | 2.38 | ng/ul | 0.01     |
| 29) 4-Chloroaniline-d4         | 0.00  | 131  | 0        | 0.00 | ng/ul |          |
| 44) Dimethylphthalate-d6       | 14.10 | 166  | 12785    | 3.57 | ng/ul | 0.00     |
| 47) Acenaphthylene-d8          | 14.40 | 160  | 13781    | 3.33 | ng/ul | 0.00     |
| 52) 4-Nitrophenol-d4           | 0.00  | 143  | 0        | 0.00 | ng/ul |          |
| 58) Fluorene-d10               | 15.69 | 176  | 12040    | 3.72 | ng/ul | 0.00     |
| 63) 4,6-Dinitro-2-methylphenol | 0.00  | 200  | 0d       | 0.00 | ng/ul |          |
| 71) Anthracene-d10             | 17.55 | 188  | 23308    | 3.66 | ng/ul | 0.00     |
| 79) Pyrene-d10                 | 19.84 | 212  | 28242    | 3.50 | ng/ul | 0.00     |
| 90) Benzo(a)pyrene-d12         | 24.75 | 264  | 28106    | 3.16 | ng/ul | -0.01    |

| Target Compounds        | R.T.  | QIon | Response | Conc  | Units  | Ovalue |
|-------------------------|-------|------|----------|-------|--------|--------|
| 28) Naphthalene         | 10.93 | 128  | 117069   | 40.74 | ng/ul  | 99     |
| 34) 2-Methylnaphthalene | 12.54 | 142  | 12686    | 5.97  | ng/ul  | 93     |
| 50) Acenaphthene        | 14.77 | 153  | 12606    | 4.28  | ng/ul  | 96     |
| 54) Dibenzofuran        | 15.10 | 168  | 9560     | 2.18  | ng/ul  | 91     |
| 59) Fluorene            | 15.75 | 166  | 8039     | 2.26  | ng/ul# | 95     |
| 70) Phenanthrene        | 17.49 | 178  | 17499    | 2.31  | ng/ul  | 98     |

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

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