

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122215\  
 Data File : BG020461.D  
 Acq On : 24 Dec 2015 2:42  
 Operator : UM/SJ  
 Sample : G4849-09DL 20X  
 Misc :  
 ALS Vial : 51 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 D9N73MEDL

Manual Integrations  
 APPROVED

Sohil  
 12/24/2015 5:31:31 PM

Quant Time: Dec 24 06:45:22 2015  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG121415.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Dec 24 02:43:25 2015  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.08  | 152  | 18452    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.90 | 136  | 81238    | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.72 | 164  | 59287    | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.46 | 188  | 144193   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.73 | 240  | 177586   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 25.00 | 264  | 163447   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |       |      |       |      |
|--------------------------------|-------|-----|-------|------|-------|------|
| 3) 1,4-Dioxane-d8              | 0.00  | 96  | 0     | 0.00 | ng/uL |      |
| 5) Phenol-d5                   | 7.24  | 99  | 2364  | 1.45 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.40  | 67  | 1418  | 1.46 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.61  | 132 | 1929  | 1.63 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.80  | 113 | 2003  | 1.44 | ng/ul | 0.01 |
| 19) Nitrobenzene-d5            | 9.25  | 128 | 1048  | 1.65 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.98  | 143 | 1085  | 1.54 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.52 | 165 | 2238  | 1.57 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.03 | 131 | 2636  | 1.64 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 14.11 | 166 | 10011 | 2.09 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.41 | 160 | 11943 | 2.14 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.92 | 143 | 54    | 0.06 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.71 | 176 | 9284  | 2.13 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.83 | 200 | 347   | 0.41 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.56 | 188 | 15099 | 2.27 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.84 | 212 | 17135 | 2.07 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 24.78 | 264 | 17660 | 2.17 | ng/ul | 0.00 |

## Target Compounds

|                          |       |     |        |       |        | Ovalue |
|--------------------------|-------|-----|--------|-------|--------|--------|
| 28) Naphthalene          | 10.95 | 128 | 95498  | 22.51 | ng/ul  | 99     |
| 34) 2-Methylnaphthalene  | 12.55 | 142 | 34910  | 10.77 | ng/ul  | 95     |
| 41) 1,1'-Biphenyl        | 13.55 | 154 | 15065  | 3.37  | ng/ul# | 93     |
| 50) Acenaphthene         | 14.78 | 153 | 70110  | 17.58 | ng/ul  | 99     |
| 54) Dibenzofuran         | 15.11 | 168 | 73157  | 12.33 | ng/ul  | 97     |
| 59) Fluorene             | 15.76 | 166 | 70544  | 14.81 | ng/ul  | 96     |
| 70) Phenanthrene         | 17.50 | 178 | 336567 | 42.67 | ng/ul  | 100    |
| 72) Anthracene           | 17.59 | 178 | 30873  | 3.86  | ng/ul  | 99     |
| 75) Carbazole            | 17.86 | 167 | 17211  | 2.56  | ng/ul  | 99     |
| 77) Fluoranthene         | 19.51 | 202 | 208586 | 22.63 | ng/ul  | 99     |
| 80) Pyrene               | 19.87 | 202 | 137908 | 12.83 | ng/ul  | 98     |
| 83) Benzo(a)anthracene   | 21.72 | 228 | 33958  | 3.28  | ng/ul  | 99     |
| 85) Chrysene             | 21.78 | 228 | 27954  | 2.86  | ng/ul  | 98     |
| 88) Benzo(b)fluoranthene | 23.95 | 252 | 13920m | 1.41  | ng/ul  |        |

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

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