

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG033115\  
 Data File : BG016207.D  
 Acq On : 31 Mar 2015 19:29  
 Operator : TP/IZ  
 Sample : G1647-15  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 C0AD4

Quant Time: Apr 01 07:01:21 2015  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM01.2-EPA-BG033115.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Apr 01 04:39:57 2015  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.77  | 152  | 92517    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.55 | 136  | 409177   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.39 | 164  | 233071   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.14 | 188  | 462282   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.32 | 240  | 425945   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.60 | 264  | 418205   | 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                   | 6.94  | 99  | 245948 | 29.00 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.10  | 67  | 149827 | 30.55 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.30  | 132 | 192158 | 28.85 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.47  | 113 | 178219 | 26.97 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.92  | 128 | 108897 | 31.75 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.64  | 143 | 116855 | 30.49 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.18 | 165 | 173201 | 27.79 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.69 | 131 | 216833 | 25.83 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.81 | 166 | 496239 | 32.21 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.09 | 160 | 669023 | 30.87 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.60 | 143 | 109636 | 28.23 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.39 | 176 | 459219 | 30.44 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.51 | 200 | 78739  | 24.93 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.24 | 188 | 632847 | 29.79 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.53 | 212 | 629524 | 32.53 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.46 | 264 | 533658 | 28.28 | ng/ul | 0.01 |

## Target Compounds

|                            |       |     |        |      |        | Qvalue |
|----------------------------|-------|-----|--------|------|--------|--------|
| 14) Acetophenone           | 8.59  | 105 | 11536  | 1.16 | ng/ul  | 94     |
| 44) Dimethylphthalate      | 13.85 | 163 | 34395  | 1.99 | ng/ul  | 98     |
| 69) Phenanthrene           | 17.18 | 178 | 33205  | 1.27 | ng/ul  | 98     |
| 74) Fluoranthene           | 19.19 | 202 | 65574  | 2.25 | ng/ul  | 99     |
| 77) Pyrene                 | 19.56 | 202 | 68348  | 2.59 | ng/ul  | 97     |
| 80) Benzo(a)anthracene     | 21.30 | 228 | 35756  | 1.45 | ng/ul  | 97     |
| 82) Chrysene               | 21.35 | 228 | 33401m | 1.46 | ng/ul  |        |
| 85) Benzo(b)fluoranthene   | 22.91 | 252 | 65902  | 2.64 | ng/ul# | 81     |
| 88) Benzo(a)pyrene         | 23.50 | 252 | 46641  | 1.96 | ng/ul# | 92     |
| 89) Indeno(1,2,3-cd)pyrene | 25.93 | 276 | 34559  | 1.24 | ng/ul  | 94     |
| 91) Benzo(g,h,i)perylene   | 26.65 | 276 | 30074  | 1.32 | ng/ul  | 96     |

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

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