

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG021016\  
 Data File : BG020836.D  
 Acq On : 10 Feb 2016 12:08  
 Operator : SJ/UM  
 Sample : H1390-05  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 C04P9

Manual Integrations  
 APPROVED

UMANGI  
 2/11/2016 10:48:48 AM

Quant Time: Feb 11 03:46:36 2016  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG020816.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Feb 11 03:24:50 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.27  | 152  | 21687    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.10 | 136  | 112119   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.89 | 164  | 63262    | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.62 | 188  | 122117   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.91 | 240  | 107995   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 25.30 | 264  | 102443   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |       |       |       |      |
|--------------------------------|-------|-----|-------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.61  | 96  | 915   | 2.05  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.40  | 99  | 25930 | 11.80 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.58  | 67  | 14485 | 12.36 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.80  | 132 | 21880 | 12.73 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.97  | 113 | 22260 | 11.66 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.44  | 128 | 11362 | 12.63 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.17 | 143 | 11592 | 12.95 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.70 | 165 | 21328 | 12.38 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.22 | 131 | 30261 | 13.20 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 14.28 | 166 | 69714 | 12.81 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.58 | 160 | 92160 | 13.86 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 15.05 | 143 | 10504 | 9.64  | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.87 | 176 | 61439 | 13.26 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.97 | 200 | 4663  | 7.88  | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.72 | 188 | 85264 | 14.13 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.99 | 212 | 82076 | 14.78 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 25.06 | 264 | 79467 | 14.00 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |      |        | Qvalue |
|--------------------------------|-------|-----|--------|------|--------|--------|
| 14) Acetophenone               | 9.10  | 105 | 6200   | 2.15 | ng/ul  | 93     |
| 45) Dimethylphthalate          | 14.32 | 163 | 28787  | 5.08 | ng/ul  | 98     |
| 70) Phenanthrene               | 17.66 | 178 | 18127  | 2.36 | ng/ul  | 98     |
| 76) Di-n-butylphthalate        | 18.56 | 149 | 18934  | 2.16 | ng/ul  | 99     |
| 77) Fluoranthene               | 19.65 | 202 | 26959  | 3.39 | ng/ul  | 97     |
| 80) Pyrene                     | 20.01 | 202 | 26403  | 3.39 | ng/ul  | 98     |
| 83) Benzo(a)anthracene         | 21.88 | 228 | 11096  | 1.58 | ng/ul# | 93     |
| 84) Bis(2-ethylhexyl)phthalate | 21.77 | 149 | 9029   | 1.65 | ng/ul# | 99     |
| 85) Chrysene                   | 21.95 | 228 | 13603m | 2.10 | ng/ul  |        |
| 88) Benzo(b)fluoranthene       | 24.22 | 252 | 17049  | 2.44 | ng/ul# | 80     |
| 91) Benzo(a)pyrene             | 25.13 | 252 | 11254  | 1.70 | ng/ul# | 78     |
| 92) Indeno(1,2,3-cd)pyrene     | 29.17 | 276 | 8461   | 1.11 | ng/ul# | 86     |

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

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