

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG011416\  
 Data File : BG020738.D  
 Acq On : 15 Jan 2016 00:27  
 Operator : SJ/IZ  
 Sample : H1101-11  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BC9F2

Manual Integrations  
 APPROVED

mohammad  
 1/15/2016 2:44:01 PM

Quant Time: Jan 15 01:30:11 2016  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG011216.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jan 14 23:48:02 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.04  | 152  | 19765    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.86 | 136  | 83531    | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.68 | 164  | 58212    | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.43 | 188  | 147548   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.70 | 240  | 174542   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 24.95 | 264  | 160109   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.41  | 96  | 884    | 2.19  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.20  | 99  | 21464  | 12.67 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.36  | 67  | 14585  | 14.32 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.57  | 132 | 17870  | 13.45 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.76  | 113 | 3840m  | 2.69  | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.22  | 128 | 9938   | 15.08 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.94  | 143 | 11307  | 14.77 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.48 | 165 | 18263  | 12.53 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.00 | 131 | 15802  | 10.95 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 14.08 | 166 | 76756  | 15.11 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.38 | 160 | 88518  | 15.52 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.90 | 143 | 7770m  | 11.39 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.68 | 176 | 71182  | 15.72 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.80 | 200 | 9307m  | 11.60 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.53 | 188 | 114294 | 15.80 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.81 | 212 | 136587 | 16.45 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 24.73 | 264 | 137532 | 16.11 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |        |      |        | Qvalue |
|----------------------------|-------|-----|--------|------|--------|--------|
| 45) Dimethylphthalate      | 14.12 | 163 | 41663  | 7.63 | ng/ul# | 98     |
| 70) Phenanthrene           | 17.47 | 178 | 42501  | 4.84 | ng/ul  | 99     |
| 72) Anthracene             | 17.56 | 178 | 9777   | 1.10 | ng/ul  | 95     |
| 77) Fluoranthene           | 19.48 | 202 | 57782  | 5.41 | ng/ul  | 99     |
| 80) Pyrene                 | 19.84 | 202 | 81282  | 7.34 | ng/ul  | 99     |
| 83) Benzo(a)anthracene     | 21.68 | 228 | 40507  | 3.72 | ng/ul  | 97     |
| 85) Chrysene               | 21.74 | 228 | 40462  | 3.93 | ng/ul  | 97     |
| 88) Benzo(b)fluoranthene   | 23.92 | 252 | 32264  | 3.13 | ng/ul# | 95     |
| 89) Benzo(k)fluoranthene   | 23.97 | 252 | 12943m | 1.27 | ng/ul  |        |
| 91) Benzo(a)pyrene         | 24.79 | 252 | 30676  | 3.09 | ng/ul  | 98     |
| 92) Indeno(1,2,3-cd)pyrene | 28.65 | 276 | 15602m | 1.33 | ng/ul  |        |
| 94) Benzo(g,h,i)perylene   | 29.84 | 276 | 15090m | 1.56 | ng/ul  |        |

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

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