

Data Path : Z:\HPCHEM1\BNA G\DATA\BG012417\  
 Data File : BG025621.D  
 Acq On : 24 Jan 2017 23:47  
 Operator : UM/SJ  
 Sample : I1316-11  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BDP21

Manual Integrations  
 APPROVED

mohammad  
 1/25/2017 3:47:56 PM

Quant Time: Jan 25 03:06:02 2017  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG011817.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jan 25 02:56:13 2017  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.18  | 152  | 64841m   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.00 | 136  | 314529   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.81 | 164  | 278541   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.55 | 188  | 625863m  | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.85 | 240  | 756006   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.23 | 264  | 779493   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.51  | 96  | 5826    | 4.50  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.34  | 99  | 139267  | 25.06 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.49  | 67  | 101722  | 28.34 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.71  | 132 | 106121m | 27.07 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.90  | 113 | 100612  | 21.30 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.36  | 128 | 61316   | 27.50 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.09 | 143 | 74540   | 28.02 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.64 | 165 | 118272  | 22.54 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.15 | 131 | 143193  | 24.41 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 14.20 | 166 | 541089  | 27.51 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.50 | 160 | 604759  | 28.46 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 15.06 | 143 | 62111m  | 23.04 | ng/ul | 0.01 |
| 57) Fluorene-d10               | 15.80 | 176 | 493099  | 26.59 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.94 | 200 | 98673   | 25.56 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.65 | 188 | 772847  | 29.00 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.93 | 212 | 932680  | 29.65 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 25.00 | 264 | 951046  | 29.03 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |         |      |        | Ovalue |
|----------------------------|-------|-----|---------|------|--------|--------|
| 6) Phenol                  | 7.36  | 94  | 20832m  | 3.57 | ng/ul  |        |
| 44) Dimethylphthalate      | 14.25 | 163 | 134773  | 6.97 | ng/ul  | 96     |
| 69) Phenanthrene           | 17.60 | 178 | 106332  | 3.50 | ng/ul  | 95     |
| 74) Fluoranthene           | 19.60 | 202 | 324115  | 9.03 | ng/ul  | 99     |
| 77) Pyrene                 | 19.96 | 202 | 278807  | 7.50 | ng/ul  | 97     |
| 80) Benzo(a)anthracene     | 21.82 | 228 | 187386  | 4.69 | ng/ul  | 96     |
| 82) Chrysene               | 21.89 | 228 | 159775  | 4.25 | ng/ul  | 97     |
| 85) Benzo(b)fluoranthene   | 24.16 | 252 | 205721  | 5.10 | ng/ul# | 98     |
| 86) Benzo(k)fluoranthene   | 24.21 | 252 | 61839   | 1.63 | ng/ul# | 91     |
| 88) Benzo(a)pyrene         | 25.07 | 252 | 143915  | 3.77 | ng/ul  | 100    |
| 89) Indeno(1,2,3-cd)pyrene | 29.16 | 276 | 100004m | 2.07 | ng/ul  |        |
| 91) Benzo(g,h,i)perylene   | 30.40 | 276 | 91839m  | 2.31 | ng/ul  |        |

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

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