

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121716\  
 Data File : BG025288.D  
 Acq On : 18 Dec 2016 00:05  
 Operator : UM/SJ  
 Sample : H5995-02DL 10X  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 DA863DL

Manual Integrations  
 APPROVED

umangi  
 12/20/2016 10:54:44 AM

Quant Time: Dec 18 04:51:56 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun Dec 18 03:57:56 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.23  | 152  | 64129    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.06 | 136  | 263088   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.85 | 164  | 198005   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.60 | 188  | 442876   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.90 | 240  | 563894   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.32 | 264  | 572241   | 20.00 | ng/ul | -0.01    |

## System Monitoring Compounds

|                                |       |     |       |      |       |       |
|--------------------------------|-------|-----|-------|------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.59  | 96  | 303   | 0.24 | ng/uL | 0.00  |
| 5) Phenol-d5                   | 7.41  | 99  | 7310  | 1.32 | ng/ul | 0.01  |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.55  | 67  | 4769  | 1.39 | ng/ul | 0.00  |
| 9) 2-Chlorophenol-d4           | 7.78  | 132 | 5962  | 1.54 | ng/ul | 0.01  |
| 13) 4-Methylphenol-d8          | 8.96  | 113 | 6165  | 1.33 | ng/ul | 0.01  |
| 19) Nitrobenzene-d5            | 9.42  | 128 | 2866  | 1.54 | ng/ul | 0.00  |
| 22) 2-Nitrophenol-d4           | 10.15 | 143 | 3603  | 1.55 | ng/ul | 0.00  |
| 26) 2,4-Dichlorophenol-d3      | 10.69 | 165 | 6579  | 1.47 | ng/ul | 0.00  |
| 29) 4-Chloroaniline-d4         | 11.23 | 131 | 771   | 0.18 | ng/ul | 0.02  |
| 43) Dimethylphthalate-d6       | 14.25 | 166 | 23067 | 1.69 | ng/ul | 0.00  |
| 46) Acenaphthylene-d8          | 14.55 | 160 | 25823 | 1.73 | ng/ul | 0.00  |
| 51) 4-Nitrophenol-d4           | 15.07 | 143 | 104   | 0.04 | ng/ul | -0.02 |
| 57) Fluorene-d10               | 15.85 | 176 | 24124 | 1.88 | ng/ul | 0.00  |
| 62) 4,6-Dinitro-2-methylphenol | 15.99 | 200 | 2143  | 0.78 | ng/ul | 0.00  |
| 70) Anthracene-d10             | 17.70 | 188 | 32371 | 1.73 | ng/ul | 0.00  |
| 76) Pyrene-d10                 | 19.98 | 212 | 43455 | 1.87 | ng/ul | 0.00  |
| 87) Benzo(a)pyrene-d12         | 25.07 | 264 | 38794 | 1.67 | ng/ul | -0.02 |

## Target Compounds

|                                |       |     |        |       |        | Ovalue |
|--------------------------------|-------|-----|--------|-------|--------|--------|
| 69) Phenanthrene               | 17.64 | 178 | 75101  | 3.55  | ng/ul# | 91     |
| 74) Fluoranthene               | 19.64 | 202 | 100781 | 4.01  | ng/ul  | 97     |
| 77) Pyrene                     | 20.00 | 202 | 85841  | 3.17  | ng/ul  | 99     |
| 80) Benzo(a)anthracene         | 21.87 | 228 | 55430  | 1.89  | ng/ul  | 99     |
| 81) Bis(2-ethylhexyl)phthalate | 21.72 | 149 | 778649 | 53.30 | ng/ul  | 99     |
| 82) Chrysene                   | 21.95 | 228 | 50199m | 1.83  | ng/ul  |        |
| 85) Benzo(b)fluoranthene       | 24.22 | 252 | 71915  | 2.53  | ng/ul# | 88     |
| 88) Benzo(a)pyrene             | 25.15 | 252 | 48687  | 1.78  | ng/ul# | 83     |
| 89) Indeno(1,2,3-cd)pyrene     | 29.27 | 276 | 34206  | 1.01  | ng/ul# | 88     |

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

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121716\  
Data File : BG025288.D  
Acq On : 18 Dec 2016 00:05  
Operator : UM/SJ  
Sample : H5995-02DL 10X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
DA863DL

Manual Integrations  
APPROVED

umangi  
12/20/2016 10:54:44 AM

Quant Time: Dec 18 04:51:56 2016  
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sun Dec 18 03:57:56 2016  
Response via : Initial Calibration

