

Data Path : Z:\HPCHEM1\BNA G\DATA\BG081316\  
 Data File : BG023688.D  
 Acq On : 13 Aug 2016 14:39  
 Operator : UM/SJ  
 Sample : H4388-13  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 PR002-SS003-0206-02

Manual Integrations  
 APPROVED

umangi  
 8/16/2016 6:59:08 PM

Quant Time: Aug 15 13:46:03 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080416.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Aug 15 11:28:11 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.11  | 152  | 95806    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.95 | 136  | 453316   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.78 | 164  | 312577   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.55 | 188  | 734731m  | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.87 | 240  | 777822m  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.28 | 264  | 782530   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.48  | 96  | 12015   | 5.33  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.26  | 99  | 283483  | 29.37 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.42  | 67  | 191776  | 31.01 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.64  | 132 | 205099  | 31.15 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.83  | 113 | 212715  | 28.23 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.29  | 128 | 118898  | 33.28 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.02 | 143 | 136913  | 33.36 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.57 | 165 | 232731  | 30.07 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.08 | 131 | 248273  | 27.34 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 14.18 | 166 | 822749  | 32.15 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.48 | 160 | 971839  | 33.74 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 15.01 | 143 | 128947  | 29.51 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.78 | 176 | 773335  | 32.70 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.91 | 200 | 135174m | 28.68 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.65 | 188 | 1118828 | 33.80 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.95 | 212 | 1236116 | 31.40 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 25.04 | 264 | 1215964 | 34.33 | ng/ul | 0.00 |

## Target Compounds

|                          |       |     |         |       |        | Ovalue |
|--------------------------|-------|-----|---------|-------|--------|--------|
| 6) Phenol                | 7.29  | 94  | 15750   | 1.57  | ng/ul# | 75     |
| 44) Dimethylphthalate    | 14.23 | 163 | 359174  | 14.17 | ng/ul  | 99     |
| 69) Phenanthrene         | 17.60 | 178 | 51142   | 1.34  | ng/ul  | 97     |
| 74) Fluoranthene         | 19.61 | 202 | 111358m | 2.86  | ng/ul  |        |
| 77) Pyrene               | 19.97 | 202 | 112359m | 2.34  | ng/ul  |        |
| 80) Benzo(a)anthracene   | 21.85 | 228 | 45274m  | 1.03  | ng/ul  |        |
| 82) Chrysene             | 21.92 | 228 | 53128m  | 1.33  | ng/ul  |        |
| 85) Benzo(b)fluoranthene | 24.20 | 252 | 65077m  | 1.46  | ng/ul  |        |
| 88) Benzo(a)pyrene       | 25.12 | 252 | 48637m  | 1.13  | ng/ul  |        |

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

Data Path : Z:\HPCHEM1\BNA G\DATA\BG081316\  
Data File : BG023688.D  
Acq On : 13 Aug 2016 14:39  
Operator : UM/SJ  
Sample : H4388-13  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
PR002-SS003-0206-02

Manual Integrations  
APPROVED

umangi  
8/16/2016 6:59:08 PM

Quant Time: Aug 15 13:46:03 2016  
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080416.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Aug 15 11:28:11 2016  
Response via : Initial Calibration

