

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG082917\  
 Data File : BG028541.D  
 Acq On : 29 Aug 2017 19:23  
 Operator : SJ/JU  
 Sample : I4960-01  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 5-POST1

Manual Integrations  
 APPROVED

Sohil  
 8/30/2017 6:48:43 PM

Quant Time: Aug 30 15:10:59 2017  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080217.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 30 01:19:11 2017  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.35  | 152  | 56253    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.35 | 136  | 220305m  | 20.00 | ng/ul | 0.18     |
| 35) Acenaphthene-d10      | 14.96 | 164  | 139553   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.71 | 188  | 309752   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 22.05 | 240  | 365942   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.56 | 264  | 362922   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.83  | 96  | 2608m  | 2.16  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.50  | 99  | 89234  | 14.44 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.66  | 67  | 71120  | 18.18 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.89  | 132 | 74268  | 19.06 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 9.05  | 113 | 82025  | 15.89 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.53  | 128 | 42697  | 24.79 | ng/ul | 0.02 |
| 22) 2-Nitrophenol-d4           | 10.29 | 143 | 54100m | 28.64 | ng/ul | 0.05 |
| 26) 2,4-Dichlorophenol-d3      | 10.86 | 165 | 87921  | 22.50 | ng/ul | 0.07 |
| 29) 4-Chloroaniline-d4         | 11.38 | 131 | 109433 | 25.16 | ng/ul | 0.08 |
| 43) Dimethylphthalate-d6       | 14.35 | 166 | 296384 | 23.88 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.65 | 160 | 338480 | 24.19 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 15.16 | 143 | 30021  | 15.21 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.95 | 176 | 245702 | 22.07 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 16.07 | 200 | 32038  | 15.92 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.80 | 188 | 358110 | 24.11 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 20.08 | 212 | 420518 | 22.41 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 25.32 | 264 | 379822 | 22.36 | ng/ul | 0.01 |

## Target Compounds

|                                |       |     |          |         | Qvalue    |
|--------------------------------|-------|-----|----------|---------|-----------|
| 6) Phenol                      | 7.53  | 94  | 9570     | 1.559   | ng/ul 90  |
| 28) Naphthalene                | 11.37 | 128 | 133684   | 12.252  | ng/ul 100 |
| 34) 2-Methylnaphthalene        | 12.81 | 142 | 9587     | 1.094   | ng/ul 99  |
| 36) 1,2,4,5-Tetrachlorobenzene | 13.19 | 216 | 2508986m | 557.778 | ng/ul     |
| 39) 2,4,5-Trichlorophenol      | 13.49 | 196 | 5769     | 1.888   | ng/ul# 92 |
| 44) Dimethylphthalate          | 14.40 | 163 | 105355   | 9.219   | ng/ul# 98 |
| 49) Acenaphthene               | 15.02 | 153 | 14041    | 1.556   | ng/ul 91  |
| 53) Dibenzofuran               | 15.35 | 168 | 20300    | 1.543   | ng/ul 89  |
| 58) Fluorene                   | 16.00 | 166 | 27426    | 2.355   | ng/ul 98  |
| 69) Phenanthrene               | 17.75 | 178 | 284596   | 18.257  | ng/ul 98  |
| 71) Anthracene                 | 17.84 | 178 | 32179m   | 2.020   | ng/ul     |
| 72) Carbazole                  | 18.11 | 167 | 18685    | 1.349   | ng/ul# 76 |
| 74) Fluoranthene               | 19.75 | 202 | 425453   | 22.458  | ng/ul 100 |
| 77) Pyrene                     | 20.11 | 202 | 345841   | 15.841  | ng/ul# 96 |
| 80) Benzo(a)anthracene         | 22.02 | 228 | 77058    | 3.644   | ng/ul# 82 |
| 81) Bis(2-ethylhexyl)phthalate | 21.86 | 149 | 62282    | 5.354   | ng/ul# 93 |
| 82) Chrysene                   | 22.09 | 228 | 146256   | 7.681   | ng/ul# 91 |
| 85) Benzo(b)fluoranthene       | 24.44 | 252 | 140273m  | 6.459   | ng/ul     |
| 86) Benzo(k)fluoranthene       | 24.49 | 252 | 43969m   | 2.073   | ng/ul     |
| 88) Benzo(a)pyrene             | 25.40 | 252 | 90085    | 4.284   | ng/ul# 91 |
| 89) Indeno(1,2,3-cd)pyrene     | 29.59 | 276 | 55037    | 2.235   | ng/ul# 96 |
| 91) Benzo(g,h,i)perylene       | 30.87 | 276 | 44953    | 2.219   | ng/ul# 86 |

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG082917\  
Data File : BG028541.D  
Acq On : 29 Aug 2017 19:23  
Operator : SJ/JU  
Sample : I4960-01  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
5-POST1

Manual Integrations  
APPROVED

Sohil  
8/30/2017 6:48:43 PM

Quant Time: Aug 30 15:10:59 2017  
Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080217.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Aug 30 01:19:11 2017  
Response via : Initial Calibration

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|--------------------|------|------|----------|------|-------|----------|
|--------------------|------|------|----------|------|-------|----------|

-----

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

