

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG082417\  
 Data File : BG028481.D  
 Acq On : 25 Aug 2017 5:21  
 Operator : SJ/JU  
 Sample : I4840-17  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 B0AN0

Quant Time: Aug 25 15:21:07 2017  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080217.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Aug 25 03:52:53 2017  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.34  | 152  | 45688    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.16 | 136  | 191951   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.95 | 164  | 135263   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.70 | 188  | 302536   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 22.03 | 240  | 350546   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.53 | 264  | 356678   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.82  | 96  | 1983   | 2.02  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.49  | 99  | 45304  | 9.03  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.65  | 67  | 30267  | 9.52  | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.87  | 132 | 35088  | 11.08 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 9.03  | 113 | 19496  | 4.65  | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.50  | 128 | 18840  | 12.55 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.23 | 143 | 22542  | 13.70 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.78 | 165 | 41006  | 12.04 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.30 | 131 | 21534  | 5.68  | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 14.34 | 166 | 142084 | 11.81 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.65 | 160 | 168164 | 12.40 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 15.16 | 143 | 13785  | 7.21  | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.94 | 176 | 126925 | 11.76 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 16.06 | 200 | 12744  | 6.48  | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.79 | 188 | 192573 | 13.27 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 20.07 | 212 | 227618 | 12.66 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 25.29 | 264 | 225931 | 13.53 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |       |              | Qvalue |
|----------------------------|-------|-----|-------|--------------|--------|
| 6) Phenol                  | 7.52  | 94  | 6318  | 1.267 ng/ul  | 98     |
| 44) Dimethylphthalate      | 14.38 | 163 | 61092 | 5.515 ng/ul  | 97     |
| 69) Phenanthrene           | 17.74 | 178 | 33474 | 2.199 ng/ul# | 93     |
| 74) Fluoranthene           | 19.74 | 202 | 98583 | 5.328 ng/ul  | 98     |
| 77) Pyrene                 | 20.10 | 202 | 87037 | 4.162 ng/ul  | 98     |
| 80) Benzo(a)anthracene     | 22.01 | 228 | 60606 | 2.992 ng/ul  | 98     |
| 82) Chrysene               | 22.08 | 228 | 58116 | 3.186 ng/ul  | 97     |
| 85) Benzo(b)fluoranthene   | 24.42 | 252 | 86917 | 4.072 ng/ul# | 92     |
| 86) Benzo(k)fluoranthene   | 24.42 | 252 | 86917 | 4.169 ng/ul# | 94     |
| 88) Benzo(a)pyrene         | 25.37 | 252 | 67451 | 3.264 ng/ul# | 93     |
| 89) Indeno(1,2,3-cd)pyrene | 29.55 | 276 | 55487 | 2.293 ng/ul# | 96     |
| 91) Benzo(g,h,i)perylene   | 30.83 | 276 | 45657 | 2.293 ng/ul  | 99     |

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

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