

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG052316\  
 Data File : BG022051.D  
 Acq On : 23 May 2016 17:57  
 Operator : UM/SJ  
 Sample : H3124-04  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 JHGG1

Manual Integrations  
 APPROVED

umangi  
 5/24/2016 7:37:12 PM

Quant Time: May 24 07:59:21 2016  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG052116.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 24 07:40:24 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.16  | 152  | 37861    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.98 | 136  | 201990   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.78 | 164  | 113273   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.53 | 188  | 208397   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.82 | 240  | 178409   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.15 | 264  | 169990   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.54  | 96  | 3042   | 4.19  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.31  | 99  | 72082  | 21.07 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.48  | 67  | 38280  | 21.34 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.69  | 132 | 62140  | 21.73 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.86  | 113 | 61936  | 21.35 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.33  | 128 | 29896  | 20.07 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.06 | 143 | 35279  | 21.86 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.60 | 165 | 58074  | 20.40 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.11 | 131 | 71787  | 23.15 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 14.18 | 166 | 190051 | 22.34 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.48 | 160 | 254484 | 21.42 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.99 | 143 | 21128m | 14.42 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.78 | 176 | 162183 | 20.36 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.89 | 200 | 15982  | 14.67 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.63 | 188 | 218325 | 21.83 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.91 | 212 | 216799 | 21.55 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 24.92 | 264 | 182459 | 23.13 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |        |      |        | Qvalue |
|----------------------------|-------|-----|--------|------|--------|--------|
| 44) Dimethylphthalate      | 14.23 | 163 | 29542  | 3.46 | ng/ul  | 99     |
| 69) Phenanthrene           | 17.57 | 178 | 13960  | 1.20 | ng/ul# | 89     |
| 74) Fluoranthene           | 19.57 | 202 | 56647  | 4.78 | ng/ul  | 98     |
| 77) Pyrene                 | 19.94 | 202 | 75955  | 6.05 | ng/ul  | 97     |
| 80) Benzo(a)anthracene     | 21.79 | 228 | 38868  | 3.71 | ng/ul  | 94     |
| 82) Chrysene               | 21.86 | 228 | 38083  | 3.79 | ng/ul  | 95     |
| 85) Benzo(b)fluoranthene   | 24.09 | 252 | 44761  | 4.50 | ng/ul# | 97     |
| 86) Benzo(k)fluoranthene   | 24.12 | 252 | 33570m | 3.47 | ng/ul  |        |
| 88) Benzo(a)pyrene         | 24.99 | 252 | 47152  | 4.91 | ng/ul  | 97     |
| 89) Indeno(1,2,3-cd)pyrene | 28.97 | 276 | 27680m | 2.52 | ng/ul  |        |
| 91) Benzo(g,h,i)perylene   | 30.16 | 276 | 22303m | 2.53 | ng/ul  |        |

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

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