

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121716\  
 Data File : BG025283.D  
 Acq On : 17 Dec 2016 20:48  
 Operator : UM/SJ  
 Sample : H6040-11  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 DA8W4

Manual Integrations  
 APPROVED

umangi  
 12/20/2016 10:54:26 AM

Quant Time: Dec 18 03:58:55 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun Dec 18 00:18:21 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.23  | 152  | 73626    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.06 | 136  | 287794   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.86 | 164  | 223994   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.60 | 188  | 459284   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.90 | 240  | 597948   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.34 | 264  | 614476   | 20.00 | ng/ul | 0.01     |

## System Monitoring Compounds

|                                |       |     |        |       |       |       |
|--------------------------------|-------|-----|--------|-------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.57  | 96  | 11014  | 7.73  | ng/uL | -0.02 |
| 5) Phenol-d5                   | 7.40  | 99  | 92068  | 14.49 | ng/ul | 0.00  |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.55  | 67  | 57397  | 14.61 | ng/ul | 0.00  |
| 9) 2-Chlorophenol-d4           | 7.77  | 132 | 63782  | 14.36 | ng/ul | 0.00  |
| 13) 4-Methylphenol-d8          | 8.95  | 113 | 76695  | 14.45 | ng/ul | 0.00  |
| 19) Nitrobenzene-d5            | 9.42  | 128 | 33832  | 16.60 | ng/ul | 0.00  |
| 22) 2-Nitrophenol-d4           | 10.14 | 143 | 38250  | 15.06 | ng/ul | 0.00  |
| 26) 2,4-Dichlorophenol-d3      | 10.69 | 165 | 76170  | 15.59 | ng/ul | 0.00  |
| 29) 4-Chloroaniline-d4         | 11.21 | 131 | 117792 | 24.53 | ng/ul | 0.00  |
| 43) Dimethylphthalate-d6       | 14.25 | 166 | 254420 | 16.51 | ng/ul | 0.00  |
| 46) Acenaphthylene-d8          | 14.56 | 160 | 296983 | 17.60 | ng/ul | 0.00  |
| 51) 4-Nitrophenol-d4           | 15.11 | 143 | 54130  | 20.44 | ng/ul | 0.02  |
| 57) Fluorene-d10               | 15.85 | 176 | 295629 | 20.38 | ng/ul | 0.00  |
| 62) 4,6-Dinitro-2-methylphenol | 15.99 | 200 | 63157  | 22.24 | ng/ul | 0.01  |
| 70) Anthracene-d10             | 17.70 | 188 | 360213 | 18.58 | ng/ul | 0.00  |
| 76) Pyrene-d10                 | 19.98 | 212 | 435790 | 17.69 | ng/ul | 0.00  |
| 87) Benzo(a)pyrene-d12         | 25.09 | 264 | 451744 | 18.14 | ng/ul | 0.01  |

## Target Compounds

|                         |       |     |         |        | Ovalue |    |
|-------------------------|-------|-----|---------|--------|--------|----|
| 6) Phenol               | 7.42  | 94  | 38003   | 5.83   | ng/ul# | 87 |
| 11) 2-Methylphenol      | 8.68  | 108 | 141415  | 29.46  | ng/ul  | 91 |
| 14) Acetophenone        | 9.07  | 105 | 27860   | 3.45   | ng/ul  | 86 |
| 16) 4-Methylphenol      | 9.02  | 108 | 383199  | 71.76  | ng/ul  | 93 |
| 24) 2,4-Dimethylphenol  | 10.22 | 107 | 832370m | 141.23 | ng/ul  |    |
| 28) Naphthalene         | 11.12 | 128 | 501752  | 37.51  | ng/ul  | 98 |
| 34) 2-Methylnaphthalene | 12.70 | 142 | 555591  | 52.63  | ng/ul  | 98 |
| 40) 1,1'-Biphenyl       | 13.69 | 154 | 127574  | 9.39   | ng/ul  | 97 |
| 44) Dimethylphthalate   | 14.31 | 163 | 72761   | 4.81   | ng/ul# | 96 |
| 49) Acenaphthene        | 14.92 | 153 | 43917   | 3.91   | ng/ul# | 76 |
| 53) Dibenzofuran        | 15.25 | 168 | 139299  | 7.84   | ng/ul  | 99 |
| 58) Fluorene            | 15.91 | 166 | 195177  | 13.49  | ng/ul# | 96 |
| 69) Phenanthrene        | 17.65 | 178 | 204284  | 9.32   | ng/ul# | 88 |
| 71) Anthracene          | 17.74 | 178 | 59691   | 2.65   | ng/ul# | 82 |

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

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