

Data Path : Z:\HPCHEM1\BNA G\DATA\BG063016\  
 Data File : BG022751.D  
 Acq On : 1 Jul 2016 7:00  
 Operator : UM/SJ  
 Sample : PB91665BL  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SBLK65

## Manual Integrations

sohil  
 7/1/2016 7:05:30 PM

Quant Time: Jul 01 08:36:14 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG063016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jul 01 08:12:10 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.16  | 152  | 103322   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.98 | 136  | 478817   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.80 | 164  | 391675   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.55 | 188  | 994236m  | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.85 | 240  | 1159380m | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.23 | 264  | 1149421m | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |          |       |       |      |
|--------------------------------|-------|-----|----------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.54  | 96  | 7087     | 3.83  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.31  | 99  | 366376   | 33.71 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.48  | 67  | 212284   | 34.64 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.69  | 132 | 236308   | 32.55 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.86  | 113 | 288423   | 34.16 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.33  | 128 | 130654   | 34.22 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.06 | 143 | 157181   | 34.13 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.60 | 165 | 317471   | 32.97 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.11 | 131 | 397687   | 48.77 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 14.19 | 166 | 1229257  | 38.63 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.49 | 160 | 1298279  | 35.60 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.99 | 143 | 181079   | 36.24 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.79 | 176 | 1114416  | 36.96 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.90 | 200 | 195655m  | 30.60 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.65 | 188 | 1748977m | 37.85 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.94 | 212 | 2069410  | 37.36 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 24.99 | 264 | 2129560  | 39.00 | ng/ul | 0.00 |

Target Compounds Ovalue

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

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