

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG072516\  
 Data File : BG023265.D  
 Acq On : 26 Jul 2016 00:00  
 Operator : UM/SJ  
 Sample : MDL-BLK-S-ML  
 Misc : MDL 4PPM  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 MDL-BLK-S-ML

## Manual Integrations

APPROVED  
 SOHIL  
 7/26/2016 11:41:20 AM

Quant Time: Jul 26 03:33:46 2016  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG072016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jul 21 04:24:30 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.14  | 152  | 169977   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.97 | 136  | 603108   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.81 | 164  | 436376   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.57 | 188  | 1002067  | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.89 | 240  | 1057637  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.30 | 264  | 1038363  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.50  | 96  | 11649   | 3.10  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.28  | 99  | 734278  | 41.90 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.45  | 67  | 564409  | 48.60 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.66  | 132 | 510078  | 43.33 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.85  | 113 | 381100  | 27.95 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.31  | 128 | 198174m | 41.08 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.05 | 143 | 218799  | 38.82 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.59 | 165 | 364259  | 34.39 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.11 | 131 | 548505  | 50.02 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 14.21 | 166 | 1404370 | 38.71 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.50 | 160 | 1655054 | 40.13 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 15.01 | 143 | 214644m | 33.52 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.80 | 176 | 1296440 | 39.52 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.92 | 200 | 247209  | 35.85 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.67 | 188 | 1908268 | 40.68 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.96 | 212 | 2119617 | 38.83 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 25.06 | 264 | 2018288 | 41.40 | ng/ul | 0.00 |

| Target Compounds | Qvalue |
|------------------|--------|
|------------------|--------|

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

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