

Data Path : Z:\HPCHEM1\BNA\_M\Data\BM121416\  
 Data File : BM008301.D  
 Acq On : 14 Dec 2016 09:00  
 Operator : UM/SJ  
 Sample : H5976-07MS  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DA8POMS

Quant Time: Dec 14 11:02:50 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM121416.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Dec 14 05:32:31 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.67  | 152  | 144084   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.45 | 136  | 561681   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.31 | 164  | 366537   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.06 | 188  | 767544   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.26 | 240  | 1005282  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.49 | 264  | 1027703  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.24  | 96  | 3136    | 1.02  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.86  | 99  | 57416   | 5.14  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.02  | 67  | 165224  | 25.72 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.21  | 132 | 177204  | 21.07 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.39  | 113 | 108481  | 12.53 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.83  | 128 | 111635  | 27.50 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.55  | 143 | 120174  | 26.72 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.09 | 165 | 202919  | 24.62 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.64 | 131 | 17505   | 1.85  | ng/ul | 0.04 |
| 43) Dimethylphthalate-d6       | 13.73 | 166 | 721764  | 29.62 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.00 | 160 | 835940  | 29.03 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.57 | 143 | 15173   | 4.23  | ng/ul | 0.02 |
| 57) Fluorene-d10               | 15.31 | 176 | 622809  | 29.41 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.47 | 200 | 114736  | 25.11 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.16 | 188 | 1026969 | 30.71 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.46 | 212 | 1292980 | 33.31 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.35 | 264 | 1374221 | 31.84 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |         |       |        | Qvalue |
|--------------------------------|-------|-----|---------|-------|--------|--------|
| 6) Phenol                      | 6.89  | 94  | 58111   | 5.01  | ng/ul  | 98     |
| 10) 2-Chlorophenol             | 7.25  | 128 | 172891  | 20.35 | ng/ul  | 97     |
| 15) N-Nitroso-di-n-propylamine | 8.47  | 70  | 192082  | 26.41 | ng/ul  | 97     |
| 33) 4-Chloro-3-methylphenol    | 11.76 | 107 | 203303  | 22.62 | ng/ul  | 100    |
| 47) Acenaphthylene             | 14.03 | 152 | 47620   | 1.63  | ng/ul# | 94     |
| 49) Acenaphthene               | 14.37 | 153 | 1562801 | 77.60 | ng/ul  | 99     |
| 52) 4-Nitrophenol              | 14.59 | 109 | 14133   | 4.10  | ng/ul  | 95     |
| 54) 2,4-Dinitrotoluene         | 14.72 | 165 | 191072  | 27.43 | ng/ul# | 72     |
| 58) Fluorene                   | 15.36 | 166 | 249401  | 10.34 | ng/ul  | 95     |
| 68) Pentachlorophenol          | 16.73 | 266 | 161357  | 31.24 | ng/ul  | 96     |
| 77) Pyrene                     | 19.49 | 202 | 1557601 | 31.56 | ng/ul  | 99     |

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

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