

Data Path : Z:\HPCHEM1\BNA M\DATA\BM021516\  
 Data File : BM004277.D  
 Acq On : 16 Feb 2016 03:46  
 Operator : SJ/UM  
 Sample : H1416-20MSD  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 D9QF3MSD

Quant Time: Feb 16 05:54:22 2016  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM021216.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Feb 16 02:45:44 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.65  | 152  | 31508    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.43 | 136  | 138587   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.31 | 164  | 72933    | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.06 | 188  | 149362   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.27 | 240  | 156620   | 20.00 | ng/ul | 0.01     |
| 86) Perylene-d12          | 23.51 | 264  | 154957   | 20.00 | ng/ul | 0.02     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.11  | 96  | 924    | 1.48  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.85  | 99  | 22958  | 9.07  | ng/ul | 0.02 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.98  | 67  | 44820  | 32.79 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.19  | 132 | 67397  | 30.03 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.38  | 113 | 42648  | 19.78 | ng/ul | 0.02 |
| 19) Nitrobenzene-d5            | 8.81  | 128 | 37924  | 35.59 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.53  | 143 | 43220  | 36.28 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.07 | 165 | 68379  | 31.83 | ng/ul | 0.01 |
| 29) 4-Chloroaniline-d4         | 10.59 | 131 | 4100   | 1.60  | ng/ul | 0.01 |
| 44) Dimethylphthalate-d6       | 13.72 | 166 | 209648 | 34.54 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.00 | 160 | 259948 | 35.25 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.57 | 143 | 7856   | 8.00  | ng/ul | 0.03 |
| 58) Fluorene-d10               | 15.30 | 176 | 175093 | 34.23 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.45 | 200 | 25469  | 28.00 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.16 | 188 | 258349 | 35.38 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.47 | 212 | 274538 | 33.02 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.37 | 264 | 288256 | 34.86 | ng/ul | 0.02 |

## Target Compounds

|                                |       |     |        |       |       | Ovalue |
|--------------------------------|-------|-----|--------|-------|-------|--------|
| 6) Phenol                      | 6.87  | 94  | 24399  | 9.05  | ng/ul | 99     |
| 10) 2-Chlorophenol             | 7.22  | 128 | 66431  | 28.47 | ng/ul | 98     |
| 15) N-Nitroso-di-n-propylamine | 8.46  | 70  | 53452  | 32.26 | ng/ul | 96     |
| 33) 4-Chloro-3-methylphenol    | 11.76 | 107 | 66876  | 26.99 | ng/ul | 99     |
| 50) Acenaphthene               | 14.37 | 153 | 178524 | 32.72 | ng/ul | 98     |
| 53) 4-Nitrophenol              | 14.59 | 109 | 6015   | 7.66  | ng/ul | 98     |
| 55) 2,4-Dinitrotoluene         | 14.69 | 165 | 60118  | 33.44 | ng/ul | 96     |
| 69) Pentachlorophenol          | 16.73 | 266 | 28603  | 31.77 | ng/ul | 100    |
| 80) Pvrene                     | 19.50 | 202 | 347874 | 30.71 | ng/ul | 99     |

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

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