

Data Path : Z:\HPCHEM1\BNA M\DATA\BM042816\  
 Data File : BM005117.D  
 Acq On : 28 Apr 2016 17:14  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD02050

Manual Integrations  
 APPROVED

sohil  
 4/29/2016 10:36:33 AM

Quant Time: Apr 29 03:55:53 2016  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM042516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Apr 29 02:20:37 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.79  | 152  | 41483    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.58 | 136  | 195500   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.43 | 164  | 124983   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.18 | 188  | 291654   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.37 | 240  | 293688   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.65 | 264  | 244373   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.29  | 96  | 7189   | 8.81  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.96  | 99  | 72908  | 21.25 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.12  | 67  | 42721  | 21.88 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.32  | 132 | 57131  | 21.04 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.50  | 113 | 61039  | 20.98 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.95  | 128 | 29004  | 21.42 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.66  | 143 | 33228  | 22.00 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.20 | 165 | 61150  | 21.31 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.72 | 131 | 83772  | 24.91 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.84 | 166 | 205457 | 20.81 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.12 | 160 | 243879 | 20.75 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.64 | 143 | 36253  | 21.01 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.43 | 176 | 176236 | 20.32 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.55 | 200 | 32027  | 19.31 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.27 | 188 | 275382 | 21.06 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.57 | 212 | 297720 | 22.27 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.50 | 264 | 227385 | 20.77 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |       |       | Ovalue |
|--------------------------------|-------|-----|--------|-------|-------|--------|
| 2) 1,4-Dioxane                 | 3.32  | 88  | 12789  | 8.44  | ng/uL | 96     |
| 4) Benzaldehyde                | 6.93  | 77  | 44492  | 26.11 | ng/ul | 99     |
| 6) Phenol                      | 6.99  | 94  | 76330  | 21.34 | ng/ul | 100    |
| 8) Bis(2-Chloroethyl)ether     | 7.22  | 93  | 58932  | 21.49 | ng/ul | 100    |
| 10) 2-Chlorophenol             | 7.36  | 128 | 58369  | 20.92 | ng/ul | 98     |
| 11) 2-Methylphenol             | 8.23  | 108 | 58772  | 21.05 | ng/ul | 98     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.32  | 45  | 76396  | 21.54 | ng/ul | 99     |
| 14) Acetophenone               | 8.61  | 105 | 96738  | 22.09 | ng/ul | 99     |
| 15) N-Nitroso-di-n-propylamine | 8.59  | 70  | 49438  | 22.54 | ng/ul | 98     |
| 16) 4-Methylphenol             | 8.56  | 108 | 66109  | 21.28 | ng/ul | 98     |
| 17) Hexachloroethane           | 8.86  | 117 | 23203  | 20.89 | ng/ul | 91     |
| 20) Nitrobenzene               | 8.99  | 77  | 70244  | 21.08 | ng/ul | 97     |
| 21) Isophorone                 | 9.51  | 82  | 140747 | 22.24 | ng/ul | 99     |
| 23) 2-Nitrophenol              | 9.70  | 139 | 35620  | 21.69 | ng/ul | 95     |
| 24) 2,4-Dimethylphenol         | 9.76  | 107 | 73679  | 20.84 | ng/ul | 99     |
| 25) Bis(2-Chloroethoxy)methane | 9.99  | 93  | 85550  | 21.76 | ng/ul | 99     |
| 27) 2,4-Dichlorophenol         | 10.23 | 162 | 62527  | 21.20 | ng/ul | 98     |
| 28) Naphthalene                | 10.63 | 128 | 201841 | 20.47 | ng/ul | 100    |
| 30) 4-Chloroaniline            | 10.74 | 127 | 83643  | 24.42 | ng/ul | 99     |
| 31) Hexachlorobutadiene        | 10.91 | 225 | 37515  | 19.27 | ng/ul | 97     |
| 32) Caprolactam                | 11.50 | 113 | 21449m | 21.03 | ng/ul |        |
| 33) 4-Chloro-3-methylphenol    | 11.87 | 107 | 73654  | 21.84 | ng/ul | 99     |
| 34) 2-Methylnaphthalene        | 12.24 | 142 | 151616 | 20.77 | ng/ul | 100    |

Data Path : Z:\HPCHEM1\BNA M\DATA\BM042816\  
 Data File : BM005117.D  
 Acq On : 28 Apr 2016 17:14  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD02050

Manual Integrations  
 APPROVED

sohil  
 4/29/2016 10:36:33 AM

Quant Time: Apr 29 03:55:53 2016  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM042516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Apr 29 02:20:37 2016  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|--------------------------------|-------|------|----------|-------|-------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 12.62 | 216  | 77972    | 19.97 | ng/ul | 99       |
| 37) Hexachlorocyclopentadiene  | 12.59 | 237  | 40274    | 17.38 | ng/ul | 96       |
| 38) 2,4,6-Trichlorophenol      | 12.86 | 196  | 53043    | 21.36 | ng/ul | 99       |
| 39) 2,4,5-Trichlorophenol      | 12.93 | 196  | 59306    | 21.32 | ng/ul | 98       |
| 40) 1,1'-Biphenyl              | 13.26 | 154  | 204969   | 20.73 | ng/ul | 98       |
| 41) 2-Chloronaphthalene        | 13.30 | 162  | 155000   | 20.61 | ng/ul | 99       |
| 42) 2-Nitroaniline             | 13.51 | 65   | 46852    | 22.97 | ng/ul | 98       |
| 44) Dimethylphthalate          | 13.89 | 163  | 205772   | 20.79 | ng/ul | 99       |
| 45) 2,6-Dinitrotoluene         | 14.01 | 165  | 41911    | 21.92 | ng/ul | 94       |
| 47) Acenaphthylene             | 14.15 | 152  | 260351   | 20.93 | ng/ul | 100      |
| 48) 3-Nitroaniline             | 14.34 | 138  | 44873    | 23.07 | ng/ul | 100      |
| 49) Acenaphthene               | 14.49 | 153  | 169738   | 20.74 | ng/ul | 99       |
| 50) 2,4-Dinitrophenol          | 14.55 | 184  | 17923    | 16.06 | ng/ul | 97       |
| 52) 4-Nitrophenol              | 14.65 | 109  | 30029    | 19.46 | ng/ul | 95       |
| 53) Dibenzofuran               | 14.83 | 168  | 246296   | 20.41 | ng/ul | 99       |
| 54) 2,4-Dinitrotoluene         | 14.80 | 165  | 61562    | 21.19 | ng/ul | 96       |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.06 | 232  | 50431    | 20.88 | ng/ul | 98       |
| 56) Diethylphthalate           | 15.25 | 149  | 208998   | 20.92 | ng/ul | 99       |
| 58) Fluorene                   | 15.48 | 166  | 196004   | 20.92 | ng/ul | 100      |
| 59) 4-Chlorophenyl-phenylether | 15.47 | 204  | 97026    | 20.53 | ng/ul | 98       |
| 60) 4-Nitroaniline             | 15.51 | 138  | 45801    | 21.84 | ng/ul | 100      |
| 63) 4,6-Dinitro-2-methylphenol | 15.56 | 198  | 33705    | 19.29 | ng/ul | 98       |
| 64) N-Nitrosodiphenylamine     | 15.69 | 169  | 174877   | 21.81 | ng/ul | 99       |
| 65) 4-Bromophenyl-phenylether  | 16.37 | 248  | 60280    | 21.18 | ng/ul | 99       |
| 66) Hexachlorobenzene          | 16.49 | 284  | 65861    | 20.31 | ng/ul | 99       |
| 67) Atrazine                   | 16.64 | 200  | 65436    | 22.15 | ng/ul | 99       |
| 68) Pentachlorophenol          | 16.83 | 266  | 37576    | 20.09 | ng/ul | 99       |
| 69) Phenanthrene               | 17.22 | 178  | 321295   | 20.70 | ng/ul | 99       |
| 71) Anthracene                 | 17.31 | 178  | 328963   | 21.02 | ng/ul | 99       |
| 72) Carbazole                  | 17.59 | 167  | 296818   | 22.12 | ng/ul | 99       |
| 73) Di-n-butylphthalate        | 18.14 | 149  | 356364   | 22.68 | ng/ul | 99       |
| 74) Fluoranthene               | 19.24 | 202  | 371383   | 21.74 | ng/ul | 100      |
| 77) Pyrene                     | 19.60 | 202  | 380141   | 22.41 | ng/ul | 99       |
| 78) Butylbenzylphthalate       | 20.50 | 149  | 153975   | 23.60 | ng/ul | 97       |
| 79) 3,3'-Dichlorobenzidine     | 21.29 | 252  | 109729   | 20.89 | ng/ul | 100      |
| 80) Benzo(a)anthracene         | 21.36 | 228  | 348754   | 20.75 | ng/ul | 99       |
| 81) Bis(2-ethylhexyl)phthalate | 21.27 | 149  | 207400   | 22.63 | ng/ul | 99       |
| 82) Chrysene                   | 21.41 | 228  | 333755   | 20.75 | ng/ul | 100      |
| 84) Di-n-octyl phthalate       | 22.17 | 149  | 332273   | 21.87 | ng/ul | 100      |
| 85) Benzo(b)fluoranthene       | 22.96 | 252  | 307504   | 20.89 | ng/ul | 99       |
| 86) Benzo(k)fluoranthene       | 23.01 | 252  | 296820   | 20.74 | ng/ul | 100      |
| 88) Benzo(a)pyrene             | 23.55 | 252  | 285148   | 20.75 | ng/ul | 99       |
| 89) Indeno(1,2,3-cd)pyrene     | 25.96 | 276  | 281019   | 21.43 | ng/ul | 99       |
| 90) Dibenzo(a,h)anthracene     | 25.97 | 278  | 236419   | 21.58 | ng/ul | 99       |
| 91) Benzo(g,h,i)perylene       | 26.67 | 276  | 231320   | 21.56 | ng/ul | 99       |

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

