

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092915\  
 Data File : BM002759.D  
 Acq On : 29 Sep 2015 23:14  
 Operator : UM/IZ  
 Sample : SSTD08094  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD08094

Quant Time: Sep 30 01:42:57 2015  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM092915.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 30 01:37:31 2015  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.70  | 152  | 67426    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.56 | 136  | 282451   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 15.29 | 164  | 140098   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 18.01 | 188  | 276630   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 22.09 | 240  | 280804   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 24.68 | 264  | 287955   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.76  | 96  | 45556   | 78.77 | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.82  | 99  | 395463  | 76.58 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.99  | 67  | 222041  | 76.86 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 8.22  | 132 | 368948  | 77.38 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 9.40  | 113 | 327584  | 76.99 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.90  | 128 | 170677  | 78.88 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.63 | 143 | 183082  | 80.10 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 11.17 | 165 | 307049  | 77.44 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.69 | 131 | 390045  | 79.39 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 14.68 | 166 | 792594  | 75.05 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 15.00 | 160 | 1067560 | 75.74 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 15.47 | 143 | 151075  | 79.50 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 16.27 | 176 | 702796  | 74.18 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 16.39 | 200 | 119415  | 91.40 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 18.11 | 188 | 988100  | 74.75 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 20.33 | 212 | 980080  | 74.37 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 24.52 | 264 | 1102663 | 75.06 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |         |       |       | Ovalue |
|--------------------------------|-------|-----|---------|-------|-------|--------|
| 2) 1,4-Dioxane                 | 3.80  | 88  | 48454   | 77.65 | ng/uL | 96     |
| 4) Benzaldehyde                | 7.82  | 77  | 221394  | 73.83 | ng/ul | 98     |
| 6) Phenol                      | 7.85  | 94  | 410185  | 76.79 | ng/ul | 99     |
| 8) Bis(2-Chloroethyl)ether     | 8.09  | 93  | 322880  | 77.11 | ng/ul | 98     |
| 10) 2-Chlorophenol             | 8.25  | 128 | 371875  | 76.98 | ng/ul | 99     |
| 11) 2-Methylphenol             | 9.13  | 108 | 317007  | 76.85 | ng/ul | 99     |
| 12) 2,2'-oxybis(1-Chloropropan | 9.23  | 45  | 441727  | 76.72 | ng/ul | 99     |
| 14) Acetophenone               | 9.54  | 105 | 487860  | 76.09 | ng/ul | 99     |
| 15) N-Nitroso-di-n-propylamine | 9.52  | 70  | 228703  | 76.71 | ng/ul | 99     |
| 16) 4-Methylphenol             | 9.47  | 108 | 341532  | 76.79 | ng/ul | 99     |
| 17) Hexachloroethane           | 9.81  | 117 | 139431  | 76.90 | ng/ul | 95     |
| 20) Nitrobenzene               | 9.94  | 77  | 330104  | 77.81 | ng/ul | 99     |
| 21) Isophorone                 | 10.46 | 82  | 616198  | 76.16 | ng/ul | 99     |
| 23) 2-Nitrophenol              | 10.66 | 139 | 201021  | 79.84 | ng/ul | 99     |
| 24) 2,4-Dimethylphenol         | 10.69 | 107 | 349830  | 76.62 | ng/ul | 98     |
| 25) Bis(2-Chloroethoxy)methane | 10.93 | 93  | 404997  | 76.47 | ng/ul | 99     |
| 27) 2,4-Dichlorophenol         | 11.20 | 162 | 307271  | 77.00 | ng/ul | 99     |
| 28) Naphthalene                | 11.61 | 128 | 1090980 | 75.85 | ng/ul | 99     |
| 30) 4-Chloroaniline            | 11.71 | 127 | 395791  | 78.27 | ng/ul | 99     |
| 31) Hexachlorobutadiene        | 11.86 | 225 | 178383  | 76.28 | ng/ul | 97     |
| 32) Caprolactam                | 12.47 | 113 | 101203  | 78.67 | ng/ul | 95     |
| 33) 4-Chloro-3-methylphenol    | 12.79 | 107 | 308420  | 76.23 | ng/ul | 98     |
| 34) 2-Methylnaphthalene        | 13.17 | 142 | 754640  | 75.37 | ng/ul | 99     |

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092915\  
 Data File : BM002759.D  
 Acq On : 29 Sep 2015 23:14  
 Operator : UM/IZ  
 Sample : SSTD08094  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD08094

Quant Time: Sep 30 01:42:57 2015  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM092915.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 30 01:37:31 2015  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|--------|--------|----------|
| 35) 1-Methylnaphthalene        | 13.39 | 142  | 731331   | 75.10  | ng/ul  | 99       |
| 37) 1,2,4,5-Tetrachlorobenzene | 13.52 | 216  | 333872   | 76.06  | ng/ul  | 99       |
| 38) Hexachlorocyclopentadiene  | 13.49 | 237  | 129591   | 117.09 | ng/ul  | 92       |
| 39) 2,4,6-Trichlorophenol      | 13.75 | 196  | 216085   | 77.27  | ng/ul  | 99       |
| 40) 2,4,5-Trichlorophenol      | 13.83 | 196  | 229914   | 77.29  | ng/ul  | 99       |
| 41) 1,1'-Biphenyl              | 14.14 | 154  | 904759   | 75.61  | ng/ul  | 100      |
| 42) 2-Chloronaphthalene        | 14.20 | 162  | 693533   | 75.93  | ng/ul  | 99       |
| 43) 2-Nitroaniline             | 14.39 | 65   | 175791   | 79.47  | ng/ul  | 97       |
| 45) Dimethylphthalate          | 14.73 | 163  | 806094   | 74.91  | ng/ul  | 100      |
| 46) 2,6-Dinitrotoluene         | 14.87 | 165  | 179773   | 80.86  | ng/ul  | 100      |
| 48) Acenaphthylene             | 15.03 | 152  | 1120959  | 75.66  | ng/ul  | 100      |
| 49) 3-Nitroaniline             | 15.20 | 138  | 183628   | 77.77  | ng/ul  | 96       |
| 50) Acenaphthene               | 15.36 | 153  | 733300   | 75.02  | ng/ul  | 100      |
| 51) 2,4-Dinitrophenol          | 15.42 | 184  | 75479    | 106.53 | ng/ul  | 97       |
| 53) 4-Nitrophenol              | 15.49 | 109  | 86074    | 79.83  | ng/ul  | 96       |
| 54) Dibenzofuran               | 15.69 | 168  | 991609   | 74.87  | ng/ul  | 100      |
| 55) 2,4-Dinitrotoluene         | 15.65 | 165  | 250429   | 80.35  | ng/ul  | 99       |
| 56) 2,3,4,6-Tetrachlorophenol  | 15.90 | 232  | 175858   | 79.41  | ng/ul  | 97       |
| 57) Diethylphthalate           | 16.06 | 149  | 811968   | 75.23  | ng/ul  | 99       |
| 59) Fluorene                   | 16.33 | 166  | 809689   | 74.35  | ng/ul  | 100      |
| 60) 4-Chlorophenyl-phenylether | 16.30 | 204  | 374383   | 74.46  | ng/ul  | 98       |
| 61) 4-Nitroaniline             | 16.34 | 138  | 196758   | 76.91  | ng/ul  | 99       |
| 64) 4,6-Dinitro-2-methylphenol | 16.41 | 198  | 127404   | 91.65  | ng/ul# | 97       |
| 65) N-Nitrosodiphenylamine     | 16.51 | 169  | 685582   | 75.70  | ng/ul  | 99       |
| 66) 4-Bromophenyl-phenylether  | 17.19 | 248  | 218664   | 75.75  | ng/ul  | 97       |
| 67) Hexachlorobenzene          | 17.32 | 284  | 246441   | 75.57  | ng/ul  | 99       |
| 68) Atrazine                   | 17.43 | 200  | 237431   | 76.55  | ng/ul  | 100      |
| 69) Pentachlorophenol          | 17.66 | 266  | 108346   | 85.40  | ng/ul  | 99       |
| 70) Phenanthrene               | 18.05 | 178  | 1190764  | 74.68  | ng/ul  | 99       |
| 72) Anthracene                 | 18.14 | 178  | 1218795  | 74.89  | ng/ul  | 99       |
| 73) 1,2,3,4-Tetrachlorobenzene | 14.12 | 216  | 318609   | 76.14  | ng/uL  | 98       |
| 74) Pentachlorobenzene         | 15.61 | 250  | 292965   | 75.32  | ng/uL  | 99       |
| 75) Carbazole                  | 18.40 | 167  | 1095694  | 75.20  | ng/ul  | 99       |
| 76) Di-n-butylphthalate        | 18.89 | 149  | 1379858  | 75.42  | ng/ul  | 100      |
| 77) Fluoranthene               | 20.01 | 202  | 1271916  | 74.64  | ng/ul  | 98       |
| 80) Pvrene                     | 20.37 | 202  | 1321241  | 74.34  | ng/ul  | 99       |
| 81) Butylbenzylphthalate       | 21.17 | 149  | 617478   | 76.12  | ng/ul  | 100      |
| 82) 3,3'-Dichlorobenzidine     | 21.98 | 252  | 389154   | 72.14  | ng/ul  | 100      |
| 83) Benzo(a)anthracene         | 22.07 | 228  | 1241336  | 74.49  | ng/ul  | 99       |
| 84) Bis(2-ethylhexyl)phthalate | 21.92 | 149  | 921356   | 76.39  | ng/ul  | 99       |
| 85) Chrysene                   | 22.13 | 228  | 1175015  | 74.94  | ng/ul  | 98       |
| 87) Di-n-octyl phthalate       | 22.89 | 149  | 1595235  | 75.59  | ng/ul  | 100      |
| 88) Benzo(b)fluoranthene       | 23.89 | 252  | 1265096  | 72.50  | ng/ul  | 99       |
| 89) Benzo(k)fluoranthene       | 23.94 | 252  | 1236242  | 76.33  | ng/ul  | 99       |
| 91) Benzo(a)pyrene             | 24.57 | 252  | 1238414  | 74.88  | ng/ul  | 99       |
| 92) Indeno(1,2,3-cd)pyrene     | 27.43 | 276  | 1447750  | 74.77  | ng/ul  | 100      |
| 93) Dibenzo(a,h)anthracene     | 27.42 | 278  | 1216977  | 74.42  | ng/ul  | 99       |
| 94) Benzo(a,h,i)perylene       | 28.27 | 276  | 1232342  | 75.68  | ng/ul  | 99       |

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092915\  
Data File : BM002759.D  
Acq On : 29 Sep 2015 23:14  
Operator : UM/IZ  
Sample : SSTD08094  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD08094

Quant Time: Sep 30 01:42:57 2015  
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM092915.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Sep 30 01:37:31 2015  
Response via : Initial Calibration

| Internal Standards                                                         | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------------------------------------------------------|------|------|----------|------|-------|----------|
| -----                                                                      |      |      |          |      |       |          |
| (#) = qualifier out of range (m) = manual integration (+) = signals summed |      |      |          |      |       |          |

