

Data Path : Z:\HPCHEM1\BNA\_M\Data\BM101617\  
 Data File : BM012216.D  
 Acq On : 16 Oct 2017 15:29  
 Operator : SJ/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleID :  
 SSTD02011

Manual Integrations  
 APPROVED

Sohil  
 10/17/2017 5:13:35 PM

Quant Time: Oct 16 23:32:55 2017  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Oct 16 23:30:15 2017  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.80  | 152  | 101309   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.60 | 136  | 450151   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.45 | 164  | 308937   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.20 | 188  | 804410   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.38 | 240  | 976519   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.69 | 264  | 924598   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.21  | 96  | 15184  | 7.12  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.97  | 99  | 149537 | 18.19 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.13  | 67  | 82570  | 16.75 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.33  | 132 | 131249 | 18.82 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.52  | 113 | 141339 | 19.97 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.97  | 128 | 66943  | 19.49 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.69  | 143 | 81338  | 20.26 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.23 | 165 | 163066 | 21.28 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.75 | 131 | 183582 | 25.60 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.86 | 166 | 565753 | 20.72 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.14 | 160 | 651222 | 19.70 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.67 | 143 | 73906  | 18.00 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.44 | 176 | 469110 | 20.93 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.59 | 200 | 97786  | 17.71 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.30 | 188 | 776396 | 19.52 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.59 | 212 | 940938 | 18.98 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.54 | 264 | 894576 | 20.07 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |        | Qvalue    |
|--------------------------------|-------|-----|--------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.25  | 88  | 15963  | 7.338  | ng/uL 100 |
| 4) Benzaldehyde                | 6.95  | 77  | 99900  | 18.188 | ng/ul 100 |
| 6) Phenol                      | 7.00  | 94  | 150346 | 17.693 | ng/ul 100 |
| 8) Bis(2-Chloroethyl)ether     | 7.22  | 93  | 115356 | 17.814 | ng/ul 100 |
| 10) 2-Chlorophenol             | 7.36  | 128 | 133301 | 18.733 | ng/ul 100 |
| 11) 2-Methylphenol             | 8.25  | 108 | 120125 | 18.796 | ng/ul 100 |
| 12) 2,2'-oxybis(1-Chloropropan | 8.33  | 45  | 119465 | 14.529 | ng/ul 100 |
| 14) Acetophenone               | 8.63  | 105 | 209155 | 19.446 | ng/ul 100 |
| 15) N-Nitroso-di-n-propylamine | 8.60  | 70  | 106498 | 18.476 | ng/ul 100 |
| 16) 4-Methylphenol             | 8.57  | 108 | 133760 | 18.993 | ng/ul 100 |
| 17) Hexachloroethane           | 8.87  | 117 | 55525  | 18.492 | ng/ul 100 |
| 20) Nitrobenzene               | 9.01  | 77  | 155082 | 18.170 | ng/ul 100 |
| 21) Isophorone                 | 9.53  | 82  | 304197 | 19.077 | ng/ul 100 |
| 23) 2-Nitrophenol              | 9.72  | 139 | 86034  | 19.987 | ng/ul 100 |
| 24) 2,4-Dimethylphenol         | 9.78  | 107 | 168804 | 19.459 | ng/ul 100 |
| 25) Bis(2-Chloroethoxy)methane | 10.01 | 93  | 175418 | 18.761 | ng/ul 100 |
| 27) 2,4-Dichlorophenol         | 10.26 | 162 | 156153 | 20.685 | ng/ul 100 |
| 28) Naphthalene                | 10.65 | 128 | 448002 | 19.236 | ng/ul 100 |
| 30) 4-Chloroaniline            | 10.77 | 127 | 179974 | 25.517 | ng/ul 100 |
| 31) Hexachlorobutadiene        | 10.92 | 225 | 121016 | 21.629 | ng/ul 100 |
| 32) Caprolactam                | 11.56 | 113 | 54878  | 23.027 | ng/ul 100 |
| 33) 4-Chloro-3-methylphenol    | 11.90 | 107 | 169030 | 21.371 | ng/ul 100 |
| 34) 2-Methylnaphthalene        | 12.27 | 142 | 361738 | 20.582 | ng/ul 100 |

Data Path : Z:\HPCHEM1\BNA\_M\Data\BM101617\  
 Data File : BM012216.D  
 Acq On : 16 Oct 2017 15:29  
 Operator : SJ/JU  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleID :  
 SSTD02011

Manual Integrations  
 APPROVED

Sohil  
 10/17/2017 5:13:35 PM

Quant Time: Oct 16 23:32:55 2017  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Oct 16 23:30:15 2017  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 12.64 | 216  | 232748   | 20.754 | ng/ul | 100      |
| 37) Hexachlorocyclopentadiene  | 12.60 | 237  | 113850   | 22.778 | ng/ul | 100      |
| 38) 2,4,6-Trichlorophenol      | 12.89 | 196  | 155410   | 20.901 | ng/ul | 100      |
| 39) 2,4,5-Trichlorophenol      | 12.97 | 196  | 159680   | 20.867 | ng/ul | 100      |
| 40) 1,1'-Biphenyl              | 13.28 | 154  | 500419   | 19.193 | ng/ul | 100      |
| 41) 2-Chloronaphthalene        | 13.33 | 162  | 397326   | 19.421 | ng/ul | 100      |
| 42) 2-Nitroaniline             | 13.54 | 65   | 105388   | 18.514 | ng/ul | 100      |
| 44) Dimethylphthalate          | 13.91 | 163  | 558219   | 20.406 | ng/ul | 100      |
| 45) 2,6-Dinitrotoluene         | 14.04 | 165  | 116831   | 21.516 | ng/ul | 100      |
| 47) Acenaphthylene             | 14.17 | 152  | 633809   | 19.528 | ng/ul | 100      |
| 48) 3-Nitroaniline             | 14.37 | 138  | 99364    | 23.520 | ng/ul | 100      |
| 49) Acenaphthene               | 14.52 | 153  | 424210   | 19.404 | ng/ul | 100      |
| 50) 2,4-Dinitrophenol          | 14.60 | 184  | 65361    | 21.197 | ng/ul | 100      |
| 52) 4-Nitrophenol              | 14.69 | 109  | 69340    | 18.796 | ng/ul | 100      |
| 53) Dibenzofuran               | 14.85 | 168  | 639030   | 20.300 | ng/ul | 100      |
| 54) 2,4-Dinitrotoluene         | 14.84 | 165  | 172524   | 21.922 | ng/ul | 100      |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.09 | 232  | 153971   | 21.987 | ng/ul | 100      |
| 56) Diethylphthalate           | 15.27 | 149  | 572137   | 19.349 | ng/ul | 100      |
| 58) Fluorene                   | 15.50 | 166  | 536421   | 20.518 | ng/ul | 100      |
| 59) 4-Chlorophenyl-phenylether | 15.49 | 204  | 299210   | 21.787 | ng/ul | 100      |
| 60) 4-Nitroaniline             | 15.54 | 138  | 90313m   | 17.313 | ng/ul |          |
| 63) 4,6-Dinitro-2-methylphenol | 15.61 | 198  | 104045   | 17.852 | ng/ul | 100      |
| 64) N-Nitrosodiphenylamine     | 15.71 | 169  | 471987   | 18.746 | ng/ul | 100      |
| 65) 4-Bromophenyl-phenylether  | 16.39 | 248  | 194505   | 20.200 | ng/ul | 100      |
| 66) Hexachlorobenzene          | 16.51 | 284  | 211759   | 19.412 | ng/ul | 100      |
| 67) Atrazine                   | 16.66 | 200  | 213479   | 20.277 | ng/ul | 100      |
| 68) Pentachlorophenol          | 16.86 | 266  | 105492   | 17.339 | ng/ul | 100      |
| 69) Phenanthrene               | 17.24 | 178  | 886774   | 19.378 | ng/ul | 100      |
| 71) Anthracene                 | 17.33 | 178  | 918627   | 19.391 | ng/ul | 100      |
| 72) Carbazole                  | 17.61 | 167  | 768828   | 19.176 | ng/ul | 100      |
| 73) Di-n-butylphthalate        | 18.15 | 149  | 1027582  | 18.460 | ng/ul | 100      |
| 74) Fluoranthene               | 19.26 | 202  | 1154882  | 20.913 | ng/ul | 100      |
| 77) Pyrene                     | 19.63 | 202  | 1200131  | 18.609 | ng/ul | 100      |
| 78) Butylbenzylphthalate       | 20.50 | 149  | 491550   | 17.122 | ng/ul | 100      |
| 79) 3,3'-Dichlorobenzidine     | 21.30 | 252  | 376677   | 19.382 | ng/ul | 100      |
| 80) Benzo(a)anthracene         | 21.37 | 228  | 1225376  | 19.285 | ng/ul | 100      |
| 81) Bis(2-ethylhexyl)phthalate | 21.27 | 149  | 712239   | 16.534 | ng/ul | 100      |
| 82) Chrysene                   | 21.42 | 228  | 1136307  | 19.047 | ng/ul | 100      |
| 84) Di-n-octyl phthalate       | 22.16 | 149  | 1226631  | 18.146 | ng/ul | 100      |
| 85) Benzo(b)fluoranthene       | 22.99 | 252  | 1174987  | 19.474 | ng/ul | 100      |
| 86) Benzo(k)fluoranthene       | 23.03 | 252  | 1188777  | 20.444 | ng/ul | 100      |
| 88) Benzo(a)pyrene             | 23.59 | 252  | 1120167  | 19.697 | ng/ul | 100      |
| 89) Indeno(1,2,3-cd)pyrene     | 26.04 | 276  | 1126248  | 18.649 | ng/ul | 100      |
| 90) Dibenzo(a,h)anthracene     | 26.05 | 278  | 915044   | 18.024 | ng/ul | 100      |
| 91) Benzo(g,h,i)perylene       | 26.77 | 276  | 900861   | 18.213 | ng/ul | 100      |

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

Data Path : Z:\HPCHEM1\BNA\_M\Data\BM101617\  
Data File : BM012216.D  
Acq On : 16 Oct 2017 15:29  
Operator : SJ/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD02011

Quant Time: Oct 16 23:32:55 2017  
Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Oct 16 23:30:15 2017  
Response via : Initial Calibration

Manual Integrations  
APPROVED

Sohil  
10/17/2017 5:13:35 PM

