

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM010617\  
 Data File : BM008732.D  
 Acq On : 06 Jan 2017 13:13  
 Operator : UM/SJ  
 Sample : SSTD02052  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD02052

Manual Integrations  
 APPROVED

Sohil  
 1/9/2017 10:44:58 AM

Quant Time: Jan 06 14:01:14 2017  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010617.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jan 06 13:54:38 2017  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.96  | 152  | 89572    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.77 | 136  | 388812   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.60 | 164  | 299941   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.34 | 188  | 695432m  | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.50 | 240  | 804712   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.87 | 264  | 761477   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.41  | 96  | 17012  | 9.39  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.11  | 99  | 143109 | 21.43 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.29  | 67  | 99778  | 25.99 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.49  | 132 | 105622 | 20.44 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.66  | 113 | 113659 | 21.90 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.13  | 128 | 48988  | 17.55 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.85  | 143 | 48406  | 15.24 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.38 | 165 | 119620 | 20.68 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.90 | 131 | 139683 | 24.80 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 14.00 | 166 | 395477 | 20.46 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.29 | 160 | 469752 | 19.78 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.77 | 143 | 61195  | 20.49 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.59 | 176 | 366632 | 21.98 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.69 | 200 | 64320  | 15.58 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.43 | 188 | 594494 | 19.87 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.72 | 212 | 704152 | 21.62 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.72 | 264 | 606584 | 19.12 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |       |        | Qvalue |
|--------------------------------|-------|-----|--------|-------|--------|--------|
| 2) 1,4-Dioxane                 | 3.45  | 88  | 19503  | 9.37  | ng/uL# | 75     |
| 4) Benzaldehyde                | 7.11  | 77  | 123288 | 25.51 | ng/ul  | 86     |
| 6) Phenol                      | 7.14  | 94  | 158146 | 23.07 | ng/ul# | 67     |
| 8) Bis(2-Chloroethyl)ether     | 7.39  | 93  | 115703 | 22.24 | ng/ul# | 84     |
| 10) 2-Chlorophenol             | 7.53  | 128 | 102944 | 19.45 | ng/ul# | 75     |
| 11) 2-Methylphenol             | 8.40  | 108 | 114823 | 22.59 | ng/ul  | 100    |
| 12) 2,2'-oxybis(1-Chloropropan | 8.50  | 45  | 185050 | 32.00 | ng/ul# | 92     |
| 14) Acetophenone               | 8.79  | 105 | 199376 | 24.02 | ng/ul# | 73     |
| 15) N-Nitroso-di-n-propylamine | 8.77  | 70  | 116935 | 26.84 | ng/ul# | 82     |
| 16) 4-Methylphenol             | 8.72  | 108 | 126272 | 23.05 | ng/ul  | 95     |
| 17) Hexachloroethane           | 9.05  | 117 | 55997  | 23.72 | ng/ul  | 94     |
| 20) Nitrobenzene               | 9.17  | 77  | 172282 | 24.82 | ng/ul# | 86     |
| 21) Isophorone                 | 9.70  | 82  | 300038 | 23.46 | ng/ul# | 93     |
| 23) 2-Nitrophenol              | 9.88  | 139 | 55364  | 17.09 | ng/ul# | 55     |
| 24) 2,4-Dimethylphenol         | 9.93  | 107 | 157934 | 22.72 | ng/ul# | 88     |
| 25) Bis(2-Chloroethoxy)methane | 10.17 | 93  | 161351 | 21.74 | ng/ul  | 96     |
| 27) 2,4-Dichlorophenol         | 10.41 | 162 | 118501 | 20.95 | ng/ul  | 91     |
| 28) Naphthalene                | 10.82 | 128 | 351837 | 19.48 | ng/ul  | 99     |
| 30) 4-Chloroaniline            | 10.93 | 127 | 143568 | 24.75 | ng/ul  | 99     |
| 31) Hexachlorobutadiene        | 11.10 | 225 | 107670 | 24.00 | ng/ul  | 90     |
| 32) Caprolactam                | 11.72 | 113 | 34573m | 21.91 | ng/ul  |        |
| 33) 4-Chloro-3-methylphenol    | 12.03 | 107 | 143054 | 23.53 | ng/ul  | 93     |
| 34) 2-Methylnaphthalene        | 12.42 | 142 | 273099 | 21.12 | ng/ul  | 99     |

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM010617\  
 Data File : BM008732.D  
 Acq On : 06 Jan 2017 13:13  
 Operator : UM/SJ  
 Sample : SSTD02052  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD02052

Manual Integrations  
 APPROVED

Sohil  
 1/9/2017 10:44:58 AM

Quant Time: Jan 06 14:01:14 2017  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM010617.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jan 06 13:54:38 2017  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 12.79 | 216  | 183617   | 19.71 | ng/ul# | 95       |
| 37) Hexachlorocyclopentadiene  | 12.76 | 237  | 124262   | 26.94 | ng/ul# | 96       |
| 38) 2,4,6-Trichlorophenol      | 13.02 | 196  | 108424   | 19.41 | ng/ul  | 98       |
| 39) 2,4,5-Trichlorophenol      | 13.09 | 196  | 118277   | 19.99 | ng/ul  | 94       |
| 40) 1,1'-Biphenyl              | 13.43 | 154  | 375255   | 18.56 | ng/ul  | 98       |
| 41) 2-Chloronaphthalene        | 13.47 | 162  | 297299   | 19.25 | ng/ul  | 96       |
| 42) 2-Nitroaniline             | 13.67 | 65   | 100570   | 22.80 | ng/ul# | 82       |
| 44) Dimethylphthalate          | 14.05 | 163  | 401038   | 21.15 | ng/ul  | 98       |
| 45) 2,6-Dinitrotoluene         | 14.17 | 165  | 69572    | 18.52 | ng/ul# | 67       |
| 47) Acenaphthylene             | 14.32 | 152  | 464281   | 19.37 | ng/ul  | 98       |
| 48) 3-Nitroaniline             | 14.50 | 138  | 69512    | 19.72 | ng/ul# | 77       |
| 49) Acenaphthene               | 14.66 | 153  | 317251   | 19.38 | ng/ul  | 98       |
| 50) 2,4-Dinitrophenol          | 14.70 | 184  | 43494    | 18.33 | ng/ul  | 94       |
| 52) 4-Nitrophenol              | 14.79 | 109  | 101199   | 36.04 | ng/ul# | 68       |
| 53) Dibenzofuran               | 14.99 | 168  | 485837   | 20.86 | ng/ul  | 93       |
| 54) 2,4-Dinitrotoluene         | 14.95 | 165  | 103107   | 19.25 | ng/ul# | 83       |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.21 | 232  | 112423   | 21.16 | ng/ul# | 97       |
| 56) Diethylphthalate           | 15.41 | 149  | 422304   | 20.81 | ng/ul  | 99       |
| 58) Fluorene                   | 15.64 | 166  | 393342   | 20.80 | ng/ul  | 99       |
| 59) 4-Chlorophenyl-phenylether | 15.63 | 204  | 226130   | 22.59 | ng/ul  | 94       |
| 60) 4-Nitroaniline             | 15.66 | 138  | 75002    | 21.22 | ng/ul# | 50       |
| 63) 4,6-Dinitro-2-methylphenol | 15.71 | 198  | 67464    | 15.89 | ng/ul# | 87       |
| 64) N-Nitrosodiphenylamine     | 15.84 | 169  | 345229   | 17.79 | ng/ul  | 97       |
| 65) 4-Bromophenyl-phenylether  | 16.53 | 248  | 140082m  | 18.07 | ng/ul  |          |
| 66) Hexachlorobenzene          | 16.63 | 284  | 161316   | 18.91 | ng/ul# | 85       |
| 67) Atrazine                   | 16.79 | 200  | 151181   | 19.88 | ng/ul  | 92       |
| 68) Pentachlorophenol          | 16.98 | 266  | 95281    | 19.30 | ng/ul  | 95       |
| 69) Phenanthrene               | 17.38 | 178  | 671617m  | 19.31 | ng/ul  |          |
| 71) Anthracene                 | 17.47 | 178  | 681010   | 19.19 | ng/ul  | 96       |
| 72) Carbazole                  | 17.74 | 167  | 596106   | 19.34 | ng/ul  | 100      |
| 73) Di-n-butylphthalate        | 18.30 | 149  | 710640   | 19.45 | ng/ul# | 97       |
| 74) Fluoranthene               | 19.39 | 202  | 836660   | 20.72 | ng/ul# | 94       |
| 77) Pyrene                     | 19.75 | 202  | 862033   | 20.82 | ng/ul# | 93       |
| 78) Butylbenzylphthalate       | 20.64 | 149  | 316285   | 19.84 | ng/ul  | 88       |
| 79) 3,3'-Dichlorobenzidine     | 21.42 | 252  | 315162   | 22.18 | ng/ul  | 98       |
| 80) Benzo(a)anthracene         | 21.49 | 228  | 856628   | 20.63 | ng/ul  | 100      |
| 81) Bis(2-ethylhexyl)phthalate | 21.41 | 149  | 427809   | 17.80 | ng/ul  | 99       |
| 82) Chrysene                   | 21.54 | 228  | 793482   | 20.23 | ng/ul  | 99       |
| 84) Di-n-octyl phthalate       | 22.33 | 149  | 724410   | 18.48 | ng/ul# | 91       |
| 85) Benzo(b)fluoranthene       | 23.15 | 252  | 786964   | 19.30 | ng/ul# | 94       |
| 86) Benzo(k)fluoranthene       | 23.20 | 252  | 766349   | 19.93 | ng/ul# | 98       |
| 88) Benzo(a)pyrene             | 23.77 | 252  | 749353   | 19.12 | ng/ul# | 95       |
| 89) Indeno(1,2,3-cd)pyrene     | 26.30 | 276  | 861192   | 18.02 | ng/ul# | 93       |
| 90) Dibenzo(a,h)anthracene     | 26.31 | 278  | 730666   | 18.13 | ng/ul# | 96       |
| 91) Benzo(g,h,i)perylene       | 27.04 | 276  | 732245   | 18.43 | ng/ul# | 95       |

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

