

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM040819\  
 Data File : BM019814.D  
 Acq On : 09 Apr 2019 02:54  
 Operator : JU/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD02081

Quant Time: Apr 09 07:13:21 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Apr 09 06:31:52 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.56  | 152  | 187655   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.33 | 136  | 734834   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.22 | 164  | 366240   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.98 | 188  | 758545   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.19 | 240  | 791536   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.39 | 264  | 929371   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.11  | 96  | 34868  | 7.27  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.75  | 99  | 305064 | 18.26 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.91  | 67  | 199621 | 17.84 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.09  | 132 | 250705 | 19.80 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.27  | 113 | 228499 | 18.22 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.73  | 128 | 111040 | 21.18 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.44  | 143 | 126830 | 21.74 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.97  | 165 | 224417 | 20.39 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.50 | 131 | 272612 | 20.62 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.64 | 166 | 552688 | 18.70 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.91 | 160 | 731439 | 19.79 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.47 | 143 | 72223  | 15.59 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.22 | 176 | 488140 | 19.17 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.37 | 200 | 75593  | 15.07 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.08 | 188 | 711996 | 19.56 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.39 | 212 | 723531 | 19.79 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.25 | 264 | 947211 | 19.23 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |        | Qvalue    |
|--------------------------------|-------|-----|--------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.15  | 88  | 39215  | 7.190  | ng/uL# 90 |
| 4) Benzaldehyde                | 6.73  | 77  | 234048 | 19.553 | ng/ul 97  |
| 6) Phenol                      | 6.77  | 94  | 318822 | 18.283 | ng/ul 99  |
| 8) Bis(2-Chloroethyl)ether     | 7.00  | 93  | 236463 | 17.725 | ng/ul 96  |
| 10) 2-Chlorophenol             | 7.12  | 128 | 253610 | 19.698 | ng/ul 98  |
| 11) 2-Methylphenol             | 8.01  | 108 | 226505 | 18.625 | ng/ul 100 |
| 12) 2,2'-oxybis(1-Chloropropan | 8.07  | 45  | 124265 | 9.827  | ng/ul# 79 |
| 14) Acetophenone               | 8.39  | 105 | 352914 | 17.276 | ng/ul# 92 |
| 15) N-Nitroso-di-n-propylamine | 8.37  | 70  | 206339 | 18.024 | ng/ul# 86 |
| 16) 4-Methylphenol             | 8.34  | 108 | 239567 | 17.961 | ng/ul 99  |
| 17) Hexachloroethane           | 8.60  | 117 | 105254 | 19.452 | ng/ul 82  |
| 20) Nitrobenzene               | 8.77  | 77  | 298155 | 19.108 | ng/ul 98  |
| 21) Isophorone                 | 9.28  | 82  | 519840 | 19.419 | ng/ul 98  |
| 23) 2-Nitrophenol              | 9.47  | 139 | 131702 | 20.465 | ng/ul 96  |
| 24) 2,4-Dimethylphenol         | 9.53  | 107 | 266435 | 19.052 | ng/ul 96  |
| 25) Bis(2-Chloroethoxy)methane | 9.77  | 93  | 301177 | 18.645 | ng/ul 97  |
| 27) 2,4-Dichlorophenol         | 10.00 | 162 | 221598 | 20.371 | ng/ul 97  |
| 28) Naphthalene                | 10.39 | 128 | 724135 | 19.022 | ng/ul 99  |
| 30) 4-Chloroaniline            | 10.52 | 127 | 275084 | 19.765 | ng/ul 99  |
| 31) Hexachlorobutadiene        | 10.64 | 225 | 138779 | 20.223 | ng/ul 98  |
| 32) Caprolactam                | 11.33 | 113 | 43483  | 13.463 | ng/ul# 78 |
| 33) 4-Chloro-3-methylphenol    | 11.66 | 107 | 219987 | 18.761 | ng/ul 97  |
| 34) 2-Methylnaphthalene        | 12.01 | 142 | 488867 | 18.327 | ng/ul 97  |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM040819\  
 Data File : BM019814.D  
 Acq On : 09 Apr 2019 02:54  
 Operator : JU/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD02081

Quant Time: Apr 09 07:13:21 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Apr 09 06:31:52 2019  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|--------|--------|----------|
| 35) 1-Methylnaphthalene        | 12.23 | 142  | 452371   | 17.961 | ng/ul  | 97       |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.38 | 216  | 236994   | 20.336 | ng/ul# | 96       |
| 38) Hexachlorocyclopentadiene  | 12.34 | 237  | 165323   | 27.656 | ng/ul# | 97       |
| 39) 2,4,6-Trichlorophenol      | 12.65 | 196  | 154543   | 21.255 | ng/ul  | 99       |
| 40) 2,4,5-Trichlorophenol      | 12.72 | 196  | 161428   | 20.702 | ng/ul  | 99       |
| 41) 1,1'-Biphenyl              | 13.04 | 154  | 613017   | 19.652 | ng/ul# | 94       |
| 42) 2-Chloronaphthalene        | 13.09 | 162  | 474749   | 19.874 | ng/ul  | 99       |
| 43) 2-Nitroaniline             | 13.32 | 65   | 129288   | 19.708 | ng/ul  | 94       |
| 45) Dimethylphthalate          | 13.69 | 163  | 562192   | 18.961 | ng/ul# | 98       |
| 46) 2,6-Dinitrotoluene         | 13.83 | 165  | 115270   | 21.177 | ng/ul  | 98       |
| 48) Acenaphthylene             | 13.94 | 152  | 708152   | 19.336 | ng/ul  | 99       |
| 49) 3-Nitroaniline             | 14.17 | 138  | 110097   | 20.670 | ng/ul# | 96       |
| 50) Acenaphthene               | 14.28 | 153  | 497473   | 19.287 | ng/ul  | 99       |
| 51) 2,4-Dinitrophenol          | 14.39 | 184  | 62778    | 19.919 | ng/ul# | 89       |
| 53) 4-Nitrophenol              | 14.49 | 109  | 55151    | 15.212 | ng/ul# | 78       |
| 54) Dibenzofuran               | 14.62 | 168  | 681658   | 18.924 | ng/ul  | 95       |
| 55) 2,4-Dinitrotoluene         | 14.63 | 165  | 164370   | 20.393 | ng/ul# | 72       |
| 56) 2,3,4,6-Tetrachlorophenol  | 14.86 | 232  | 130349   | 19.844 | ng/ul# | 80       |
| 57) Diethylphthalate           | 15.06 | 149  | 550339   | 18.871 | ng/ul  | 97       |
| 59) Fluorene                   | 15.27 | 166  | 546302   | 19.139 | ng/ul  | 100      |
| 60) 4-Chlorophenyl-phenylether | 15.27 | 204  | 268267   | 18.840 | ng/ul  | 94       |
| 61) 4-Nitroaniline             | 15.34 | 138  | 116630   | 20.060 | ng/ul  | 92       |
| 64) 4,6-Dinitro-2-methylphenol | 15.39 | 198  | 80322    | 15.124 | ng/ul# | 91       |
| 65) N-Nitrosodiphenylamine     | 15.50 | 169  | 467128   | 19.794 | ng/ul  | 99       |
| 66) 4-Bromophenyl-phenylether  | 16.17 | 248  | 164393   | 19.956 | ng/ul  | 97       |
| 67) Hexachlorobenzene          | 16.27 | 284  | 192870   | 19.335 | ng/ul  | 92       |
| 68) Atrazine                   | 16.46 | 200  | 159356   | 19.296 | ng/ul  | 95       |
| 69) Pentachlorophenol          | 16.64 | 266  | 97324    | 18.667 | ng/ul  | 98       |
| 70) Phenanthrene               | 17.02 | 178  | 811205   | 18.862 | ng/ul  | 99       |
| 72) Anthracene                 | 17.12 | 178  | 843724   | 19.229 | ng/ul  | 98       |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.00 | 216  | 225987   | 20.549 | ng/uL  | 99       |
| 74) Pentachlorobenzene         | 14.53 | 250  | 228533   | 19.867 | ng/uL  | 98       |
| 75) Carbazole                  | 17.40 | 167  | 727143   | 18.479 | ng/ul  | 98       |
| 76) Di-n-butylphthalate        | 17.95 | 149  | 944867   | 21.662 | ng/ul  | 100      |
| 77) Fluoranthene               | 19.05 | 202  | 912361   | 18.513 | ng/ul# | 87       |
| 80) Pyrene                     | 19.42 | 202  | 924909   | 19.423 | ng/ul# | 86       |
| 81) Butylbenzylphthalate       | 20.33 | 149  | 445711   | 23.639 | ng/ul  | 93       |
| 82) 3,3'-Dichlorobenzidine     | 21.12 | 252  | 364796   | 21.215 | ng/ul# | 98       |
| 83) Benzo(a)anthracene         | 21.17 | 228  | 942397   | 19.781 | ng/ul  | 99       |
| 84) Bis(2-ethylhexyl)phthalate | 21.09 | 149  | 624580   | 22.854 | ng/ul  | 96       |
| 85) Chrysene                   | 21.23 | 228  | 881254   | 19.278 | ng/ul  | 99       |
| 87) Di-n-octyl phthalate       | 21.96 | 149  | 1161625  | 20.039 | ng/ul  | 100      |
| 88) Benzo(b)fluoranthene       | 22.73 | 252  | 1037151  | 18.321 | ng/ul# | 96       |
| 89) Benzo(k)fluoranthene       | 22.77 | 252  | 998562   | 18.368 | ng/ul# | 97       |
| 91) Benzo(a)pyrene             | 23.30 | 252  | 1029901  | 19.031 | ng/ul# | 96       |
| 92) Indeno(1,2,3-cd)pyrene     | 25.61 | 276  | 1266942  | 18.695 | ng/ul# | 89       |
| 93) Dibenzo(a,h)anthracene     | 25.61 | 278  | 1079772  | 18.689 | ng/ul# | 96       |
| 94) Benzo(g,h,i)perylene       | 26.28 | 276  | 1056710  | 18.659 | ng/ul# | 93       |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM040819\  
Data File : BM019814.D  
Acq On : 09 Apr 2019 02:54  
Operator : JU/SJ  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD02081

Quant Time: Apr 09 07:13:21 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Apr 09 06:31:52 2019  
Response via : Initial Calibration

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

