

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031620\  
 Data File : BM025584.D  
 Acq On : 16 Mar 2020 17:32  
 Operator : CG/JU  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD02009

Quant Time: Mar 16 18:11:32 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM031420MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Mar 16 04:27:51 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.64  | 152  | 196184   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.40 | 136  | 825276   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.27 | 164  | 521001   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.02 | 188  | 1098819  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.20 | 240  | 1087665  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.39 | 264  | 1251049  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.22  | 96  | 36126   | 7.73  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.82  | 99  | 326596  | 21.09 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.98  | 67  | 189206  | 21.41 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.17  | 132 | 275724  | 21.94 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.35  | 113 | 271383  | 21.33 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.78  | 128 | 135187  | 22.04 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.50  | 143 | 146609  | 22.48 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.03 | 165 | 290354  | 21.77 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.54 | 131 | 317747  | 20.52 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.69 | 166 | 816390  | 20.62 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.96 | 160 | 1021392 | 21.53 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.47 | 143 | 155661  | 20.65 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.27 | 176 | 719405  | 20.70 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.39 | 200 | 123378  | 18.40 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.11 | 188 | 1090175 | 21.24 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.41 | 212 | 1151725 | 22.05 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.25 | 264 | 1348677 | 21.06 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |        | Ovalue   |
|--------------------------------|-------|-----|--------|--------|----------|
| 2) 1,4-Dioxane                 | 3.25  | 88  | 37908  | 7.529  | ng/uL 92 |
| 4) Benzaldehyde                | 6.79  | 77  | 124739 | 15.211 | ng/ul 97 |
| 6) Phenol                      | 6.85  | 94  | 343022 | 21.215 | ng/ul 95 |
| 8) Bis(2-Chloroethyl)ether     | 7.07  | 93  | 271183 | 21.179 | ng/ul 99 |
| 10) 2-Chlorophenol             | 7.21  | 128 | 283527 | 21.434 | ng/ul 98 |
| 11) 2-Methylphenol             | 8.08  | 108 | 266750 | 21.561 | ng/ul 99 |
| 12) 2,2'-oxybis(1-Chloropropan | 8.17  | 45  | 369566 | 21.632 | ng/ul 97 |
| 14) Acetophenone               | 8.45  | 105 | 416600 | 21.229 | ng/ul 98 |
| 15) N-Nitroso-di-n-propylamine | 8.45  | 70  | 212764 | 21.642 | ng/ul 98 |
| 16) 4-Methylphenol             | 8.40  | 108 | 290091 | 21.479 | ng/ul 97 |
| 17) Hexachloroethane           | 8.71  | 117 | 117370 | 21.343 | ng/ul 96 |
| 20) Nitrobenzene               | 8.82  | 77  | 316341 | 21.169 | ng/ul 98 |
| 21) Isophorone                 | 9.35  | 82  | 595413 | 21.576 | ng/ul 99 |
| 23) 2-Nitrophenol              | 9.53  | 139 | 160133 | 21.865 | ng/ul 93 |
| 24) 2,4-Dimethylphenol         | 9.60  | 107 | 319101 | 21.487 | ng/ul 98 |
| 25) Bis(2-Chloroethoxy)methane | 9.83  | 93  | 374377 | 21.167 | ng/ul 99 |
| 27) 2,4-Dichlorophenol         | 10.06 | 162 | 285867 | 21.607 | ng/ul 98 |
| 28) Naphthalene                | 10.46 | 128 | 938650 | 20.884 | ng/ul 99 |
| 30) 4-Chloroaniline            | 10.56 | 127 | 323044 | 20.561 | ng/ul 97 |
| 31) Hexachlorobutadiene        | 10.76 | 225 | 177261 | 20.507 | ng/ul 98 |
| 32) Caprolactam                | 11.32 | 113 | 86657  | 19.806 | ng/ul 92 |
| 33) 4-Chloro-3-methylphenol    | 11.70 | 107 | 297609 | 21.638 | ng/ul 98 |
| 34) 2-Methylnaphthalene        | 12.07 | 142 | 666058 | 20.748 | ng/ul 99 |

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031620\  
 Data File : BM025584.D  
 Acq On : 16 Mar 2020 17:32  
 Operator : CG/JU  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleID :  
 SSTD02009

Quant Time: Mar 16 18:11:32 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM031420MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Mar 16 04:27:51 2020  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 35) 1-Methylnaphthalene        | 12.29 | 142  | 661197   | 20.654 | ng/ul | 99       |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.45 | 216  | 344751   | 20.922 | ng/ul | 98       |
| 38) Hexachlorocyclopentadiene  | 12.44 | 237  | 199266   | 17.720 | ng/ul | 99       |
| 39) 2,4,6-Trichlorophenol      | 12.69 | 196  | 224934   | 21.471 | ng/ul | 99       |
| 40) 2,4,5-Trichlorophenol      | 12.76 | 196  | 245555   | 21.534 | ng/ul | 98       |
| 41) 1,1'-Biphenyl              | 13.10 | 154  | 876038   | 21.044 | ng/ul | 99       |
| 42) 2-Chloronaphthalene        | 13.13 | 162  | 687703   | 20.974 | ng/ul | 99       |
| 43) 2-Nitroaniline             | 13.34 | 65   | 183356   | 22.376 | ng/ul | 100      |
| 45) Dimethylphthalate          | 13.73 | 163  | 852756   | 20.689 | ng/ul | 99       |
| 46) 2,6-Dinitrotoluene         | 13.85 | 165  | 177421   | 21.885 | ng/ul | 96       |
| 48) Acenaphthylene             | 13.99 | 152  | 1101784  | 21.444 | ng/ul | 99       |
| 49) 3-Nitroaniline             | 14.17 | 138  | 147187   | 20.125 | ng/ul | 92       |
| 50) Acenaphthene               | 14.33 | 153  | 733281   | 20.994 | ng/ul | 99       |
| 51) 2,4-Dinitrophenol          | 14.38 | 184  | 85856    | 17.653 | ng/ul | 98       |
| 53) 4-Nitrophenol              | 14.49 | 109  | 114349   | 19.404 | ng/ul | 98       |
| 54) Dibenzofuran               | 14.67 | 168  | 1023601  | 20.659 | ng/ul | 98       |
| 55) 2,4-Dinitrotoluene         | 14.63 | 165  | 258486   | 21.809 | ng/ul | 96       |
| 56) 2,3,4,6-Tetrachlorophenol  | 14.90 | 232  | 213347   | 20.642 | ng/ul | 96       |
| 57) Diethylphthalate           | 15.11 | 149  | 854120   | 20.217 | ng/ul | 100      |
| 59) Fluorene                   | 15.32 | 166  | 864193   | 20.735 | ng/ul | 99       |
| 60) 4-Chlorophenyl-phenylether | 15.32 | 204  | 430373   | 20.176 | ng/ul | 98       |
| 61) 4-Nitroaniline             | 15.33 | 138  | 178719   | 20.536 | ng/ul | 98       |
| 64) 4,6-Dinitro-2-methylphenol | 15.40 | 198  | 137669   | 18.778 | ng/ul | 97       |
| 65) N-Nitrosodiphenylamine     | 15.53 | 169  | 732422   | 21.570 | ng/ul | 99       |
| 66) 4-Bromophenyl-phenylether  | 16.22 | 248  | 283585   | 20.927 | ng/ul | 99       |
| 67) Hexachlorobenzene          | 16.33 | 284  | 335307   | 20.153 | ng/ul | 97       |
| 68) Atrazine                   | 16.49 | 200  | 263822   | 20.601 | ng/ul | 99       |
| 69) Pentachlorophenol          | 16.67 | 266  | 196543   | 20.226 | ng/ul | 99       |
| 70) Phenanthrene               | 17.06 | 178  | 1354256  | 20.938 | ng/ul | 100      |
| 72) Anthracene                 | 17.14 | 178  | 1404700  | 21.311 | ng/ul | 99       |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.06 | 216  | 363715   | 21.601 | ng/uL | 98       |
| 74) Pentachlorobenzene         | 14.60 | 250  | 397920   | 21.231 | ng/uL | 99       |
| 75) Carbazole                  | 17.42 | 167  | 1205452  | 20.926 | ng/ul | 99       |
| 76) Di-n-butylphthalate        | 18.00 | 149  | 1478316  | 21.980 | ng/ul | 99       |
| 77) Fluoranthene               | 19.07 | 202  | 1551291  | 20.348 | ng/ul | 99       |
| 80) Pvrene                     | 19.44 | 202  | 1588158  | 21.741 | ng/ul | 98       |
| 81) Butylbenzylphthalate       | 20.36 | 149  | 651283   | 23.410 | ng/ul | 100      |
| 82) 3,3'-Dichlorobenzidine     | 21.13 | 252  | 445705   | 18.034 | ng/ul | 99       |
| 83) Benzo(a)anthracene         | 21.19 | 228  | 1527784  | 21.305 | ng/ul | 100      |
| 84) Bis(2-ethylhexyl)phthalate | 21.14 | 149  | 980539   | 22.721 | ng/ul | 100      |
| 85) Chrysene                   | 21.24 | 228  | 1470440  | 20.900 | ng/ul | 100      |
| 87) Di-n-octyl phthalate       | 22.01 | 149  | 1676330  | 22.362 | ng/ul | 100      |
| 88) Benzo(b)fluoranthene       | 22.74 | 252  | 1613633  | 20.672 | ng/ul | 100      |
| 89) Benzo(k)fluoranthene       | 22.78 | 252  | 1629175  | 21.187 | ng/ul | 100      |
| 91) Benzo(a)pyrene             | 23.29 | 252  | 1497340  | 21.079 | ng/ul | 99       |
| 92) Indeno(1,2,3-cd)pyrene     | 25.55 | 276  | 1926677  | 20.538 | ng/ul | 99       |
| 93) Dibenzo(a,h)anthracene     | 25.56 | 278  | 1621134  | 20.407 | ng/ul | 99       |
| 94) Benzo(a,h,i)perylene       | 26.21 | 276  | 1589425  | 20.557 | ng/ul | 99       |

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031620\  
Data File : BM025584.D  
Acq On : 16 Mar 2020 17:32  
Operator : CG/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD02009

Quant Time: Mar 16 18:11:32 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM031420MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Mar 16 04:27:51 2020  
Response via : Initial Calibration

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

