

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG091423\  
 Data File : BG059062.D  
 Acq On : 14 Sep 2023 13:07  
 Operator : MA/JU  
 Sample : SSTD02015  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD020416

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 09/15/2023  
 Supervised By :mohammad ahmed 09/15/2023

Quant Time: Sep 15 05:33:31 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG091423.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Sep 14 15:06:26 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.299  | 152  | 85705    | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 11.148 | 136  | 410772   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.932 | 164  | 325914   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.682 | 188  | 889832   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 22.006 | 240  | 838259   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 25.555 | 264  | 952376   | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.604  | 96   | 16369    | 8.255  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 4.039  | 84   | 128711   | 20.226 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.412  | 99   | 176790   | 20.479 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.617  | 67   | 104586   | 20.614 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.811  | 132  | 125772   | 22.533 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.986  | 113  | 148805   | 21.896 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.497  | 128  | 65536    | 21.556 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.220 | 143  | 84854    | 24.347 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.737 | 165  | 157523   | 22.018 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.284 | 131  | 220434   | 22.823 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.327 | 166  | 579903   | 22.464 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.633 | 160  | 632219   | 22.098 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 15.103 | 143  | 105674   | 23.997 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.919 | 176  | 534070   | 23.066 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 16.025 | 200  | 113286   | 24.259 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.782 | 188  | 869352   | 21.664 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 20.056 | 212  | 1007038  | 22.088 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 25.308 | 264  | 1019622  | 21.575 | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.645  | 88   | 20099    | 7.830  | ng/uL  | 100      |
| 5) Pyridine                   | 4.063  | 79   | 133480   | 19.779 | ng/u1  | 100      |
| 6) Benzaldehyde               | 7.441  | 77   | 116431   | 21.737 | ng/u1  | 100      |
| 8) Phenol                     | 7.435  | 94   | 189817   | 20.497 | ng/u1  | 100      |
| 10) Bis(2-Chloroethyl)ether   | 7.717  | 93   | 138177   | 20.680 | ng/u1  | 100      |
| 12) 2-Chlorophenol            | 7.846  | 128  | 118781   | 20.806 | ng/u1  | 100      |
| 13) 2-Methylphenol            | 8.716  | 108  | 149219   | 21.181 | ng/u1  | 100      |
| 14) 2,2'-oxybis(1-Chloropr... | 8.816  | 45   | 166417   | 20.543 | ng/u1  | 100      |
| 16) Acetophenone              | 9.139  | 105  | 249859   | 22.155 | ng/u1  | 100      |
| 17) N-Nitroso-di-n-propyla... | 9.104  | 70   | 139350   | 21.945 | ng/u1  | 100      |
| 18) 4-Methylphenol            | 9.045  | 108  | 168716   | 21.898 | ng/u1  | 100      |
| 19) Hexachloroethane          | 9.386  | 117  | 51235    | 21.057 | ng/u1  | 100      |
| 22) Nitrobenzene              | 9.538  | 77   | 196234   | 21.570 | ng/u1  | 100      |
| 23) Isophorone                | 10.050 | 82   | 425407   | 21.909 | ng/u1  | 100      |
| 25) 2-Nitrophenol             | 10.249 | 139  | 79452    | 23.055 | ng/u1  | 100      |
| 26) 2,4-Dimethylphenol        | 10.267 | 107  | 189891   | 22.311 | ng/u1  | 100      |
| 27) Bis(2-Chloroethoxy)met... | 10.531 | 93   | 212011   | 20.822 | ng/u1  | 100      |
| 29) 2,4-Dichlorophenol        | 10.766 | 162  | 151741   | 22.466 | ng/u1  | 100      |
| 30) Naphthalene               | 11.195 | 128  | 474125   | 21.661 | ng/u1  | 100      |
| 32) 4-Chloroaniline           | 11.313 | 127  | 206822   | 22.732 | ng/u1  | 100      |
| 33) Hexachlorobutadiene       | 11.430 | 225  | 116001   | 22.198 | ng/u1  | 100      |
| 34) Caprolactam               | 12.088 | 113  | 58347    | 23.860 | ng/u1  | 100      |
| 35) 4-Chloro-3-methylphenol   | 12.382 | 107  | 198146   | 23.692 | ng/u1  | 100      |
| 36) 2-Methylnaphthalene       | 12.776 | 142  | 364415   | 22.011 | ng/u1  | 100      |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG091423\  
 Data File : BG059062.D  
 Acq On : 14 Sep 2023 13:07  
 Operator : MA/JU  
 Sample : SST02015  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

## Instrument :

BNA\_G

## ClientSampleId :

SSTD020416

## Manual Integrations

## APPROVED

Reviewed By :Yogesh Patel 09/15/2023

Supervised By :mohammad ahmed 09/15/2023

Quant Time: Sep 15 05:33:31 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG091423.MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Sep 14 15:06:26 2023

Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 37) 1-Methylnaphthalene       | 12.993 | 142  | 368646   | 22.563 | ng/ul | 100      |
| 39) 1,2,4,5-Tetrachloroben... | 13.117 | 216  | 226069   | 21.177 | ng/ul | 100      |
| 40) Hexachlorocyclopentadiene | 13.070 | 237  | 161051   | 19.736 | ng/ul | 100      |
| 41) 2,4,6-Trichlorophenol     | 13.358 | 196  | 148509   | 21.796 | ng/ul | 100      |
| 42) 2,4,5-Trichlorophenol     | 13.428 | 196  | 164686   | 22.103 | ng/ul | 100      |
| 43) 1,1'-Biphenyl             | 13.769 | 154  | 530529   | 21.564 | ng/ul | 100      |
| 44) 2-Chloronaphthalene       | 13.822 | 162  | 400567   | 21.454 | ng/ul | 100      |
| 45) 2-Nitroaniline            | 14.033 | 65   | 140025   | 23.572 | ng/ul | 100      |
| 47) Dimethylphthalate         | 14.374 | 163  | 559121   | 22.759 | ng/ul | 100      |
| 48) 2,6-Dinitrotoluene        | 14.515 | 165  | 111531   | 23.780 | ng/ul | 100      |
| 50) Acenaphthylene            | 14.662 | 152  | 681353   | 21.937 | ng/ul | 100      |
| 51) 3-Nitroaniline            | 14.850 | 138  | 111176   | 24.016 | ng/ul | 100      |
| 52) Acenaphthene              | 14.991 | 153  | 452662   | 21.682 | ng/ul | 100      |
| 53) 2,4-Dinitrophenol         | 15.044 | 184  | 64389    | 20.640 | ng/ul | 100      |
| 55) 4-Nitrophenol             | 15.120 | 109  | 123243   | 23.673 | ng/ul | 100      |
| 56) Dibenzofuran              | 15.326 | 168  | 672850   | 22.114 | ng/ul | 100      |
| 57) 2,4-Dinitrotoluene        | 15.291 | 165  | 170036   | 25.483 | ng/ul | 100      |
| 58) 2,3,4,6-Tetrachlorophenol | 15.531 | 232  | 180252   | 24.587 | ng/ul | 100      |
| 59) Diethylphthalate          | 15.719 | 149  | 570810   | 23.352 | ng/ul | 100      |
| 61) Fluorene                  | 15.972 | 166  | 567562   | 22.701 | ng/ul | 100      |
| 62) 4-Chlorophenyl-phenyle... | 15.954 | 204  | 308501   | 23.020 | ng/ul | 100      |
| 63) 4-Nitroaniline            | 16.013 | 138  | 105233m  | 23.946 | ng/ul |          |
| 66) 4,6-Dinitro-2-methylph... | 16.037 | 198  | 118055   | 23.398 | ng/ul | 100      |
| 67) N-Nitrosodiphenylamine    | 16.178 | 169  | 484800   | 20.645 | ng/ul | 100      |
| 68) 4-Bromophenyl-phenylether | 16.859 | 248  | 219980   | 22.039 | ng/ul | 100      |
| 69) Hexachlorobenzene         | 16.947 | 284  | 241815   | 21.438 | ng/ul | 100      |
| 70) Atrazine                  | 17.106 | 200  | 225592   | 23.901 | ng/ul | 100      |
| 71) Pentachlorophenol         | 17.300 | 266  | 152062   | 21.445 | ng/ul | 100      |
| 72) Phenanthrene              | 17.723 | 178  | 971328   | 21.246 | ng/ul | 100      |
| 74) Anthracene                | 17.817 | 178  | 1007843  | 21.608 | ng/ul | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 13.728 | 216  | 252624   | 19.898 | ng/ul | 100      |
| 76) Pentachlorobenzene        | 15.220 | 250  | 261987   | 21.190 | ng/ul | 100      |
| 77) Carbazole                 | 18.087 | 167  | 847039   | 21.810 | ng/ul | 100      |
| 78) Di-n-butylphthalate       | 18.598 | 149  | 974844   | 22.276 | ng/ul | 100      |
| 80) Fluoranthene              | 19.715 | 202  | 1172929  | 22.090 | ng/ul | 100      |
| 82) Pyrene                    | 20.085 | 202  | 1205311  | 21.886 | ng/ul | 100      |
| 83) Butylbenzylphthalate      | 20.943 | 149  | 419434   | 23.752 | ng/ul | 100      |
| 84) 3,3'-Dichlorobenzidine    | 21.895 | 252  | 432444   | 22.683 | ng/ul | 100      |
| 85) Benzo(a)anthracene        | 21.983 | 228  | 1220619  | 21.428 | ng/ul | 100      |
| 86) Bis(2-ethylhexyl)phtha... | 21.818 | 149  | 612714   | 23.232 | ng/ul | 100      |
| 87) Chrysene                  | 22.053 | 228  | 1148525  | 21.606 | ng/ul | 100      |
| 89) Di-n-octyl phthalate      | 23.146 | 149  | 1037472  | 23.012 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 24.403 | 252  | 1208440  | 21.145 | ng/ul | 100      |
| 91) Benzo(k)fluoranthene      | 24.486 | 252  | 1220427  | 21.026 | ng/ul | 100      |
| 93) Benzo(a)pyrene            | 25.390 | 252  | 1129975  | 21.057 | ng/ul | 100      |
| 94) Indeno(1,2,3-cd)pyrene    | 29.668 | 276  | 1412598m | 20.909 | ng/ul |          |
| 95) Dibenzo(a,h)anthracene    | 29.756 | 278  | 1119792  | 20.083 | ng/ul | 100      |
| 96) Benzo(g,h,i)perylene      | 30.984 | 276  | 1140618  | 21.320 | ng/ul | 100      |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG091423\  
 Data File : BG059062.D  
 Acq On : 14 Sep 2023 13:07  
 Operator : MA/JU  
 Sample : SSTD02015  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

## Instrument :

BNA\_G

## ClientSampleId :

SSTD020416

## Manual Integrations

## APPROVED

Reviewed By :Yogesh Patel 09/15/2023

Supervised By :mohammad ahmed 09/15/2023

Quant Time: Sep 15 05:33:31 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG091423.MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Sep 14 15:06:26 2023

Response via : Initial Calibration

