

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070523\  
 Data File : BG058054.D  
 Acq On : 6 Jul 2023 17:44  
 Operator : CG/JU  
 Sample : 03258-10MSD  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 DCGW7MSD

Quant Time: Jul 06 22:50:06 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG070123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jul 05 23:20:13 2023  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Yogesh Patel 07/07/2023  
 Supervised By :mohammad ahmed 07/07/2023

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.921  | 152  | 39478    | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.724 | 136  | 178869   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.554 | 164  | 129774   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.298 | 188  | 322476   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.558 | 240  | 291365   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.707 | 264  | 367412   | 20.000 | ng/u1  | -0.01    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.362  | 96   | 4518     | 4.007  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.767  | 84   | 27867    | 8.142  | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.081  | 99   | 21109    | 5.236  | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.245  | 67   | 60681    | 24.975 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.451  | 132  | 51615    | 18.562 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.626  | 113  | 38474    | 12.647 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.084  | 128  | 36916    | 25.322 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.801  | 143  | 36421    | 23.770 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.342 | 165  | 64724    | 22.020 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.859 | 131  | 90306    | 20.264 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.961 | 166  | 277082   | 27.061 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.249 | 160  | 290344   | 25.985 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.760 | 143  | 8767     | 4.846  | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.547 | 176  | 232816   | 26.773 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.665 | 200  | 37372    | 21.343 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.398 | 188  | 405064   | 27.070 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.690 | 212  | 499357   | 27.212 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.496 | 264  | 522644   | 27.921 | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.391  | 88   | 4102     | 2.945  | ng/uL# | 71       |
| 5) Pyridine                   | 3.791  | 79   | 26398    | 7.527  | ng/u1  | 96       |
| 6) Benzaldehyde               | 7.057  | 77   | 58606    | 43.527 | ng/u1  | 97       |
| 8) Phenol                     | 7.110  | 94   | 22760    | 5.563  | ng/u1  | 96       |
| 10) Bis(2-Chloroethyl)ether   | 7.339  | 93   | 71246    | 22.553 | ng/u1  | 99       |
| 12) 2-Chlorophenol            | 7.486  | 128  | 50132    | 17.325 | ng/u1  | 96       |
| 13) 2-Methylphenol            | 8.362  | 108  | 44935    | 13.985 | ng/u1  | 100      |
| 14) 2,2'-oxybis(1-Chloropr... | 8.444  | 45   | 116307m  | 23.918 | ng/u1  |          |
| 16) Acetophenone              | 8.738  | 105  | 120200   | 23.815 | ng/u1  | 100      |
| 17) N-Nitroso-di-n-propyla... | 8.720  | 70   | 64226    | 23.803 | ng/u1  | 98       |
| 18) 4-Methylphenol            | 8.685  | 108  | 42741    | 12.302 | ng/u1  | 98       |
| 19) Hexachloroethane          | 9.002  | 117  | 18987    | 17.117 | ng/u1  | 95       |
| 22) Nitrobenzene              | 9.120  | 77   | 92264    | 24.379 | ng/u1  | 98       |
| 23) Isophorone                | 9.643  | 82   | 190823   | 23.630 | ng/u1  | 99       |
| 25) 2-Nitrophenol             | 9.836  | 139  | 34617    | 21.055 | ng/u1# | 92       |
| 26) 2,4-Dimethylphenol        | 9.895  | 107  | 56789    | 15.361 | ng/u1  | 96       |
| 27) Bis(2-Chloroethoxy)met... | 10.124 | 93   | 107233   | 23.564 | ng/u1  | 96       |
| 29) 2,4-Dichlorophenol        | 10.371 | 162  | 58784    | 20.570 | ng/u1  | 89       |
| 30) Naphthalene               | 10.771 | 128  | 214097   | 21.299 | ng/u1  | 100      |
| 32) 4-Chloroaniline           | 10.882 | 127  | 78095    | 17.139 | ng/u1  | 100      |
| 33) Hexachlorobutadiene       | 11.053 | 225  | 36763    | 18.170 | ng/u1  | 98       |
| 34) Caprolactam               | 11.629 | 113  | 4038     | 3.514  | ng/u1  | 92       |
| 35) 4-Chloro-3-methylphenol   | 12.005 | 107  | 72580    | 20.675 | ng/u1  | 98       |
| 36) 2-Methylnaphthalene       | 12.381 | 142  | 150011   | 22.138 | ng/u1  | 97       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070523\  
 Data File : BG058054.D  
 Acq On : 6 Jul 2023 17:44  
 Operator : CG/JU  
 Sample : 03258-10MSD  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

## Instrument :

BNA\_G

## ClientSampleId :

DCGW7MSD

## Manual Integrations

## APPROVED

Reviewed By :Yogesh Patel 07/07/2023

Supervised By :mohammad ahmed 07/07/2023

Quant Time: Jul 06 22:50:06 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG070123.MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Jul 05 23:20:13 2023

Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 37) 1-Methylnaphthalene       | 12.604 | 142  | 156136   | 22.978 | ng/ul  | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 12.751 | 216  | 89828    | 22.714 | ng/ul  | 99       |
| 40) Hexachlorocyclopentadiene | 12.721 | 237  | 33156    | 13.310 | ng/ul  | 98       |
| 41) 2,4,6-Trichlorophenol     | 12.992 | 196  | 54113    | 20.188 | ng/ul  | 96       |
| 42) 2,4,5-Trichlorophenol     | 13.062 | 196  | 61593    | 21.248 | ng/ul  | 100      |
| 43) 1,1'-Biphenyl             | 13.391 | 154  | 218310   | 23.343 | ng/ul  | 98       |
| 44) 2-Chloronaphthalene       | 13.432 | 162  | 171406   | 22.837 | ng/ul  | 100      |
| 45) 2-Nitroaniline            | 13.638 | 65   | 65111    | 24.951 | ng/ul  | 94       |
| 47) Dimethylphthalate         | 14.008 | 163  | 257117   | 24.994 | ng/ul  | 99       |
| 48) 2,6-Dinitrotoluene        | 14.131 | 165  | 55090    | 26.089 | ng/ul  | 96       |
| 50) Acenaphthylene            | 14.278 | 152  | 289210   | 23.969 | ng/ul  | 99       |
| 51) 3-Nitroaniline            | 14.460 | 138  | 49275    | 27.613 | ng/ul  | 99       |
| 52) Acenaphthene              | 14.619 | 153  | 200790   | 24.251 | ng/ul  | 97       |
| 53) 2,4-Dinitrophenol         | 14.666 | 184  | 18588    | 16.885 | ng/ul# | 89       |
| 55) 4-Nitrophenol             | 14.778 | 109  | 6014     | 4.488  | ng/ul# | 74       |
| 56) Dibenzofuran              | 14.954 | 168  | 291187   | 24.740 | ng/ul  | 99       |
| 57) 2,4-Dinitrotoluene        | 14.919 | 165  | 82174    | 27.467 | ng/ul# | 98       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.183 | 232  | 54517    | 20.659 | ng/ul# | 99       |
| 59) Diethylphthalate          | 15.365 | 149  | 270136   | 25.160 | ng/ul  | 99       |
| 61) Fluorene                  | 15.600 | 166  | 244357   | 25.303 | ng/ul  | 98       |
| 62) 4-Chlorophenyl-phenyle... | 15.594 | 204  | 130099   | 25.461 | ng/ul  | 99       |
| 63) 4-Nitroaniline            | 15.624 | 138  | 50419    | 31.540 | ng/ul  | 96       |
| 66) 4,6-Dinitro-2-methylph... | 15.683 | 198  | 34941    | 19.512 | ng/ul# | 98       |
| 67) N-Nitrosodiphenylamine    | 15.806 | 169  | 184549   | 21.504 | ng/ul  | 96       |
| 68) 4-Bromophenyl-phenylether | 16.488 | 248  | 92523    | 26.138 | ng/ul  | 96       |
| 69) Hexachlorobenzene         | 16.605 | 284  | 106106   | 26.773 | ng/ul  | 99       |
| 70) Atrazine                  | 16.752 | 200  | 69747    | 19.874 | ng/ul  | 97       |
| 71) Pentachlorophenol         | 16.952 | 266  | 45818    | 19.257 | ng/ul  | 96       |
| 72) Phenanthrene              | 17.345 | 178  | 417668   | 25.462 | ng/ul  | 99       |
| 74) Anthracene                | 17.433 | 178  | 418383   | 25.205 | ng/ul  | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 13.356 | 216  | 95260    | 22.274 | ng/uL  | 99       |
| 76) Pentachlorobenzene        | 14.872 | 250  | 112168   | 24.557 | ng/uL  | 98       |
| 77) Carbazole                 | 17.704 | 167  | 397311   | 26.465 | ng/ul  | 99       |
| 78) Di-n-butylphthalate       | 18.256 | 149  | 499763   | 26.133 | ng/ul  | 99       |
| 80) Fluoranthene              | 19.355 | 202  | 532781   | 25.056 | ng/ul  | 98       |
| 82) Pyrene                    | 19.719 | 202  | 533001   | 24.902 | ng/ul  | 99       |
| 83) Butylbenzylphthalate      | 20.600 | 149  | 223632   | 25.766 | ng/ul  | 98       |
| 84) 3,3'-Dichlorobenzidine    | 21.452 | 252  | 120401   | 17.069 | ng/ul  | 96       |
| 85) Benzo(a)anthracene        | 21.535 | 228  | 544113   | 26.009 | ng/ul  | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.441 | 149  | 327988   | 26.370 | ng/ul  | 96       |
| 87) Chrysene                  | 21.605 | 228  | 516532   | 26.338 | ng/ul  | 100      |
| 89) Di-n-octyl phthalate      | 22.621 | 149  | 568934   | 25.202 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.708 | 252  | 570695   | 25.086 | ng/ul  | 98       |
| 91) Benzo(k)fluoranthene      | 23.773 | 252  | 575279   | 25.799 | ng/ul  | 99       |
| 93) Benzo(a)pyrene            | 24.560 | 252  | 507383   | 25.199 | ng/ul  | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.303 | 276  | 683979   | 25.581 | ng/ul  | 99       |
| 95) Dibenzo(a,h)anthracene    | 28.368 | 278  | 573504   | 26.011 | ng/ul  | 98       |
| 96) Benzo(g,h,i)perylene      | 29.431 | 276  | 558397   | 25.568 | ng/ul  | 99       |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070523\  
 Data File : BG058054.D  
 Acq On : 6 Jul 2023 17:44  
 Operator : CG/JU  
 Sample : 03258-10MSD  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

## Instrument :

BNA\_G

## ClientSampleId :

DCGW7MSD

## Manual Integrations

## APPROVED

Reviewed By :Yogesh Patel 07/07/2023

Supervised By :mohammad ahmed 07/07/2023

Quant Time: Jul 06 22:50:06 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG070123.MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Jul 05 23:20:13 2023

Response via : Initial Calibration

