

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG071224\  
 Data File : BG062045.D  
 Acq On : 12 Jul 2024 17:58  
 Operator : MA/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

## Instrument :

BNA\_G

## ClientSampleId :

SSTD020570

## Manual Integrations

## APPROVED

Reviewed By :Yogesh Patel 07/13/2024

Supervised By :mohammad ahmed 07/16/2024

Supervised By :mohammad

ahmed

07/16/2024

Quant Time: Jul 12 19:17:56 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG070224.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jul 11 12:11:32 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.038  | 152  | 45130    | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.852 | 136  | 203553   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.672 | 164  | 138856   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.421 | 188  | 342427   | 20.000 | ng/u1  | # 0.00   |
| 79) Chrysene-d12              | 21.681 | 240  | 368741   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.889 | 264  | 384534   | 20.000 | ng/u1  | 0.00     |
|                               |        |      |          |        |        |          |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.438  | 96   | 8863     | 7.380  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.855  | 84   | 63632    | 17.232 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.204  | 99   | 72789    | 14.738 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.362  | 67   | 41948    | 13.412 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.574  | 132  | 57527    | 16.981 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.749  | 113  | 65985    | 16.969 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.207  | 128  | 32832    | 21.208 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.930  | 143  | 38155    | 21.582 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.471 | 165  | 75542    | 22.153 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.988 | 131  | 94814    | 19.152 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.072 | 166  | 212770   | 18.638 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.366 | 160  | 232538   | 19.481 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.877 | 143  | 33194    | 18.208 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.664 | 176  | 187520   | 19.512 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.782 | 200  | 37982    | 21.120 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.515 | 188  | 313662   | 19.619 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.801 | 212  | 346192   | 15.962 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.666 | 264  | 361031   | 18.931 | ng/u1  | 0.00     |
|                               |        |      |          |        |        |          |
| Target Compounds              |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.473  | 88   | 8761m    | 6.609  | ng/uL  |          |
| 5) Pyridine                   | 3.878  | 79   | 66339    | 17.517 | ng/u1  | 95       |
| 6) Benzaldehyde               | 7.180  | 77   | 44973    | 17.205 | ng/u1  | 90       |
| 8) Phenol                     | 7.227  | 94   | 74946    | 14.822 | ng/u1  | 91       |
| 10) Bis(2-Chloroethyl)ether   | 7.456  | 93   | 60733    | 15.635 | ng/u1  | 95       |
| 12) 2-Chlorophenol            | 7.603  | 128  | 58221    | 16.869 | ng/u1# | 86       |
| 13) 2-Methylphenol            | 8.479  | 108  | 63916    | 17.710 | ng/u1  | 87       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.567  | 45   | 132794   | 15.648 | ng/u1# | 96       |
| 16) Acetophenone              | 8.867  | 105  | 112261   | 17.477 | ng/u1  | 94       |
| 17) N-Nitroso-di-n-propyla... | 8.843  | 70   | 59741    | 16.613 | ng/u1# | 85       |
| 18) 4-Methylphenol            | 8.814  | 108  | 69582    | 17.216 | ng/u1  | 94       |
| 19) Hexachloroethane          | 9.125  | 117  | 27912    | 18.642 | ng/u1# | 72       |
| 22) Nitrobenzene              | 9.248  | 77   | 91427    | 20.805 | ng/u1  | 98       |
| 23) Isophorone                | 9.765  | 82   | 176840   | 17.813 | ng/u1# | 94       |
| 25) 2-Nitrophenol             | 9.959  | 139  | 38155    | 21.163 | ng/u1# | 91       |
| 26) 2,4-Dimethylphenol        | 10.018 | 107  | 88397    | 20.487 | ng/u1  | 96       |
| 27) Bis(2-Chloroethoxy)met... | 10.253 | 93   | 93083    | 17.026 | ng/u1  | 98       |
| 29) 2,4-Dichlorophenol        | 10.500 | 162  | 65796    | 20.541 | ng/u1  | 96       |
| 30) Naphthalene               | 10.905 | 128  | 221312   | 19.793 | ng/u1# | 94       |
| 32) 4-Chloroaniline           | 11.011 | 127  | 93976    | 19.346 | ng/u1  | 98       |
| 33) Hexachlorobutadiene       | 11.176 | 225  | 52354    | 21.720 | ng/u1# | 86       |
| 34) Caprolactam               | 11.775 | 113  | 24927m   | 19.863 | ng/u1  |          |
| 35) 4-Chloro-3-methylphenol   | 12.133 | 107  | 89730    | 21.034 | ng/u1  | 94       |
| 36) 2-Methylnaphthalene       | 12.503 | 142  | 158155   | 19.524 | ng/u1  | 98       |

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 Data File : BG062045.D  
 Acq On : 12 Jul 2024 17:58  
 Operator : MA/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

## Instrument :

BNA\_G

## ClientSampleId :

SSTD020570

## Manual Integrations

## APPROVED

Reviewed By :Yogesh Patel 07/13/2024

Supervised By :mohammad ahmed 07/16/2024

Supervised By :mohammad

ahmed

07/16/2024

Quant Time: Jul 12 19:17:56 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG070224.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jul 11 12:11:32 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 37) 1-Methylnaphthalene       | 12.721 | 142  | 170259   | 20.225 | ng/ul  | 90       |
| 39) 1,2,4,5-Tetrachloroben... | 12.868 | 216  | 96159    | 23.060 | ng/ul# | 97       |
| 40) Hexachlorocyclopentadiene | 12.838 | 237  | 55404    | 19.959 | ng/ul  | 96       |
| 41) 2,4,6-Trichlorophenol     | 13.109 | 196  | 62961    | 22.733 | ng/ul  | 99       |
| 42) 2,4,5-Trichlorophenol     | 13.185 | 196  | 70802    | 23.271 | ng/ul  | 89       |
| 43) 1,1'-Biphenyl             | 13.508 | 154  | 224329   | 20.968 | ng/ul  | 99       |
| 44) 2-Chloronaphthalene       | 13.555 | 162  | 172160   | 21.441 | ng/ul  | 95       |
| 45) 2-Nitroaniline            | 13.755 | 65   | 79595    | 24.754 | ng/ul  | 90       |
| 47) Dimethylphthalate         | 14.119 | 163  | 202680   | 18.458 | ng/ul  | 99       |
| 48) 2,6-Dinitrotoluene        | 14.243 | 165  | 44351    | 21.291 | ng/ul  | 98       |
| 50) Acenaphthylene            | 14.395 | 152  | 250094   | 19.089 | ng/ul  | 98       |
| 51) 3-Nitroaniline            | 14.583 | 138  | 42701    | 21.059 | ng/ul# | 98       |
| 52) Acenaphthene              | 14.736 | 153  | 174796   | 19.322 | ng/ul  | 98       |
| 53) 2,4-Dinitrophenol         | 14.789 | 184  | 29802    | 27.385 | ng/ul  | 93       |
| 55) 4-Nitrophenol             | 14.895 | 109  | 45535    | 21.884 | ng/ul  | 91       |
| 56) Dibenzofuran              | 15.071 | 168  | 249657   | 19.526 | ng/ul  | 94       |
| 57) 2,4-Dinitrotoluene        | 15.036 | 165  | 68190    | 22.322 | ng/ul# | 91       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.294 | 232  | 56140    | 20.435 | ng/ul# | 95       |
| 59) Diethylphthalate          | 15.482 | 149  | 224433   | 18.968 | ng/ul  | 95       |
| 61) Fluorene                  | 15.717 | 166  | 207550   | 19.088 | ng/ul  | 99       |
| 62) 4-Chlorophenyl-phenyle... | 15.706 | 204  | 109565   | 20.061 | ng/ul  | 91       |
| 63) 4-Nitroaniline            | 15.741 | 138  | 43539    | 21.162 | ng/ul  | 82       |
| 66) 4,6-Dinitro-2-methylph... | 15.800 | 198  | 38869    | 20.083 | ng/ul# | 95       |
| 67) N-Nitrosodiphenylamine    | 15.923 | 169  | 174920   | 18.705 | ng/ul  | 97       |
| 68) 4-Bromophenyl-phenylether | 16.605 | 248  | 74297    | 19.125 | ng/ul  | 95       |
| 69) Hexachlorobenzene         | 16.716 | 284  | 91314    | 20.442 | ng/ul  | 96       |
| 70) Atrazine                  | 16.869 | 200  | 75337    | 20.408 | ng/ul  | 98       |
| 71) Pentachlorophenol         | 17.063 | 266  | 50588    | 20.603 | ng/ul  | 92       |
| 72) Phenanthrene              | 17.462 | 178  | 344919   | 19.103 | ng/ul  | 100      |
| 74) Anthracene                | 17.550 | 178  | 357956   | 19.184 | ng/ul  | 97       |
| 75) 1,2,3,4-Tetrachloroben... | 13.473 | 216  | 93250    | 21.926 | ng/ul  | 90       |
| 76) Pentachlorobenzene        | 14.989 | 250  | 100612   | 20.742 | ng/ul  | 96       |
| 77) Carbazole                 | 17.821 | 167  | 307519   | 19.583 | ng/ul  | 100      |
| 78) Di-n-butylphthalate       | 18.367 | 149  | 381795   | 18.937 | ng/ul  | 98       |
| 80) Fluoranthene              | 19.466 | 202  | 413485   | 16.109 | ng/ul# | 94       |
| 82) Pyrene                    | 19.830 | 202  | 415890   | 15.686 | ng/ul# | 91       |
| 83) Butylbenzylphthalate      | 20.706 | 149  | 204265   | 18.357 | ng/ul  | 91       |
| 84) 3,3'-Dichlorobenzidine    | 21.575 | 252  | 191256   | 21.802 | ng/ul  | 93       |
| 85) Benzo(a)anthracene        | 21.657 | 228  | 509188   | 19.072 | ng/ul  | 98       |
| 86) Bis(2-ethylhexyl)phtha... | 21.552 | 149  | 294566   | 18.448 | ng/ul  | 98       |
| 87) Chrysene                  | 21.722 | 228  | 467184   | 18.692 | ng/ul  | 99       |
| 89) Di-n-octyl phthalate      | 22.762 | 149  | 507931   | 21.621 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.867 | 252  | 455136m  | 18.786 | ng/ul  |          |
| 91) Benzo(k)fluoranthene      | 23.931 | 252  | 457097   | 18.920 | ng/ul# | 92       |
| 93) Benzo(a)pyrene            | 24.742 | 252  | 435492   | 18.983 | ng/ul# | 97       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.549 | 276  | 650966   | 22.204 | ng/ul# | 93       |
| 95) Dibenzo(a,h)anthracene    | 28.614 | 278  | 529109   | 21.852 | ng/ul  | 98       |
| 96) Benzo(g,h,i)perylene      | 29.689 | 276  | 511624   | 21.936 | ng/ul# | 93       |

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

