

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG100923\  
 Data File : BG059415.D  
 Acq On : 10 Oct 2023 3:27  
 Operator : MA/JU  
 Sample : PB156170BS  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SLCS170

Quant Time: Oct 10 05:38:08 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG100723.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Oct 07 22:45:22 2023  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Yogesh Patel 10/11/2023  
 Supervised By :mohammad ahmed 10/11/2023

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.230  | 152  | 37053    | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 11.068 | 136  | 181990   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.863 | 164  | 123197   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.607 | 188  | 273898   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.914 | 240  | 240580   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 25.392 | 264  | 279029   | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.530  | 96   | 8823m    | 8.490  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.959  | 84   | 115793   | 38.711 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.355  | 99   | 168340   | 41.146 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.548  | 67   | 94761    | 40.978 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.748  | 132  | 123088   | 43.017 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.923  | 113  | 129012   | 40.472 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.429  | 128  | 65762    | 42.448 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.145 | 143  | 74139    | 44.215 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.674 | 165  | 134352   | 44.392 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.215 | 131  | 191256   | 39.783 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.258 | 166  | 414823   | 41.100 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.564 | 160  | 501720   | 41.771 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 15.057 | 143  | 76516    | 40.455 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.850 | 176  | 384625   | 39.923 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.968 | 200  | 74355    | 41.117 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.707 | 188  | 589066   | 43.842 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.987 | 212  | 628784   | 43.144 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 25.151 | 264  | 645105   | 43.935 | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.565  | 88   | 18458m   | 16.638 | ng/uL  |          |
| 5) Pyridine                   | 3.988  | 79   | 103104   | 34.139 | ng/u1  | 95       |
| 6) Benzaldehyde               | 7.372  | 77   | 76661    | 45.172 | ng/u1  | 93       |
| 8) Phenol                     | 7.384  | 94   | 170402   | 42.119 | ng/u1  | 97       |
| 10) Bis(2-Chloroethyl)ether   | 7.642  | 93   | 130912   | 42.017 | ng/u1  | 95       |
| 12) 2-Chlorophenol            | 7.778  | 128  | 128861   | 42.738 | ng/u1  | 95       |
| 13) 2-Methylphenol            | 8.653  | 108  | 125316   | 41.453 | ng/u1  | 98       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.729  | 45   | 269429m  | 41.330 | ng/u1  |          |
| 16) Acetophenone              | 9.070  | 105  | 196411   | 40.547 | ng/u1  | 95       |
| 17) N-Nitroso-di-n-propyla... | 9.035  | 70   | 94511    | 39.157 | ng/u1  | 92       |
| 18) 4-Methylphenol            | 8.988  | 108  | 136711   | 41.124 | ng/u1  | 95       |
| 19) Hexachloroethane          | 9.311  | 117  | 48737    | 44.439 | ng/u1# | 83       |
| 22) Nitrobenzene              | 9.470  | 77   | 139038   | 40.938 | ng/u1# | 97       |
| 23) Isophorone                | 9.975  | 82   | 299071   | 41.238 | ng/u1  | 98       |
| 25) 2-Nitrophenol             | 10.175 | 139  | 74891    | 43.306 | ng/u1# | 91       |
| 26) 2,4-Dimethylphenol        | 10.204 | 107  | 130504   | 38.255 | ng/u1  | 95       |
| 27) Bis(2-Chloroethoxy)met... | 10.457 | 93   | 183692   | 42.150 | ng/u1  | 99       |
| 29) 2,4-Dichlorophenol        | 10.704 | 162  | 136189   | 44.687 | ng/u1  | 95       |
| 30) Naphthalene               | 11.121 | 128  | 446770   | 42.446 | ng/u1  | 98       |
| 32) 4-Chloroaniline           | 11.238 | 127  | 175565   | 38.603 | ng/u1  | 99       |
| 33) Hexachlorobutadiene       | 11.350 | 225  | 82605    | 44.891 | ng/u1  | 93       |
| 34) Caprolactam               | 12.026 | 113  | 44029m   | 39.785 | ng/u1  |          |
| 35) 4-Chloro-3-methylphenol   | 12.319 | 107  | 139757   | 42.736 | ng/u1  | 96       |
| 36) 2-Methylnaphthalene       | 12.707 | 142  | 300656   | 42.847 | ng/u1  | 94       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG100923\  
 Data File : BG059415.D  
 Acq On : 10 Oct 2023 3:27  
 Operator : MA/JU  
 Sample : PB156170BS  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SLCS170

Quant Time: Oct 10 05:38:08 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG100723.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Oct 07 22:45:22 2023  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Yogesh Patel 10/11/2023  
 Supervised By :mohammad ahmed 10/11/2023

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 37) 1-Methylnaphthalene       | 12.924 | 142  | 300510   | 41.248 | ng/ul  | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 13.048 | 216  | 161348   | 43.968 | ng/ul  | 97       |
| 40) Hexachlorocyclopentadiene | 13.001 | 237  | 86340    | 39.191 | ng/ul# | 96       |
| 41) 2,4,6-Trichlorophenol     | 13.295 | 196  | 110304   | 44.299 | ng/ul  | 91       |
| 42) 2,4,5-Trichlorophenol     | 13.365 | 196  | 118773   | 43.920 | ng/ul  | 99       |
| 43) 1,1'-Biphenyl             | 13.700 | 154  | 441308   | 43.523 | ng/ul  | 96       |
| 44) 2-Chloronaphthalene       | 13.753 | 162  | 331049   | 42.474 | ng/ul  | 97       |
| 45) 2-Nitroaniline            | 13.970 | 65   | 107614   | 42.165 | ng/ul  | 93       |
| 47) Dimethylphthalate         | 14.305 | 163  | 418960   | 41.302 | ng/ul  | 99       |
| 48) 2,6-Dinitrotoluene        | 14.452 | 165  | 89189    | 43.582 | ng/ul  | 95       |
| 50) Acenaphthylene            | 14.593 | 152  | 549656   | 42.573 | ng/ul  | 99       |
| 51) 3-Nitroaniline            | 14.793 | 138  | 91413    | 45.618 | ng/ul  | 90       |
| 52) Acenaphthene              | 14.928 | 153  | 368844   | 42.223 | ng/ul  | 99       |
| 53) 2,4-Dinitrophenol         | 14.993 | 184  | 33676    | 22.909 | ng/ul  | 91       |
| 55) 4-Nitrophenol             | 15.075 | 109  | 50679    | 39.713 | ng/ul  | 86       |
| 56) Dibenzofuran              | 15.257 | 168  | 502691   | 41.705 | ng/ul  | 98       |
| 57) 2,4-Dinitrotoluene        | 15.228 | 165  | 124777   | 41.632 | ng/ul# | 100      |
| 58) 2,3,4,6-Tetrachlorophenol | 15.469 | 232  | 108529   | 41.824 | ng/ul  | 95       |
| 59) Diethylphthalate          | 15.651 | 149  | 409367   | 40.577 | ng/ul  | 98       |
| 61) Fluorene                  | 15.909 | 166  | 417024   | 40.664 | ng/ul  | 98       |
| 62) 4-Chlorophenyl-phenyle... | 15.886 | 204  | 214731   | 42.374 | ng/ul  | 96       |
| 63) 4-Nitroaniline            | 15.956 | 138  | 90748    | 50.302 | ng/ul  | 91       |
| 66) 4,6-Dinitro-2-methylph... | 15.980 | 198  | 78139    | 40.626 | ng/ul  | 95       |
| 67) N-Nitrosodiphenylamine    | 16.109 | 169  | 348703   | 44.988 | ng/ul  | 99       |
| 68) 4-Bromophenyl-phenylether | 16.785 | 248  | 129424   | 44.499 | ng/ul  | 98       |
| 69) Hexachlorobenzene         | 16.879 | 284  | 145434   | 43.807 | ng/ul  | 97       |
| 70) Atrazine                  | 17.043 | 200  | 126945   | 41.308 | ng/ul  | 98       |
| 71) Pentachlorophenol         | 17.237 | 266  | 71217    | 36.910 | ng/ul  | 93       |
| 72) Phenanthrene              | 17.654 | 178  | 718719   | 44.307 | ng/ul  | 99       |
| 74) Anthracene                | 17.748 | 178  | 728175   | 44.137 | ng/ul  | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 13.659 | 216  | 162053   | 45.541 | ng/uL  | 93       |
| 76) Pentachlorobenzene        | 15.151 | 250  | 158945   | 43.159 | ng/uL  | 96       |
| 77) Carbazole                 | 18.024 | 167  | 612477   | 44.610 | ng/ul  | 98       |
| 78) Di-n-butylphthalate       | 18.524 | 149  | 734676   | 45.229 | ng/ul  | 99       |
| 80) Fluoranthene              | 19.646 | 202  | 803180   | 44.289 | ng/ul  | 98       |
| 82) Pyrene                    | 20.016 | 202  | 830515   | 44.066 | ng/ul  | 97       |
| 83) Butylbenzylphthalate      | 20.862 | 149  | 312540   | 44.881 | ng/ul  | 99       |
| 84) 3,3'-Dichlorobenzidine    | 21.808 | 252  | 280435   | 46.726 | ng/ul  | 98       |
| 85) Benzo(a)anthracene        | 21.890 | 228  | 810408   | 44.107 | ng/ul  | 97       |
| 86) Bis(2-ethylhexyl)phtha... | 21.720 | 149  | 461052   | 44.268 | ng/ul  | 99       |
| 87) Chrysene                  | 21.961 | 228  | 753085   | 43.636 | ng/ul  | 96       |
| 89) Di-n-octyl phthalate      | 23.013 | 149  | 797333   | 45.287 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 24.264 | 252  | 823833   | 44.811 | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene      | 24.340 | 252  | 854464   | 45.774 | ng/ul# | 95       |
| 93) Benzo(a)pyrene            | 25.228 | 252  | 794365   | 45.360 | ng/ul  | 98       |
| 94) Indeno(1,2,3-cd)pyrene    | 29.417 | 276  | 906537   | 44.939 | ng/ul  | 95       |
| 95) Dibenzo(a,h)anthracene    | 29.493 | 278  | 738032   | 44.994 | ng/ul  | 98       |
| 96) Benzo(g,h,i)perylene      | 30.715 | 276  | 727405m  | 45.214 | ng/ul  |          |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG100923\  
Data File : BG059415.D  
Acq On : 10 Oct 2023 3:27  
Operator : MA/JU  
Sample : PB156170BS  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

## Instrument :

BNA\_G

## ClientSampleId :

SLCS170

## Manual Integrations

## APPROVED

Reviewed By :Yogesh Patel 10/11/2023

Supervised By :mohammad ahmed 10/11/2023

Quant Time: Oct 10 05:38:08 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG100723.MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Sat Oct 07 22:45:22 2023

Response via : Initial Calibration

