

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG111124\  
 Data File : BG063323.D  
 Acq On : 12 Nov 2024 13:22  
 Operator : RC/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD020529

Quant Time: Nov 12 13:02:46 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG110624.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 06 15:45:52 2024  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Yogesh Patel 11/13/2024  
 Supervised By :mohammad ahmed 11/14/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.832  | 152  | 106021   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.623 | 136  | 497797   | 20.000 | ng/u1  | # 0.00   |
| 38) Acenaphthene-d10          | 14.472 | 164  | 441034   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.221 | 188  | 1043856  | 20.000 | ng/u1  | # 0.00   |
| 79) Chrysene-d12              | 21.475 | 240  | 1065059  | 20.000 | ng/u1  | # 0.00   |
| 88) Perylene-d12              | 24.513 | 264  | 1126676  | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.296  | 96   | 20532    | 8.757  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.702  | 84   | 139085   | 17.947 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.010  | 99   | 194140   | 20.293 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.168  | 67   | 112087   | 19.975 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.368  | 132  | 135386   | 21.803 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.543  | 113  | 171357   | 21.132 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 8.990  | 128  | 72815    | 20.806 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.707  | 143  | 93043    | 20.570 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.253 | 165  | 195390   | 22.012 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.758 | 131  | 227018   | 19.635 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.878 | 166  | 591277   | 17.115 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.166 | 160  | 706031   | 20.880 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.683 | 143  | 82290    | 17.398 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.464 | 176  | 556946   | 19.102 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.594 | 200  | 115294   | 17.574 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.321 | 188  | 952558   | 21.232 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.618 | 212  | 1148380  | 21.506 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.301 | 264  | 1138444  | 20.928 | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.332  | 88   | 23807    | 8.248  | ng/uL# | 81       |
| 5) Pyridine                   | 3.725  | 79   | 144881   | 18.323 | ng/u1  | 99       |
| 6) Benzaldehyde               | 6.974  | 77   | 133282   | 25.207 | ng/u1  | 95       |
| 8) Phenol                     | 7.033  | 94   | 208978   | 20.045 | ng/u1  | 94       |
| 10) Bis(2-Chloroethyl)ether   | 7.256  | 93   | 138479   | 19.531 | ng/u1  | 87       |
| 12) 2-Chlorophenol            | 7.397  | 128  | 144647   | 21.777 | ng/u1  | 95       |
| 13) 2-Methylphenol            | 8.279  | 108  | 154958   | 20.348 | ng/u1  | 94       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.349  | 45   | 194848   | 21.742 | ng/u1# | 94       |
| 16) Acetophenone              | 8.649  | 105  | 252334   | 19.327 | ng/u1  | 93       |
| 17) N-Nitroso-di-n-propyla... | 8.637  | 70   | 144742   | 19.006 | ng/u1  | 99       |
| 18) 4-Methylphenol            | 8.602  | 108  | 178802   | 20.648 | ng/u1  | 98       |
| 19) Hexachloroethane          | 8.913  | 117  | 58762    | 21.831 | ng/u1  | 90       |
| 22) Nitrobenzene              | 9.031  | 77   | 218867   | 20.283 | ng/u1# | 93       |
| 23) Isophorone                | 9.548  | 82   | 404575   | 17.764 | ng/u1  | 96       |
| 25) 2-Nitrophenol             | 9.742  | 139  | 92698    | 20.732 | ng/u1# | 87       |
| 26) 2,4-Dimethylphenol        | 9.806  | 107  | 200304   | 20.538 | ng/u1  | 93       |
| 27) Bis(2-Chloroethoxy)met... | 10.036 | 93   | 225641   | 20.552 | ng/u1# | 97       |
| 29) 2,4-Dichlorophenol        | 10.276 | 162  | 184129   | 22.004 | ng/u1  | 93       |
| 30) Naphthalene               | 10.676 | 128  | 537717   | 21.111 | ng/u1  | 99       |
| 32) 4-Chloroaniline           | 10.782 | 127  | 218659   | 19.760 | ng/u1  | 99       |
| 33) Hexachlorobutadiene       | 10.964 | 225  | 160744   | 20.070 | ng/u1  | 99       |
| 34) Caprolactam               | 11.534 | 113  | 57831m   | 16.627 | ng/u1  |          |
| 35) 4-Chloro-3-methylphenol   | 11.916 | 107  | 202549   | 20.172 | ng/u1  | 94       |
| 36) 2-Methylnaphthalene       | 12.286 | 142  | 406198   | 20.121 | ng/u1  | 100      |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG111124\  
 Data File : BG063323.D  
 Acq On : 12 Nov 2024 13:22  
 Operator : RC/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD020529

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 11/13/2024  
 Supervised By :mohammad ahmed 11/14/2024

Quant Time: Nov 12 13:02:46 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG110624.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 06 15:45:52 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 37) 1-Methylnaphthalene       | 12.509 | 142  | 415183   | 19.777 | ng/ul  | 93       |
| 39) 1,2,4,5-Tetrachloroben... | 12.662 | 216  | 321746   | 20.695 | ng/ul  | 96       |
| 40) Hexachlorocyclopentadiene | 12.638 | 237  | 149867   | 17.653 | ng/ul  | 93       |
| 41) 2,4,6-Trichlorophenol     | 12.903 | 196  | 179058   | 21.336 | ng/ul  | 99       |
| 42) 2,4,5-Trichlorophenol     | 12.979 | 196  | 197117   | 21.660 | ng/ul  | 99       |
| 43) 1,1'-Biphenyl             | 13.302 | 154  | 589799   | 21.098 | ng/ul  | 97       |
| 44) 2-Chloronaphthalene       | 13.343 | 162  | 455134   | 20.912 | ng/ul  | 96       |
| 45) 2-Nitroaniline            | 13.549 | 65   | 146277   | 19.703 | ng/ul  | 96       |
| 47) Dimethylphthalate         | 13.925 | 163  | 579153   | 17.484 | ng/ul  | 99       |
| 48) 2,6-Dinitrotoluene        | 14.049 | 165  | 123913   | 19.549 | ng/ul  | 89       |
| 50) Acenaphthylene            | 14.195 | 152  | 690507   | 20.500 | ng/ul  | 98       |
| 51) 3-Nitroaniline            | 14.378 | 138  | 120137   | 21.714 | ng/ul  | 90       |
| 52) Acenaphthene              | 14.536 | 153  | 512980   | 21.279 | ng/ul  | 98       |
| 53) 2,4-Dinitrophenol         | 14.595 | 184  | 67908    | 16.250 | ng/ul  | 91       |
| 55) 4-Nitrophenol             | 14.695 | 109  | 95026    | 15.458 | ng/ul  | 100      |
| 56) Dibenzofuran              | 14.871 | 168  | 754878   | 20.455 | ng/ul  | 99       |
| 57) 2,4-Dinitrotoluene        | 14.842 | 165  | 191750   | 19.127 | ng/ul# | 99       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.106 | 232  | 179271   | 19.021 | ng/ul# | 97       |
| 59) Diethylphthalate          | 15.294 | 149  | 586971   | 17.284 | ng/ul  | 98       |
| 61) Fluorene                  | 15.523 | 166  | 621490   | 19.898 | ng/ul  | 97       |
| 62) 4-Chlorophenyl-phenyle... | 15.517 | 204  | 383800   | 19.120 | ng/ul  | 100      |
| 63) 4-Nitroaniline            | 15.541 | 138  | 114433   | 20.435 | ng/ul  | 91       |
| 66) 4,6-Dinitro-2-methylph... | 15.611 | 198  | 121141   | 17.793 | ng/ul# | 99       |
| 67) N-Nitrosodiphenylamine    | 15.729 | 169  | 525925   | 22.585 | ng/ul  | 98       |
| 68) 4-Bromophenyl-phenylether | 16.410 | 248  | 256116   | 21.462 | ng/ul  | 98       |
| 69) Hexachlorobenzene         | 16.534 | 284  | 271744   | 21.375 | ng/ul  | 97       |
| 70) Atrazine                  | 16.681 | 200  | 249767   | 19.606 | ng/ul  | 98       |
| 71) Pentachlorophenol         | 16.880 | 266  | 106996   | 16.788 | ng/ul  | 89       |
| 72) Phenanthrene              | 17.262 | 178  | 1019574  | 21.378 | ng/ul  | 98       |
| 74) Anthracene                | 17.356 | 178  | 1045101  | 21.391 | ng/ul  | 97       |
| 75) 1,2,3,4-Tetrachloroben... | 13.267 | 216  | 326974   | 24.223 | ng/ul  | 97       |
| 76) Pentachlorobenzene        | 14.795 | 250  | 333741   | 23.185 | ng/ul  | 97       |
| 77) Carbazole                 | 17.627 | 167  | 861856   | 21.247 | ng/ul  | 99       |
| 78) Di-n-butylphthalate       | 18.191 | 149  | 958063   | 20.764 | ng/ul  | 99       |
| 80) Fluoranthene              | 19.284 | 202  | 1285112  | 22.265 | ng/ul# | 98       |
| 82) Pyrene                    | 19.648 | 202  | 1336073  | 21.838 | ng/ul# | 96       |
| 83) Butylbenzylphthalate      | 20.547 | 149  | 416202   | 23.277 | ng/ul  | 97       |
| 84) 3,3'-Dichlorobenzidine    | 21.381 | 252  | 505898   | 22.090 | ng/ul  | 99       |
| 85) Benzo(a)anthracene        | 21.457 | 228  | 1439776  | 21.371 | ng/ul  | 98       |
| 86) Bis(2-ethylhexyl)phtha... | 21.375 | 149  | 617349   | 23.704 | ng/ul  | 97       |
| 87) Chrysene                  | 21.522 | 228  | 1339308  | 21.323 | ng/ul  | 97       |
| 89) Di-n-octyl phthalate      | 22.521 | 149  | 1051145  | 23.675 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.549 | 252  | 1444875  | 21.314 | ng/ul  | 98       |
| 91) Benzo(k)fluoranthene      | 23.614 | 252  | 1404153  | 20.796 | ng/ul  | 99       |
| 93) Benzo(a)pyrene            | 24.372 | 252  | 1299639  | 20.727 | ng/ul  | 97       |
| 94) Indeno(1,2,3-cd)pyrene    | 27.938 | 276  | 1592749  | 20.708 | ng/ul  | 97       |
| 95) Dibenzo(a,h)anthracene    | 27.979 | 278  | 1314603  | 20.964 | ng/ul  | 98       |
| 96) Benzo(g,h,i)perylene      | 28.996 | 276  | 1209941  | 20.748 | ng/ul# | 97       |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG111124\  
 Data File : BG063323.D  
 Acq On : 12 Nov 2024 13:22  
 Operator : RC/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :

BNA\_G

ClientSampleId :

SSTD020529

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 11/13/2024

Supervised By :mohammad ahmed 11/14/2024

Quant Time: Nov 12 13:02:46 2024

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG110624.MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 06 15:45:52 2024

Response via : Initial Calibration

