

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121523\  
 Data File : BG060046.D  
 Acq On : 15 Dec 2023 22:35  
 Operator : MA/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

## Instrument :

BNA\_G

## ClientSampleId :

SSTDCCC040

## Manual Integrations

## APPROVED

Reviewed By :Yogesh Patel 12/18/2023

Supervised By :mohammad ahmed 12/18/2023

Quant Time: Dec 16 02:06:44 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG121323.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Thu Dec 14 01:47:12 2023

Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.958  | 152  | 63905    | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8            | 10.766 | 136  | 280164   | 20.000 | ng    | # 0.00   |
| 39) Acenaphthene-d10          | 14.591 | 164  | 233898   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10          | 17.341 | 188  | 722381   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12              | 21.612 | 240  | 869262   | 20.000 | ng    | 0.00     |
| 86) Perylene-d12              | 24.803 | 264  | 996959   | 20.000 | ng    | 0.00     |
|                               |        |      |          |        |       |          |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 5) 2-Fluorophenol             | 5.525  | 112  | 326174   | 87.374 | ng    | 0.00     |
| 7) Phenol-d6                  | 7.112  | 99   | 466842   | 88.902 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 9.121  | 82   | 490290   | 79.465 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 16.084 | 330  | 460546   | 84.490 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 13.216 | 172  | 1526248  | 79.863 | ng    | 0.00     |
| 79) Terphenyl-d14             | 19.961 | 244  | 3793087  | 79.666 | ng    | 0.00     |
|                               |        |      |          |        |       |          |
| Target Compounds              |        |      |          |        |       | Qvalue   |
| 2) 1,4-Dioxane                | 3.440  | 88   | 75837    | 42.931 | ng    | 90       |
| 3) Pyridine                   | 3.827  | 79   | 206623   | 41.766 | ng    | 94       |
| 4) n-Nitrosodimethylamine     | 3.745  | 42   | 90949    | 37.376 | ng    | 84       |
| 6) Aniline                    | 7.282  | 93   | 242845   | 45.404 | ng    | 95       |
| 8) 2-Chlorophenol             | 7.523  | 128  | 154393   | 43.644 | ng    | 94       |
| 9) Benzaldehyde               | 7.094  | 77   | 137082   | 42.933 | ng    | 99       |
| 10) Phenol                    | 7.141  | 94   | 240940   | 45.674 | ng    | 96       |
| 11) bis(2-Chloroethyl)ether   | 7.376  | 93   | 152355   | 40.343 | ng    | 91       |
| 12) 1,3-Dichlorobenzene       | 7.846  | 146  | 206144   | 43.577 | ng    | 99       |
| 13) 1,4-Dichlorobenzene       | 7.993  | 146  | 207235   | 42.099 | ng    | 96       |
| 14) 1,2-Dichlorobenzene       | 8.310  | 146  | 209074   | 44.145 | ng    | 94       |
| 15) Benzyl Alcohol            | 8.193  | 79   | 197164   | 46.716 | ng    | 99       |
| 16) 2,2'-oxybis(1-Chloropr... | 8.487  | 45   | 338860m  | 50.144 | ng    |          |
| 17) 2-Methylphenol            | 8.393  | 107  | 164967   | 46.906 | ng    | 92       |
| 18) Hexachloroethane          | 9.045  | 117  | 68369    | 45.849 | ng    | 97       |
| 19) n-Nitroso-di-n-propyla... | 8.763  | 70   | 162572   | 47.986 | ng    | # 94     |
| 20) 3+4-Methylphenols         | 8.722  | 107  | 235697   | 47.979 | ng    | 93       |
| 22) Acetophenone              | 8.780  | 105  | 321730   | 39.407 | ng    | 97       |
| 24) Nitrobenzene              | 9.162  | 77   | 247714   | 39.546 | ng    | 99       |
| 25) Isophorone                | 9.679  | 82   | 460992   | 39.038 | ng    | 99       |
| 26) 2-Nitrophenol             | 9.873  | 139  | 97798    | 42.453 | ng    | 90       |
| 27) 2,4-Dimethylphenol        | 9.926  | 122  | 129940   | 41.150 | ng    | 93       |
| 28) bis(2-Chloroethoxy)met... | 10.167 | 93   | 227160   | 37.258 | ng    | 98       |
| 29) 2,4-Dichlorophenol        | 10.408 | 162  | 231982   | 40.904 | ng    | 93       |
| 30) 1,2,4-Trichlorobenzene    | 10.625 | 180  | 297304   | 38.728 | ng    | 100      |
| 31) Naphthalene               | 10.813 | 128  | 592339   | 40.614 | ng    | 96       |
| 32) Benzoic acid              | 10.061 | 122  | 144826   | 50.420 | ng    | 95       |
| 33) 4-Chloroaniline           | 10.919 | 127  | 234875   | 42.324 | ng    | 97       |
| 34) Hexachlorobutadiene       | 11.101 | 225  | 258511   | 36.792 | ng    | 93       |
| 35) Caprolactam               | 11.700 | 113  | 70081    | 48.643 | ng    | 96       |
| 36) 4-Chloro-3-methylphenol   | 12.035 | 107  | 225355   | 43.690 | ng    | 99       |
| 37) 2-Methylnaphthalene       | 12.417 | 142  | 462618   | 40.442 | ng    | 96       |
| 38) 1-Methylnaphthalene       | 12.640 | 142  | 460445   | 40.749 | ng    | 99       |
| 40) 1,2,4,5-Tetrachloroben... | 12.782 | 216  | 465411   | 38.168 | ng    | 99       |
| 41) Hexachlorocyclopentadiene | 12.770 | 237  | 277058   | 54.992 | ng    | 97       |
| 43) 2,4,6-Trichlorophenol     | 13.022 | 196  | 270819   | 39.611 | ng    | 95       |

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## Manual Integrations

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Quant Time: Dec 16 02:06:44 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG121323.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Thu Dec 14 01:47:12 2023

Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 13.093 | 196  | 308934   | 41.542 | ng    | 99       |
| 46) 1,1'-Biphenyl             | 13.428 | 154  | 718317   | 41.411 | ng    | 99       |
| 47) 2-Chloronaphthalene       | 13.469 | 162  | 598006   | 39.808 | ng    | 97       |
| 48) 2-Nitroaniline            | 13.669 | 65   | 162114   | 49.992 | ng    | 87       |
| 49) Acenaphthylene            | 14.315 | 152  | 872150   | 42.443 | ng    | 98       |
| 50) Dimethylphthalate         | 14.051 | 163  | 762069   | 41.974 | ng    | 98       |
| 51) 2,6-Dinitrotoluene        | 14.168 | 165  | 167360   | 44.500 | ng    | 98       |
| 52) Acenaphthene              | 14.662 | 154  | 638101m  | 45.111 | ng    |          |
| 53) 3-Nitroaniline            | 14.497 | 138  | 144599   | 48.115 | ng    | 86       |
| 54) 2,4-Dinitrophenol         | 14.697 | 184  | 183420   | 66.195 | ng    | # 89     |
| 55) Dibenzofuran              | 14.991 | 168  | 966643   | 41.986 | ng    | 99       |
| 56) 4-Nitrophenol             | 14.797 | 139  | 141436   | 55.997 | ng    | 94       |
| 57) 2,4-Dinitrotoluene        | 14.950 | 165  | 245099   | 46.361 | ng    | 94       |
| 58) Fluorene                  | 15.643 | 166  | 810416   | 41.744 | ng    | 98       |
| 59) 2,3,4,6-Tetrachlorophenol | 15.214 | 232  | 319505m  | 41.274 | ng    |          |
| 60) Diethylphthalate          | 15.414 | 149  | 714747   | 42.786 | ng    | 100      |
| 61) 4-Chlorophenyl-phenyle... | 15.637 | 204  | 568391   | 40.652 | ng    | 95       |
| 62) 4-Nitroaniline            | 15.660 | 138  | 153681   | 46.962 | ng    | 96       |
| 63) Azobenzene                | 15.931 | 77   | 680766   | 41.334 | ng    | 96       |
| 65) 4,6-Dinitro-2-methylph... | 15.713 | 198  | 205505   | 45.373 | ng    | 88       |
| 66) n-Nitrosodiphenylamine    | 15.849 | 169  | 735796   | 41.694 | ng    | 98       |
| 67) 4-Bromophenyl-phenylether | 16.530 | 248  | 412019   | 40.425 | ng    | # 96     |
| 68) Hexachlorobenzene         | 16.648 | 284  | 477851   | 40.105 | ng    | 98       |
| 69) Atrazine                  | 16.794 | 200  | 230098   | 36.277 | ng    | 98       |
| 70) Pentachlorophenol         | 16.988 | 266  | 390236   | 50.556 | ng    | 92       |
| 71) Phenanthrene              | 17.388 | 178  | 1472886  | 42.094 | ng    | 98       |
| 72) Anthracene                | 17.476 | 178  | 1480290  | 42.077 | ng    | 99       |
| 73) Carbazole                 | 17.746 | 167  | 1290385  | 43.163 | ng    | 97       |
| 74) Di-n-butylphthalate       | 18.310 | 149  | 1216881  | 44.620 | ng    | 99       |
| 75) Fluoranthene              | 19.397 | 202  | 2116093  | 41.444 | ng    | 97       |
| 77) Benzidine                 | 19.579 | 184  | 605706   | 54.248 | ng    | 99       |
| 78) Pyrene                    | 19.762 | 202  | 2171253  | 41.205 | ng    | 99       |
| 80) Butylbenzylphthalate      | 20.655 | 149  | 481909   | 50.922 | ng    | 95       |
| 81) Benzo(a)anthracene        | 21.595 | 228  | 2413056  | 41.326 | ng    | 99       |
| 82) 3,3'-Dichlorobenzidine    | 21.512 | 252  | 838973   | 43.665 | ng    | 99       |
| 83) Chrysene                  | 21.659 | 228  | 2258015  | 40.818 | ng    | 99       |
| 84) Bis(2-ethylhexyl)phtha... | 21.512 | 149  | 709455   | 49.460 | ng    | 98       |
| 85) Di-n-octyl phthalate      | 22.717 | 149  | 1247422  | 50.275 | ng    | 97       |
| 87) Indeno(1,2,3-cd)pyrene    | 28.440 | 276  | 2969538  | 41.736 | ng    | 99       |
| 88) Benzo(b)fluoranthene      | 23.792 | 252  | 2504054  | 40.669 | ng    | 99       |
| 89) Benzo(k)fluoranthene      | 23.857 | 252  | 2510843  | 42.041 | ng    | 99       |
| 90) Benzo(a)pyrene            | 24.656 | 252  | 2254378  | 41.326 | ng    | # 98     |
| 91) Dibenzo(a,h)anthracene    | 28.510 | 278  | 2512171  | 42.224 | ng    | 97       |
| 92) Benzo(g,h,i)perylene      | 29.591 | 276  | 2451741  | 41.428 | ng    | 98       |

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

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Quant Time: Dec 16 02:06:44 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG121323.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Dec 14 01:47:12 2023  
 Response via : Initial Calibration

