

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG040725\  
 Data File : BG064203.D  
 Acq On : 7 Apr 2025 13:40  
 Operator : RC/JU  
 Sample : SSTDICC080  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTDICC080

Quant Time: Apr 07 16:06:21 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG040725.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Apr 07 15:51:43 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-------------------------------|--------|------|----------|---------|-------|----------|
| -----                         |        |      |          |         |       |          |
| Internal Standards            |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.862  | 152  | 33840    | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8            | 10.647 | 136  | 159501   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10          | 14.483 | 164  | 109813   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10          | 17.227 | 188  | 240363   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12              | 21.458 | 240  | 243539   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12              | 24.472 | 264  | 262382   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds   |        |      |          |         |       |          |
| 5) 2-Fluorophenol             | 5.447  | 112  | 350169   | 166.424 | ng    | 0.00     |
| 7) Phenol-d6                  | 7.033  | 99   | 502790   | 168.760 | ng    | 0.01     |
| 23) Nitrobenzene-d5           | 9.019  | 82   | 536875   | 167.976 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 15.976 | 330  | 240046   | 162.805 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 13.115 | 172  | 1200425  | 158.964 | ng    | 0.00     |
| 79) Terphenyl-d14             | 19.848 | 244  | 1876050  | 153.677 | ng    | 0.00     |
| Target Compounds              |        |      |          |         |       |          |
|                               |        |      |          |         |       | Qvalue   |
| 2) 1,4-Dioxane                | 3.367  | 88   | 64205    | 76.990  | ng    | 96       |
| 3) Pyridine                   | 3.755  | 79   | 191392m  | 86.770  | ng    |          |
| 4) n-Nitrosodimethylamine     | 3.673  | 42   | 145523   | 84.769  | ng    | 99       |
| 6) Aniline                    | 7.192  | 93   | 188337   | 68.471  | ng    | 97       |
| 8) 2-Chlorophenol             | 7.427  | 128  | 199947   | 85.096  | ng    | 96       |
| 10) Phenol                    | 7.057  | 94   | 254929   | 86.327  | ng    | 98       |
| 11) bis(2-Chloroethyl)ether   | 7.286  | 93   | 186900   | 81.760  | ng    | 98       |
| 12) 1,3-Dichlorobenzene       | 7.750  | 146  | 206550   | 80.992  | ng    | 97       |
| 13) 1,4-Dichlorobenzene       | 7.897  | 146  | 208657   | 81.198  | ng    | 97       |
| 14) 1,2-Dichlorobenzene       | 8.214  | 146  | 207551   | 81.483  | ng    | 95       |
| 15) Benzyl Alcohol            | 8.097  | 79   | 219688   | 84.758  | ng    | 98       |
| 16) 2,2'-oxybis(1-Chloropr... | 8.391  | 45   | 412631   | 80.855  | ng    | 99       |
| 17) 2-Methylphenol            | 8.303  | 107  | 178284   | 83.901  | ng    | 97       |
| 18) Hexachloroethane          | 8.937  | 117  | 82602    | 82.316  | ng    | 90       |
| 19) n-Nitroso-di-n-propyla... | 8.673  | 70   | 181387   | 81.303  | ng    | 97       |
| 20) 3+4-Methylphenols         | 8.632  | 107  | 247869   | 82.699  | ng    | 95       |
| 22) Acetophenone              | 8.679  | 105  | 349116   | 80.040  | ng    | # 97     |
| 24) Nitrobenzene              | 9.060  | 77   | 260278   | 81.752  | ng    | 97       |
| 25) Isophorone                | 9.583  | 82   | 464378   | 79.259  | ng    | 99       |
| 26) 2-Nitrophenol             | 9.760  | 139  | 117470   | 88.118  | ng    | 95       |
| 27) 2,4-Dimethylphenol        | 9.824  | 122  | 150111   | 82.771  | ng    | 98       |
| 28) bis(2-Chloroethoxy)met... | 10.059 | 93   | 271994   | 80.399  | ng    | 96       |
| 29) 2,4-Dichlorophenol        | 10.300 | 162  | 201289   | 84.370  | ng    | 93       |
| 30) 1,2,4-Trichlorobenzene    | 10.512 | 180  | 215522   | 81.932  | ng    | 98       |
| 31) Naphthalene               | 10.700 | 128  | 693364   | 79.730  | ng    | 99       |
| 32) Benzoic acid              | 10.024 | 122  | 173451m  | 91.655  | ng    |          |
| 33) 4-Chloroaniline           | 10.811 | 127  | 248615   | 79.012  | ng    | 98       |
| 34) Hexachlorobutadiene       | 10.988 | 225  | 148722   | 80.974  | ng    | 96       |
| 35) Caprolactam               | 11.599 | 113  | 70815m   | 78.295  | ng    |          |
| 36) 4-Chloro-3-methylphenol   | 11.934 | 107  | 255215   | 79.858  | ng    | 98       |
| 37) 2-Methylnaphthalene       | 12.310 | 142  | 505049   | 79.834  | ng    | 99       |
| 38) 1-Methylnaphthalene       | 12.527 | 142  | 499387   | 79.697  | ng    | 100      |
| 40) 1,2,4,5-Tetrachloroben... | 12.674 | 216  | 267142   | 82.114  | ng    | 98       |
| 41) Hexachlorocyclopentadiene | 12.662 | 237  | 104502   | 89.351  | ng    | 98       |
| 43) 2,4,6-Trichlorophenol     | 12.915 | 196  | 175337   | 84.494  | ng    | 94       |
| 44) 2,4,5-Trichlorophenol     | 12.985 | 196  | 197508   | 84.845  | ng    | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG040725\  
 Data File : BG064203.D  
 Acq On : 7 Apr 2025 13:40  
 Operator : RC/JU  
 Sample : SSTDICC080  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTDICC080

Quant Time: Apr 07 16:06:21 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG040725.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Apr 07 15:51:43 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-------------------------------|--------|------|----------|---------|-------|----------|
| 46) 1,1'-Biphenyl             | 13.320 | 154  | 696748   | 81.643  | ng    | 99       |
| 47) 2-Chloronaphthalene       | 13.361 | 162  | 500937   | 81.456  | ng    | 99       |
| 48) 2-Nitroaniline            | 13.561 | 65   | 194294   | 86.256  | ng    | 94       |
| 49) Acenaphthylene            | 14.207 | 152  | 785207   | 81.304  | ng    | 99       |
| 50) Dimethylphthalate         | 13.949 | 163  | 660580   | 78.442  | ng    | 100      |
| 51) 2,6-Dinitrotoluene        | 14.060 | 165  | 144452   | 83.955  | ng    | 89       |
| 52) Acenaphthene              | 14.554 | 154  | 517867   | 80.055  | ng    | 100      |
| 53) 3-Nitroaniline            | 14.390 | 138  | 143073   | 84.974  | ng    | 97       |
| 54) 2,4-Dinitrophenol         | 14.595 | 184  | 91943    | 81.153  | ng    | 86       |
| 55) Dibenzofuran              | 14.883 | 168  | 828407   | 79.766  | ng    | 99       |
| 56) 4-Nitrophenol             | 14.701 | 139  | 121549   | 89.313  | ng    | 93       |
| 57) 2,4-Dinitrotoluene        | 14.848 | 165  | 203658   | 83.305  | ng    | # 92     |
| 58) Fluorene                  | 15.535 | 166  | 641397   | 77.874  | ng    | 98       |
| 59) 2,3,4,6-Tetrachlorophenol | 15.112 | 232  | 186667   | 82.878  | ng    | # 98     |
| 60) Diethylphthalate          | 15.312 | 149  | 715830   | 79.610  | ng    | 98       |
| 61) 4-Chlorophenyl-phenyle... | 15.529 | 204  | 325474   | 78.835  | ng    | 98       |
| 62) 4-Nitroaniline            | 15.559 | 138  | 148768   | 86.812  | ng    | 95       |
| 63) Azobenzene                | 15.823 | 77   | 701864   | 79.601  | ng    | 98       |
| 65) 4,6-Dinitro-2-methylph... | 15.612 | 198  | 131471   | 93.176  | ng    | 95       |
| 66) n-Nitrosodiphenylamine    | 15.747 | 169  | 560284   | 79.816  | ng    | 99       |
| 67) 4-Bromophenyl-phenylether | 16.422 | 248  | 217335   | 80.655  | ng    | 95       |
| 68) Hexachlorobenzene         | 16.534 | 284  | 231179   | 79.285  | ng    | 96       |
| 70) Pentachlorophenol         | 16.875 | 266  | 165277   | 90.564  | ng    | 99       |
| 71) Phenanthrene              | 17.268 | 178  | 1011827  | 78.009  | ng    | 100      |
| 72) Anthracene                | 17.357 | 178  | 1020965  | 79.132  | ng    | 99       |
| 73) Carbazole                 | 17.627 | 167  | 937240   | 79.474  | ng    | 98       |
| 74) Di-n-butylphthalate       | 18.197 | 149  | 1164790  | 78.761  | ng    | 99       |
| 75) Fluoranthene              | 19.278 | 202  | 1197551  | 77.109  | ng    | 99       |
| 77) Benzidine                 | 19.460 | 184  | 350296   | 109.854 | ng    | 100      |
| 78) Pyrene                    | 19.642 | 202  | 1241726  | 79.331  | ng    | 98       |
| 80) Butylbenzylphthalate      | 20.535 | 149  | 526361   | 82.702  | ng    | 95       |
| 81) Benzo(a)anthracene        | 21.440 | 228  | 1270039  | 80.624  | ng    | 100      |
| 82) 3,3'-Dichlorobenzidine    | 21.364 | 252  | 430769   | 81.111  | ng    | 99       |
| 83) Chrysene                  | 21.505 | 228  | 1226196  | 79.353  | ng    | 99       |
| 84) Bis(2-ethylhexyl)phtha... | 21.370 | 149  | 760796   | 80.029  | ng    | 99       |
| 85) Di-n-octyl phthalate      | 22.515 | 149  | 1348999  | 81.830  | ng    | 99       |
| 87) Indeno(1,2,3-cd)pyrene    | 27.880 | 276  | 1479341  | 81.452  | ng    | # 93     |
| 88) Benzo(b)fluoranthene      | 23.526 | 252  | 1294381  | 80.112  | ng    | 98       |
| 89) Benzo(k)fluoranthene      | 23.590 | 252  | 1286440  | 81.452  | ng    | 97       |
| 90) Benzo(a)pyrene            | 24.337 | 252  | 1157946  | 82.062  | ng    | 100      |
| 91) Dibenzo(a,h)anthracene    | 27.950 | 278  | 1225232  | 81.848  | ng    | 98       |
| 92) Benzo(g,h,i)perylene      | 28.949 | 276  | 1236276  | 81.599  | ng    | 97       |

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

Quant Time: Apr 07 16:06:21 2025  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG040725.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Mon Apr 07 15:51:43 2025  
Response via : Initial Calibration

