

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG031522\  
 Data File : BG052666.D  
 Acq On : 15 Mar 2022 14:56  
 Operator : CG/JU  
 Sample : SSTDICV040  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 ICVBG031522

Manual Integrations  
 APPROVED

Reviewed By : Christian Giraldo 03/16/2022  
 Supervised By : Jagrut Upadhyay 03/16/2022

Quant Time: Mar 15 17:29:42 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG031522.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Mar 15 15:56:10 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.154  | 152  | 40770    | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8            | 10.974 | 136  | 165873   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10          | 14.781 | 164  | 122795   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10          | 17.518 | 188  | 292468   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12              | 21.812 | 240  | 281426   | 20.000 | ng    | 0.00     |
| 86) Perylene-d12              | 25.137 | 264  | 291778   | 20.000 | ng    | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 5) 2-Fluorophenol             | 5.705  | 112  | 185746   | 79.342 | ng    | 0.00     |
| 7) Phenol-d6                  | 7.297  | 99   | 278743   | 78.977 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 9.317  | 82   | 272294   | 82.391 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 16.261 | 330  | 134113   | 81.279 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 13.406 | 172  | 643244   | 77.258 | ng    | 0.00     |
| 79) Terphenyl-d14             | 20.126 | 244  | 1228257  | 77.636 | ng    | 0.00     |
| Target Compounds              |        |      |          |        |       |          |
|                               |        |      |          |        |       | Qvalue   |
| 2) 1,4-Dioxane                | 3.596  | 88   | 45119    | 38.528 | ng    | 94       |
| 3) Pyridine                   | 4.001  | 79   | 124169   | 41.363 | ng    | 90       |
| 4) n-Nitrosodimethylamine     | 3.907  | 42   | 51728    | 38.843 | ng    | 95       |
| 6) Aniline                    | 7.473  | 93   | 165122   | 39.900 | ng    | 99       |
| 8) 2-Chlorophenol             | 7.720  | 128  | 96574    | 39.260 | ng    | 97       |
| 9) Benzaldehyde               | 7.285  | 77   | 76013    | 38.850 | ng    | 97       |
| 10) Phenol                    | 7.326  | 94   | 138318   | 39.701 | ng    | 94       |
| 11) bis(2-Chloroethyl)ether   | 7.567  | 93   | 105622   | 40.145 | ng    | 94       |
| 12) 1,3-Dichlorobenzene       | 8.049  | 146  | 116232   | 39.244 | ng    | 97       |
| 13) 1,4-Dichlorobenzene       | 8.190  | 146  | 114984   | 38.640 | ng    | 94       |
| 14) 1,2-Dichlorobenzene       | 8.513  | 146  | 108988   | 38.726 | ng    | 96       |
| 15) Benzyl Alcohol            | 8.383  | 79   | 119392   | 40.175 | ng    | 100      |
| 16) 2,2'-oxybis(1-Chloropr... | 8.683  | 45   | 166171   | 36.267 | ng    | 97       |
| 17) 2-Methylphenol            | 8.583  | 107  | 90870    | 39.125 | ng    | 97       |
| 18) Hexachloroethane          | 9.253  | 117  | 42750    | 39.155 | ng    | 99       |
| 19) n-Nitroso-di-n-propyla... | 8.965  | 70   | 98402    | 39.365 | ng    | 97       |
| 20) 3+4-Methylphenols         | 8.912  | 107  | 130689   | 40.207 | ng    | 96       |
| 22) Acetophenone              | 8.977  | 105  | 164753   | 38.381 | ng    | # 95     |
| 24) Nitrobenzene              | 9.359  | 77   | 136247   | 41.202 | ng    | 93       |
| 25) Isophorone                | 9.887  | 82   | 253792   | 39.660 | ng    | 98       |
| 26) 2-Nitrophenol             | 10.075 | 139  | 47422    | 41.296 | ng    | 99       |
| 27) 2,4-Dimethylphenol        | 10.122 | 122  | 90802    | 38.651 | ng    | 92       |
| 28) bis(2-Chloroethoxy)met... | 10.363 | 93   | 147590   | 39.005 | ng    | 98       |
| 29) 2,4-Dichlorophenol        | 10.610 | 162  | 108666   | 41.107 | ng    | 96       |
| 30) 1,2,4-Trichlorobenzene    | 10.833 | 180  | 124284   | 39.854 | ng    | 96       |
| 31) Naphthalene               | 11.021 | 128  | 327144   | 38.140 | ng    | 100      |
| 32) Benzoic acid              | 10.240 | 122  | 65780    | 38.923 | ng    | 96       |
| 33) 4-Chloroaniline           | 11.115 | 127  | 140282   | 38.587 | ng    | 95       |
| 34) Hexachlorobutadiene       | 11.315 | 225  | 87407    | 39.093 | ng    | 97       |
| 35) Caprolactam               | 11.896 | 113  | 31037    | 42.998 | ng    | # 85     |
| 36) 4-Chloro-3-methylphenol   | 12.225 | 107  | 124617   | 40.595 | ng    | 98       |
| 37) 2-Methylnaphthalene       | 12.619 | 142  | 237824   | 37.860 | ng    | 95       |
| 38) 1-Methylnaphthalene       | 12.836 | 142  | 234153m  | 38.194 | ng    |          |
| 40) 1,2,4,5-Tetrachloroben... | 12.983 | 216  | 149895   | 39.072 | ng    | 99       |
| 41) Hexachlorocyclopentadiene | 12.971 | 237  | 88528    | 38.930 | ng    | 90       |
| 43) 2,4,6-Trichlorophenol     | 13.212 | 196  | 99509    | 40.503 | ng    | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG031522\  
 Data File : BG052666.D  
 Acq On : 15 Mar 2022 14:56  
 Operator : CG/JU  
 Sample : SSTDICV040  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 ICVBG031522

Manual Integrations  
 APPROVED

Reviewed By : Christian Giraldo 03/16/2022  
 Supervised By : Jagrut Upadhyay 03/16/2022

Quant Time: Mar 15 17:29:42 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG031522.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Mar 15 15:56:10 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 13.283 | 196  | 104184   | 41.780 | ng    | 93       |
| 46) 1,1'-Biphenyl             | 13.617 | 154  | 328620   | 38.267 | ng    | 99       |
| 47) 2-Chloronaphthalene       | 13.659 | 162  | 258732   | 38.258 | ng    | 97       |
| 48) 2-Nitroaniline            | 13.847 | 65   | 80002    | 43.255 | ng    | 97       |
| 49) Acenaphthylene            | 14.505 | 152  | 413327   | 38.340 | ng    | 96       |
| 50) Dimethylphthalate         | 14.228 | 163  | 360805   | 39.535 | ng    | 98       |
| 51) 2,6-Dinitrotoluene        | 14.340 | 165  | 66192    | 42.958 | ng    | # 84     |
| 52) Acenaphthene              | 14.845 | 154  | 269786   | 39.063 | ng    | 96       |
| 53) 3-Nitroaniline            | 14.663 | 138  | 78140    | 43.338 | ng    | # 98     |
| 54) 2,4-Dinitrophenol         | 14.869 | 184  | 33841    | 40.292 | ng    | # 76     |
| 55) Dibenzofuran              | 15.174 | 168  | 436620   | 38.713 | ng    | 98       |
| 56) 4-Nitrophenol             | 14.957 | 139  | 62037    | 42.180 | ng    | # 82     |
| 57) 2,4-Dinitrotoluene        | 15.121 | 165  | 94435    | 44.171 | ng    | 91       |
| 58) Fluorene                  | 15.820 | 166  | 346878   | 37.730 | ng    | 98       |
| 59) 2,3,4,6-Tetrachlorophenol | 15.392 | 232  | 108855   | 40.840 | ng    | 99       |
| 60) Diethylphthalate          | 15.591 | 149  | 374724   | 38.898 | ng    | 99       |
| 61) 4-Chlorophenyl-phenyle... | 15.815 | 204  | 192094   | 37.886 | ng    | 97       |
| 62) 4-Nitroaniline            | 15.826 | 138  | 78288    | 41.092 | ng    | 88       |
| 63) Azobenzene                | 16.108 | 77   | 382287   | 37.128 | ng    | 99       |
| 65) 4,6-Dinitro-2-methylph... | 15.885 | 198  | 54253    | 41.609 | ng    | 95       |
| 66) n-Nitrosodiphenylamine    | 16.020 | 169  | 320361   | 38.907 | ng    | 99       |
| 67) 4-Bromophenyl-phenylether | 16.707 | 248  | 133480   | 37.855 | ng    | 95       |
| 68) Hexachlorobenzene         | 16.831 | 284  | 147396   | 38.890 | ng    | 93       |
| 69) Atrazine                  | 16.972 | 200  | 127375   | 41.953 | ng    | 99       |
| 70) Pentachlorophenol         | 17.166 | 266  | 96472    | 41.029 | ng    | 92       |
| 71) Phenanthrene              | 17.565 | 178  | 600008   | 38.348 | ng    | 98       |
| 72) Anthracene                | 17.653 | 178  | 591023   | 38.211 | ng    | 100      |
| 73) Carbazole                 | 17.918 | 167  | 582299   | 38.012 | ng    | 99       |
| 74) Di-n-butylphthalate       | 18.481 | 149  | 660464   | 39.466 | ng    | 99       |
| 75) Fluoranthene              | 19.568 | 202  | 760808   | 38.307 | ng    | 99       |
| 77) Benzidine                 | 19.739 | 184  | 262012   | 36.097 | ng    | 99       |
| 78) Pyrene                    | 19.927 | 202  | 762988   | 39.084 | ng    | 99       |
| 80) Butylbenzylphthalate      | 20.814 | 149  | 272081   | 39.291 | ng    | 98       |
| 81) Benzo(a)anthracene        | 21.795 | 228  | 721766   | 38.301 | ng    | 98       |
| 82) 3,3'-Dichlorobenzidine    | 21.701 | 252  | 253924   | 38.157 | ng    | 98       |
| 83) Chrysene                  | 21.859 | 228  | 667647   | 37.686 | ng    | 100      |
| 84) Bis(2-ethylhexyl)phtha... | 21.707 | 149  | 369359   | 39.221 | ng    | 99       |
| 85) Di-n-octyl phthalate      | 22.964 | 149  | 621445   | 39.639 | ng    | 98       |
| 87) Indeno(1,2,3-cd)pyrene    | 28.950 | 276  | 765902   | 38.350 | ng    | # 91     |
| 88) Benzo(b)fluoranthene      | 24.080 | 252  | 679582   | 37.511 | ng    | 99       |
| 89) Benzo(k)fluoranthene      | 24.156 | 252  | 685193   | 38.441 | ng    | 98       |
| 90) Benzo(a)pyrene            | 24.984 | 252  | 588181   | 38.355 | ng    | 98       |
| 91) Dibenzo(a,h)anthracene    | 29.026 | 278  | 626412   | 38.065 | ng    | 98       |
| 92) Benzo(g,h,i)perylene      | 30.148 | 276  | 626576   | 38.482 | ng    | 99       |

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

## Manual Integrations

Reviewed By :Christian Giraldo 03/16/2022  
Supervised By :Jagrut Upadhyay 03/16/2022

Quant Time: Mar 15 17:29:42 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG031522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Tue Mar 15 15:56:10 2022  
Response via : Initial Calibration

