

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM080822\  
 Data File : BM036155.D  
 Acq On : 08 Aug 2022 21:49  
 Operator : CG/JU  
 Sample : N3963-11MS  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 RVE-216MS

Manual Integrations  
 APPROVED

Reviewed By : Christian Giraldo 08/09/2022  
 Supervised By : Jagrut Upadhyay 08/09/2022

Quant Time: Aug 08 23:41:50 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM080122.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Aug 08 23:36:56 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-------------------------------|--------|------|----------|---------|-------|----------|
| -----                         |        |      |          |         |       |          |
| Internal Standards            |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.845  | 152  | 274336   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8            | 10.651 | 136  | 1054753  | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10          | 14.486 | 164  | 558431   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10          | 17.227 | 188  | 1011997  | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12              | 21.409 | 240  | 1036810  | 20.000  | ng    | -0.01    |
| 86) Perylene-d12              | 23.768 | 264  | 1190878  | 20.000  | ng    | 0.00     |
| System Monitoring Compounds   |        |      |          |         |       |          |
| 5) 2-Fluorophenol             | 5.434  | 112  | 2115270  | 125.007 | ng    | 0.00     |
| 7) Phenol-d6                  | 7.057  | 99   | 2548291  | 118.355 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 9.016  | 82   | 1541190  | 81.793  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 15.980 | 330  | 778756   | 118.873 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 13.116 | 172  | 3124190  | 82.603  | ng    | 0.00     |
| 79) Terphenyl-d14             | 19.856 | 244  | 3945069  | 75.052  | ng    | 0.00     |
| Target Compounds              |        |      |          |         |       |          |
| 2) 1,4-Dioxane                | 3.275  | 88   | 305461   | 44.799  | ng    | Qvalue   |
| 3) Pyridine                   | 3.687  | 79   | 823795   | 45.055  | ng    | # 91     |
| 4) n-Nitrosodimethylamine     | 3.599  | 42   | 365486   | 48.644  | ng    | # 89     |
| 6) Aniline                    | 7.181  | 93   | 910034   | 38.168  | ng    | # 62     |
| 8) 2-Chlorophenol             | 7.422  | 128  | 870035   | 48.224  | ng    | 94       |
| 9) Benzaldehyde               | 6.987  | 77   | 567983   | 49.829  | ng    | 95       |
| 10) Phenol                    | 7.081  | 94   | 1036487  | 47.810  | ng    | 93       |
| 11) bis(2-Chloroethyl)ether   | 7.275  | 93   | 842961   | 47.201  | ng    | 92       |
| 12) 1,3-Dichlorobenzene       | 7.734  | 146  | 986171   | 49.327  | ng    | 96       |
| 13) 1,4-Dichlorobenzene       | 7.881  | 146  | 1006434  | 49.364  | ng    | 96       |
| 14) 1,2-Dichlorobenzene       | 8.198  | 146  | 954025   | 48.496  | ng    | 99       |
| 15) Benzyl Alcohol            | 8.104  | 79   | 687707m  | 47.597  | ng    | 98       |
| 16) 2,2'-oxybis(1-Chloropr... | 8.386  | 45   | 1110926  | 46.781  | ng    | 96       |
| 17) 2-Methylphenol            | 8.322  | 107  | 676730   | 47.156  | ng    | 95       |
| 18) Hexachloroethane          | 8.922  | 117  | 371679   | 50.641  | ng    | 97       |
| 19) n-Nitroso-di-n-propyla... | 8.663  | 70   | 581109   | 44.433  | ng    | 97       |
| 20) 3+4-Methylphenols         | 8.657  | 107  | 890655   | 45.681  | ng    | 88       |
| 22) Acetophenone              | 8.681  | 105  | 1231581  | 49.326  | ng    | 93       |
| 24) Nitrobenzene              | 9.057  | 77   | 897712   | 50.679  | ng    | # 92     |
| 25) Isophorone                | 9.586  | 82   | 1578055  | 51.481  | ng    | 98       |
| 26) 2-Nitrophenol             | 9.769  | 139  | 444145   | 53.160  | ng    | 98       |
| 27) 2,4-Dimethylphenol        | 9.851  | 122  | 731092   | 56.487  | ng    | 93       |
| 28) bis(2-Chloroethoxy)met... | 10.069 | 93   | 1007566  | 46.137  | ng    | 97       |
| 29) 2,4-Dichlorophenol        | 10.316 | 162  | 713652   | 49.032  | ng    | 100      |
| 30) 1,2,4-Trichlorobenzene    | 10.510 | 180  | 780391   | 47.314  | ng    | 99       |
| 31) Naphthalene               | 10.698 | 128  | 2548012  | 48.589  | ng    | 98       |
| 32) Benzoic acid              | 10.022 | 122  | 520944   | 50.777  | ng    | 99       |
| 33) 4-Chloroaniline           | 10.828 | 127  | 488305   | 23.105  | ng    | 95       |
| 34) Hexachlorobutadiene       | 10.975 | 225  | 475959   | 49.108  | ng    | 96       |
| 35) Caprolactam               | 11.627 | 113  | 206439   | 46.048  | ng    | 98       |
| 36) 4-Chloro-3-methylphenol   | 11.974 | 107  | 703049   | 45.956  | ng    | 95       |
| 37) 2-Methylnaphthalene       | 12.316 | 142  | 1652898  | 47.366  | ng    | 99       |
| 38) 1-Methylnaphthalene       | 12.533 | 142  | 1561009  | 45.679  | ng    | 97       |
| 40) 1,2,4,5-Tetrachloroben... | 12.680 | 216  | 788470   | 52.080  | ng    | 98       |
| 41) Hexachlorocyclopentadiene | 12.651 | 237  | 834289   | 104.047 | ng    | 99       |
| 43) 2,4,6-Trichlorophenol     | 12.933 | 196  | 516640   | 50.150  | ng    | 98       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM080822\  
 Data File : BM036155.D  
 Acq On : 08 Aug 2022 21:49  
 Operator : CG/JU  
 Sample : N3963-11MS  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 RVE-216MS

Quant Time: Aug 08 23:41:50 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM080122.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Aug 08 23:36:56 2022  
 Response via : Initial Calibration

Manual Integrations  
 APPROVED

Reviewed By :Christian Giraldo 08/09/2022  
 Supervised By :Jagrut Upadhyay 08/09/2022

| Compound                      | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-------------------------------|--------|------|----------|---------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 13.016 | 196  | 549349   | 50.515  | ng    | 96       |
| 46) 1,1'-Biphenyl             | 13.327 | 154  | 2045359  | 50.258  | ng    | 99       |
| 47) 2-Chloronaphthalene       | 13.369 | 162  | 1553423  | 49.873  | ng    | 100      |
| 48) 2-Nitroaniline            | 13.586 | 65   | 437158   | 51.500  | ng    | 92       |
| 49) Acenaphthylene            | 14.210 | 152  | 2425868  | 50.384  | ng    | 100      |
| 50) Dimethylphthalate         | 13.957 | 163  | 1760474  | 46.358  | ng    | 99       |
| 51) 2,6-Dinitrotoluene        | 14.080 | 165  | 385144   | 50.548  | ng    | 89       |
| 52) Acenaphthene              | 14.551 | 154  | 1690950m | 53.688  | ng    |          |
| 53) 3-Nitroaniline            | 14.416 | 138  | 307044   | 34.539  | ng    | 93       |
| 54) 2,4-Dinitrophenol         | 14.616 | 184  | 396836   | 98.421  | ng    | 89       |
| 55) Dibenzofuran              | 14.886 | 168  | 2213138  | 46.923  | ng    | 98       |
| 56) 4-Nitrophenol             | 14.757 | 139  | 730184   | 102.641 | ng    | 87       |
| 57) 2,4-Dinitrotoluene        | 14.863 | 165  | 508734   | 50.940  | ng    | 99       |
| 58) Fluorene                  | 15.533 | 166  | 1753271  | 47.125  | ng    | 99       |
| 59) 2,3,4,6-Tetrachlorophenol | 15.121 | 232  | 428655   | 46.532  | ng    | 97       |
| 60) Diethylphthalate          | 15.315 | 149  | 1745659  | 44.731  | ng    | 99       |
| 61) 4-Chlorophenyl-phenyle... | 15.533 | 204  | 835712   | 45.048  | ng    | 96       |
| 62) 4-Nitroaniline            | 15.574 | 138  | 421541m  | 46.094  | ng    |          |
| 63) Azobenzene                | 15.821 | 77   | 1675870  | 45.172  | ng    | 93       |
| 65) 4,6-Dinitro-2-methylph... | 15.621 | 198  | 236259   | 45.659  | ng    | 93       |
| 66) n-Nitrosodiphenylamine    | 15.751 | 169  | 1450416  | 50.770  | ng    | 99       |
| 67) 4-Bromophenyl-phenylether | 16.421 | 248  | 491761   | 48.580  | ng    | 96       |
| 68) Hexachlorobenzene         | 16.527 | 284  | 578220   | 49.784  | ng    | 95       |
| 69) Atrazine                  | 16.704 | 200  | 472328   | 59.087  | ng    | 97       |
| 70) Pentachlorophenol         | 16.886 | 266  | 709736   | 104.339 | ng    | 98       |
| 71) Phenanthrene              | 17.274 | 178  | 2614032  | 50.701  | ng    | 99       |
| 72) Anthracene                | 17.362 | 178  | 2584493  | 51.778  | ng    | 99       |
| 73) Carbazole                 | 17.645 | 167  | 2428754  | 48.023  | ng    | 99       |
| 74) Di-n-butylphthalate       | 18.203 | 149  | 2892496  | 47.910  | ng    | 99       |
| 75) Fluoranthene              | 19.286 | 202  | 3136859  | 54.353  | ng    | 98       |
| 77) Benzidine                 | 19.486 | 184  | 1585460  | 101.965 | ng    | 99       |
| 78) Pyrene                    | 19.650 | 202  | 3277259  | 50.974  | ng    | 99       |
| 80) Butylbenzylphthalate      | 20.550 | 149  | 1337879  | 48.214  | ng    | 98       |
| 81) Benzo(a)anthracene        | 21.397 | 228  | 3294986  | 51.596  | ng    | 99       |
| 82) 3,3'-Dichlorobenzidine    | 21.333 | 252  | 1137653  | 73.366  | ng    | 100      |
| 83) Chrysene                  | 21.450 | 228  | 3136194  | 51.662  | ng    | 99       |
| 84) Bis(2-ethylhexyl)phtha... | 21.327 | 149  | 1977061  | 47.227  | ng    | # 99     |
| 85) Di-n-octyl phthalate      | 22.233 | 149  | 3494061  | 51.141  | ng    | # 95     |
| 87) Indeno(1,2,3-cd)pyrene    | 26.221 | 276  | 4301816  | 51.403  | ng    | 99       |
| 88) Benzo(b)fluoranthene      | 23.050 | 252  | 3850829  | 55.191  | ng    | 99       |
| 89) Benzo(k)fluoranthene      | 23.097 | 252  | 3505900  | 49.750  | ng    | 99       |
| 90) Benzo(a)pyrene            | 23.662 | 252  | 3383798  | 57.802  | ng    | 99       |
| 91) Dibenzo(a,h)anthracene    | 26.232 | 278  | 3628976  | 51.804  | ng    | 99       |
| 92) Benzo(g,h,i)perylene      | 26.974 | 276  | 3719140  | 54.841  | ng    | 98       |

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

