

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF081121\  
 Data File : BF125055.D  
 Acq On : 11 Aug 2021 13:42  
 Operator : JU/CG  
 Sample : M3318-01MSD  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 20210536-COM-9MSD

Manual Integrations  
 APPROVED

mohammad  
 8/12/2021 1:43:47 PM

Quant Time: Aug 11 17:21:12 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF080421.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Aug 10 11:32:36 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-------------------------------|--------|------|----------|---------|-------|----------|
| -----                         |        |      |          |         |       |          |
| Internal Standards            |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 6.810  | 152  | 7177     | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8            | 8.098  | 136  | 26375    | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10          | 9.851  | 164  | 9321     | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10          | 11.339 | 188  | 17202    | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12              | 13.974 | 240  | 11091    | 20.000  | ng    | 0.00     |
| 86) Perylene-d12              | 15.427 | 264  | 10164    | 20.000  | ng    | 0.01     |
| System Monitoring Compounds   |        |      |          |         |       |          |
| 5) 2-Fluorophenol             | 5.440  | 112  | 37034    | 84.528  | ng    | 0.00     |
| 7) Phenol-d6                  | 6.451  | 99   | 43757    | 79.206  | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 7.381  | 82   | 26661    | 52.880  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 10.645 | 330  | 8252     | 99.095  | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 9.169  | 172  | 58883    | 92.436  | ng    | 0.00     |
| 79) Terphenyl-d14             | 12.916 | 244  | 36874    | 56.081  | ng    | 0.00     |
| Target Compounds              |        |      |          |         |       |          |
|                               |        |      |          |         |       | Qvalue   |
| 2) 1,4-Dioxane                | 2.587  | 88   | 6024     | 36.752  | ng    | # 100    |
| 3) Pyridine                   | 3.340  | 79   | 12976    | 34.069  | ng    | 85       |
| 4) n-Nitrosodimethylamine     | 3.293  | 42   | 11329    | 41.886  | ng    | # 81     |
| 6) Aniline                    | 6.475  | 93   | 15781    | 27.043  | ng    | # 81     |
| 8) 2-Chlorophenol             | 6.598  | 128  | 19922    | 41.453  | ng    | 88       |
| 9) Benzaldehyde               | 6.363  | 77   | 11390    | 37.896  | ng    | 91       |
| 10) Phenol                    | 6.469  | 94   | 24376    | 42.763  | ng    | 89       |
| 11) bis(2-Chloroethyl)ether   | 6.551  | 93   | 18840    | 44.058  | ng    | 98       |
| 12) 1,3-Dichlorobenzene       | 6.751  | 146  | 22189    | 42.420  | ng    | 99       |
| 13) 1,4-Dichlorobenzene       | 6.828  | 146  | 22437    | 42.908  | ng    | 95       |
| 14) 1,2-Dichlorobenzene       | 6.981  | 146  | 20795    | 42.434  | ng    | 90       |
| 15) Benzyl Alcohol            | 6.963  | 79   | 16065    | 40.314  | ng    | 93       |
| 16) 2,2'-oxybis(1-Chloropr... | 7.087  | 45   | 30371    | 44.331  | ng    | 88       |
| 17) 2-Methylphenol            | 7.075  | 107  | 15615    | 43.406  | ng    | # 93     |
| 18) Hexachloroethane          | 7.322  | 117  | 8152     | 39.819  | ng    | 87       |
| 19) n-Nitroso-di-n-propyla... | 7.228  | 70   | 13292    | 40.549  | ng    | # 84     |
| 20) 3+4-Methylphenols         | 7.228  | 107  | 18806    | 39.380  | ng    | # 70     |
| 22) Acetophenone              | 7.228  | 105  | 26166    | 40.508  | ng    | # 91     |
| 24) Nitrobenzene              | 7.398  | 77   | 19315    | 39.591  | ng    | 91       |
| 25) Isophorone                | 7.639  | 82   | 34589    | 40.196  | ng    | # 91     |
| 26) 2-Nitrophenol             | 7.716  | 139  | 11523    | 46.513  | ng    | # 90     |
| 27) 2,4-Dimethylphenol        | 7.751  | 122  | 18676    | 53.180  | ng    | 90       |
| 28) bis(2-Chloroethoxy)met... | 7.845  | 93   | 23208    | 47.387  | ng    | 96       |
| 29) 2,4-Dichlorophenol        | 7.957  | 162  | 15710    | 40.152  | ng    | 94       |
| 30) 1,2,4-Trichlorobenzene    | 8.039  | 180  | 17854    | 40.718  | ng    | 98       |
| 31) Naphthalene               | 8.122  | 128  | 51682    | 38.139  | ng    | # 96     |
| 32) Benzoic acid              | 7.892  | 122  | 9911     | 47.676  | ng    | 95       |
| 33) 4-Chloroaniline           | 8.169  | 127  | 4523     | 8.957   | ng    | 96       |
| 34) Hexachlorobutadiene       | 8.233  | 225  | 10642    | 40.642  | ng    | 97       |
| 35) Caprolactam               | 8.551  | 113  | 4448m    | 42.810  | ng    |          |
| 36) 4-Chloro-3-methylphenol   | 8.657  | 107  | 17072    | 41.712  | ng    | # 88     |
| 37) 2-Methylnaphthalene       | 8.810  | 142  | 37659    | 41.333  | ng    | 99       |
| 38) 1-Methylnaphthalene       | 8.910  | 142  | 38133    | 44.730  | ng    | 98       |
| 40) 1,2,4,5-Tetrachloroben... | 8.975  | 216  | 18787    | 70.220  | ng    | 98       |
| 41) Hexachlorocyclopentadiene | 8.957  | 237  | 14000    | 133.637 | ng    | 98       |
| 43) 2,4,6-Trichlorophenol     | 9.092  | 196  | 13303    | 74.693  | ng    | 97       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF081121\  
 Data File : BF125055.D  
 Acq On : 11 Aug 2021 13:42  
 Operator : JU/CG  
 Sample : M3318-01MSD  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 20210536-COM-9MSD

Manual Integrations  
 APPROVED

mohammad  
 8/12/2021 1:43:47 PM

Quant Time: Aug 11 17:21:12 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF080421.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Aug 10 11:32:36 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 9.133  | 196  | 13448    | 74.591 | ng    | 96       |
| 46) 1,1'-Biphenyl             | 9.275  | 154  | 49696    | 71.547 | ng    | 98       |
| 47) 2-Chloronaphthalene       | 9.298  | 162  | 38047    | 68.849 | ng    | 98       |
| 48) 2-Nitroaniline            | 9.398  | 65   | 11452    | 70.548 | ng    | 95       |
| 49) Acenaphthylene            | 9.716  | 152  | 34710    | 41.350 | ng    | 98       |
| 50) Dimethylphthalate         | 9.575  | 163  | 32724    | 51.808 | ng    | 99       |
| 51) 2,6-Dinitrotoluene        | 9.639  | 165  | 6063     | 43.106 | ng    | # 58     |
| 52) Acenaphthene              | 9.886  | 154  | 20515    | 40.314 | ng    | 96       |
| 53) 3-Nitroaniline            | 9.810  | 138  | 2610     | 17.618 | ng    | # 61     |
| 54) 2,4-Dinitrophenol         | 9.922  | 184  | 5349     | 75.338 | ng    | # 84     |
| 55) Dibenzofuran              | 10.057 | 168  | 32410    | 40.752 | ng    | 92       |
| 56) 4-Nitrophenol             | 9.986  | 139  | 7551     | 77.324 | ng    | # 51     |
| 57) 2,4-Dinitrotoluene        | 10.051 | 165  | 8611     | 45.093 | ng    | 96       |
| 58) Fluorene                  | 10.398 | 166  | 26074    | 42.758 | ng    | 94       |
| 59) 2,3,4,6-Tetrachlorophenol | 10.180 | 232  | 5970     | 44.134 | ng    | # 96     |
| 60) Diethylphthalate          | 10.275 | 149  | 30843    | 47.582 | ng    | 97       |
| 61) 4-Chlorophenyl-phenyle... | 10.392 | 204  | 13005    | 45.160 | ng    | 92       |
| 62) 4-Nitroaniline            | 10.427 | 138  | 5895     | 40.986 | ng    | 83       |
| 63) Azobenzene                | 10.551 | 77   | 32574    | 51.388 | ng    | 88       |
| 65) 4,6-Dinitro-2-methylph... | 10.457 | 198  | 4316     | 40.050 | ng    | 94       |
| 66) n-Nitrosodiphenylamine    | 10.516 | 169  | 22813    | 40.917 | ng    | 97       |
| 67) 4-Bromophenyl-phenylether | 10.880 | 248  | 8441     | 44.801 | ng    | # 88     |
| 68) Hexachlorobenzene         | 10.945 | 284  | 9155     | 44.663 | ng    | 95       |
| 69) Atrazine                  | 11.039 | 200  | 8161     | 48.712 | ng    | 98       |
| 70) Pentachlorophenol         | 11.145 | 266  | 8282     | 85.233 | ng    | 89       |
| 71) Phenanthrene              | 11.363 | 178  | 38467    | 41.223 | ng    | 98       |
| 72) Anthracene                | 11.416 | 178  | 39590    | 42.292 | ng    | 98       |
| 73) Carbazole                 | 11.569 | 167  | 33440    | 38.953 | ng    | 98       |
| 74) Di-n-butylphthalate       | 11.892 | 149  | 51144    | 46.877 | ng    | 99       |
| 75) Fluoranthene              | 12.551 | 202  | 39051    | 40.453 | ng    | 97       |
| 77) Benzidine                 | 12.669 | 184  | 10694    | 36.036 | ng    | 98       |
| 78) Pyrene                    | 12.780 | 202  | 38701    | 44.301 | ng    | 97       |
| 80) Butylbenzylphthalate      | 13.392 | 149  | 17689    | 45.500 | ng    | # 84     |
| 81) Benzo(a)anthracene        | 13.963 | 228  | 27520    | 38.076 | ng    | 99       |
| 82) 3,3'-Dichlorobenzidine    | 13.921 | 252  | 4857     | 25.490 | ng    | 97       |
| 83) Chrysene                  | 14.004 | 228  | 27736    | 39.696 | ng    | 98       |
| 84) Bis(2-ethylhexyl)phtha... | 13.945 | 149  | 26157    | 48.501 | ng    | 99       |
| 85) Di-n-octyl phthalate      | 14.557 | 149  | 37557    | 44.141 | ng    | 100      |
| 87) Indeno(1,2,3-cd)pyrene    | 16.886 | 276  | 29321    | 48.443 | ng    | # 89     |
| 88) Benzo(b)fluoranthene      | 15.004 | 252  | 27319    | 38.279 | ng    | # 94     |
| 89) Benzo(k)fluoranthene      | 15.033 | 252  | 23574    | 36.312 | ng    | 99       |
| 90) Benzo(a)pyrene            | 15.368 | 252  | 24260    | 39.920 | ng    | 93       |
| 91) Dibenzo(a,h)anthracene    | 16.892 | 278  | 25357    | 47.648 | ng    | # 92     |
| 92) Benzo(g,h,i)perylene      | 17.321 | 276  | 24431    | 47.859 | ng    | # 87     |

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

