

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG032221\  
 Data File : BG048444.D  
 Acq On : 22 Mar 2021 14:27  
 Operator : CG/JU  
 Sample : M1732-07MS  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SB-3B-(2.5-3.0)MS

Manual Integrations  
 APPROVED

mohammad  
 3/23/2021 11:21:51 AM

Quant Time: Mar 22 15:23:44 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG031821.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Mar 22 11:22:24 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.106  | 152  | 45413    | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8            | 10.926 | 136  | 179581   | 20.00  | ng    | # 0.00   |
| 39) Acenaphthene-d10          | 14.745 | 164  | 149957   | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10          | 17.495 | 188  | 386691   | 20.00  | ng    | # 0.00   |
| 76) Chrysene-d12              | 21.790 | 240  | 384239   | 20.00  | ng    | # 0.00   |
| 86) Perylene-d12              | 25.121 | 264  | 394558   | 20.00  | ng    | 0.00     |
|                               |        |      |          |        |       |          |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 5) 2-Fluorophenol             | 5.650  | 112  | 319087   | 116.19 | ng    | 0.00     |
| 7) Phenol-d6                  | 7.248  | 99   | 443126   | 110.89 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 9.269  | 82   | 325163   | 72.40  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 16.238 | 330  | 276161   | 113.24 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 13.371 | 172  | 728585   | 69.89  | ng    | 0.00     |
| 79) Terphenyl-d14             | 20.104 | 244  | 1481669  | 69.54  | ng    | 0.00     |
|                               |        |      |          |        |       |          |
| Target Compounds              |        |      |          |        |       | Qvalue   |
| 2) 1,4-Dioxane                | 3.517  | 88   | 39445    | 34.04  | ng    | # 89     |
| 3) Pyridine                   | 3.923  | 79   | 122377   | 37.89  | ng    | 88       |
| 4) n-Nitrosodimethylamine     | 3.835  | 42   | 78371    | 41.75  | ng    | # 68     |
| 6) Aniline                    | 7.419  | 93   | 63113    | 13.35  | ng    | 95       |
| 8) 2-Chlorophenol             | 7.660  | 128  | 136616   | 47.71  | ng    | 91       |
| 9) Benzaldehyde               | 7.231  | 77   | 129322   | 65.13  | ng    | # 81     |
| 10) Phenol                    | 7.278  | 94   | 202990   | 49.57  | ng    | # 68     |
| 11) bis(2-Chloroethyl)ether   | 7.513  | 93   | 125482   | 43.73  | ng    | 93       |
| 12) 1,3-Dichlorobenzene       | 7.994  | 146  | 147468   | 43.29  | ng    | # 89     |
| 13) 1,4-Dichlorobenzene       | 8.135  | 146  | 148189   | 43.82  | ng    | 93       |
| 14) 1,2-Dichlorobenzene       | 8.459  | 146  | 145512   | 44.85  | ng    | 94       |
| 15) Benzyl Alcohol            | 8.335  | 79   | 166030   | 46.80  | ng    | # 85     |
| 16) 2,2'-oxybis(1-Chloropr... | 8.635  | 45   | 141127   | 43.02  | ng    | 89       |
| 17) 2-Methylphenol            | 8.541  | 107  | 129003   | 49.68  | ng    | # 84     |
| 18) Hexachloroethane          | 9.199  | 117  | 55333    | 42.54  | ng    | 93       |
| 19) n-Nitroso-di-n-propyla... | 8.905  | 70   | 127472   | 42.98  | ng    | # 95     |
| 20) 3+4-Methylphenols         | 8.864  | 107  | 179998   | 50.61  | ng    | 88       |
| 22) Acetophenone              | 8.929  | 105  | 221911   | 43.94  | ng    | # 91     |
| 24) Nitrobenzene              | 9.311  | 77   | 197802   | 41.11  | ng    | # 77     |
| 25) Isophorone                | 9.839  | 82   | 357127   | 41.54  | ng    | # 92     |
| 26) 2-Nitrophenol             | 10.027 | 139  | 79720    | 43.59  | ng    | # 64     |
| 27) 2,4-Dimethylphenol        | 10.080 | 122  | 148618   | 52.08  | ng    | 86       |
| 28) bis(2-Chloroethoxy)met... | 10.321 | 93   | 187199   | 48.04  | ng    | # 92     |
| 29) 2,4-Dichlorophenol        | 10.562 | 162  | 159319   | 48.11  | ng    | 96       |
| 30) 1,2,4-Trichlorobenzene    | 10.785 | 180  | 165967   | 41.46  | ng    | 97       |
| 31) Naphthalene               | 10.979 | 128  | 415962   | 41.78  | ng    | 99       |
| 32) Benzoic acid              | 10.215 | 122  | 127079   | 57.22  | ng    | # 77     |
| 33) 4-Chloroaniline           | 11.073 | 127  | 24623    | 5.78   | ng    | # 86     |
| 34) Hexachlorobutadiene       | 11.261 | 225  | 126693   | 39.93  | ng    | 95       |
| 35) Caprolactam               | 11.855 | 113  | 63254m   | 53.50  | ng    |          |
| 36) 4-Chloro-3-methylphenol   | 12.190 | 107  | 188536   | 49.05  | ng    | # 81     |
| 37) 2-Methylnaphthalene       | 12.577 | 142  | 330355   | 43.47  | ng    | # 93     |
| 38) 1-Methylnaphthalene       | 12.795 | 142  | 310060   | 43.38  | ng    | # 99     |
| 40) 1,2,4,5-Tetrachloroben... | 12.942 | 216  | 231395   | 42.26  | ng    | # 97     |
| 41) Hexachlorocyclopentadiene | 12.924 | 237  | 313274   | 88.49  | ng    | 99       |
| 43) 2,4,6-Trichlorophenol     | 13.177 | 196  | 168147   | 45.30  | ng    | 97       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG032221\  
 Data File : BG048444.D  
 Acq On : 22 Mar 2021 14:27  
 Operator : CG/JU  
 Sample : M1732-07MS  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SB-3B-(2.5-3.0)MS

Manual Integrations  
 APPROVED

mohammad  
 3/23/2021 11:21:51 AM

Quant Time: Mar 22 15:23:44 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG031821.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Mar 22 11:22:24 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 13.247 | 196  | 179398   | 43.92  | ng    | # 92     |
| 46) 1,1'-Biphenyl             | 13.582 | 154  | 463939   | 43.11  | ng    | 97       |
| 47) 2-Chloronaphthalene       | 13.623 | 162  | 360998   | 41.60  | ng    | 98       |
| 48) 2-Nitroaniline            | 13.817 | 65   | 145565   | 45.32  | ng    | # 70     |
| 49) Acenaphthylene            | 14.469 | 152  | 625505   | 44.52  | ng    | 99       |
| 50) Dimethylphthalate         | 14.193 | 163  | 567113   | 46.47  | ng    | 99       |
| 51) 2,6-Dinitrotoluene        | 14.316 | 165  | 116303   | 44.28  | ng    | # 64     |
| 52) Acenaphthene              | 14.810 | 154  | 480942m  | 50.93  | ng    |          |
| 53) 3-Nitroaniline            | 14.640 | 138  | 18548    | 7.37   | ng    | # 87     |
| 54) 2,4-Dinitrophenol         | 14.851 | 184  | 174113   | 108.61 | ng    | # 72     |
| 55) Dibenzofuran              | 15.145 | 168  | 582561   | 41.42  | ng    | 93       |
| 56) 4-Nitrophenol             | 14.945 | 139  | 211058   | 93.78  | ng    | # 45     |
| 57) 2,4-Dinitrotoluene        | 15.098 | 165  | 165694   | 44.43  | ng    | # 72     |
| 58) Fluorene                  | 15.797 | 166  | 510708   | 43.26  | ng    | 96       |
| 59) 2,3,4,6-Tetrachlorophenol | 15.368 | 232  | 183273   | 45.47  | ng    | # 96     |
| 60) Diethylphthalate          | 15.556 | 149  | 534616   | 44.81  | ng    | 97       |
| 61) 4-Chlorophenyl-phenyle... | 15.785 | 204  | 296919   | 43.72  | ng    | 98       |
| 62) 4-Nitroaniline            | 15.809 | 138  | 117278   | 42.57  | ng    | # 53     |
| 63) Azobenzene                | 16.079 | 77   | 538858   | 41.51  | ng    | 88       |
| 65) 4,6-Dinitro-2-methylph... | 15.862 | 198  | 110469   | 45.36  | ng    | 85       |
| 66) n-Nitrosodiphenylamine    | 15.997 | 169  | 478450   | 42.71  | ng    | 96       |
| 67) 4-Bromophenyl-phenylether | 16.678 | 248  | 208391   | 42.79  | ng    | 95       |
| 68) Hexachlorobenzene         | 16.796 | 284  | 224185   | 42.96  | ng    | 95       |
| 69) Atrazine                  | 16.943 | 200  | 218266   | 48.44  | ng    | 95       |
| 70) Pentachlorophenol         | 17.143 | 266  | 312158   | 89.98  | ng    | 99       |
| 71) Phenanthrene              | 17.542 | 178  | 879739   | 43.07  | ng    | 98       |
| 72) Anthracene                | 17.630 | 178  | 897014   | 44.26  | ng    | 98       |
| 73) Carbazole                 | 17.895 | 167  | 804314   | 41.62  | ng    | 97       |
| 74) Di-n-butylphthalate       | 18.453 | 149  | 928563   | 43.41  | ng    | # 97     |
| 75) Fluoranthene              | 19.552 | 202  | 1127743  | 41.46  | ng    | 96       |
| 77) Benzidine                 | 19.722 | 184  | 471830   | 40.99  | ng    | 99       |
| 78) Pyrene                    | 19.910 | 202  | 1132422  | 42.91  | ng    | 98       |
| 80) Butylbenzylphthalate      | 20.791 | 149  | 408473   | 43.55  | ng    | # 86     |
| 81) Benzo(a)anthracene        | 21.772 | 228  | 1157037  | 43.42  | ng    | 99       |
| 82) 3,3'-Dichlorobenzidine    | 21.678 | 252  | 93297    | 9.70   | ng    | # 95     |
| 83) Chrysene                  | 21.837 | 228  | 1085749  | 42.50  | ng    | 98       |
| 84) Bis(2-ethylhexyl)phtha... | 21.673 | 149  | 572836   | 44.20  | ng    | 96       |
| 85) Di-n-octyl phthalate      | 22.924 | 149  | 968852   | 44.65  | ng    | # 96     |
| 87) Indeno(1,2,3-cd)pyrene    | 28.946 | 276  | 1327667  | 44.12  | ng    | # 84     |
| 88) Benzo(b)fluoranthene      | 24.064 | 252  | 1156354  | 41.59  | ng    | # 97     |
| 89) Benzo(k)fluoranthene      | 24.134 | 252  | 1084278  | 41.33  | ng    | # 95     |
| 90) Benzo(a)pyrene            | 24.969 | 252  | 1016556  | 40.10  | ng    | # 96     |
| 91) Dibenzo(a,h)anthracene    | 29.029 | 278  | 1098930  | 42.81  | ng    | # 92     |
| 92) Benzo(g,h,i)perylene      | 30.151 | 276  | 1086824  | 43.64  | ng    | # 92     |

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

