

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102921\  
 Data File : BF126007.D  
 Acq On : 29 Oct 2021 16:08  
 Operator : JU/CG  
 Sample : SSTDICC060  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SSTDICC060

Manual Integrations  
 APPROVED

mohammad  
 11/1/2021 1:37:51 PM

Quant Time: Oct 29 17:20:42 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF102921.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Oct 29 17:15:51 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-------------------------------|--------|------|----------|---------|-------|----------|
| -----                         |        |      |          |         |       |          |
| Internal Standards            |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 6.857  | 152  | 231371   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8            | 8.139  | 136  | 853098   | 20.000  | ng    | # 0.00   |
| 39) Acenaphthene-d10          | 9.892  | 164  | 478537   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10          | 11.374 | 188  | 845990   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12              | 14.015 | 240  | 497489   | 20.000  | ng    | # 0.00   |
| 86) Perylene-d12              | 15.474 | 264  | 415275   | 20.000  | ng    | 0.00     |
|                               |        |      |          |         |       |          |
| System Monitoring Compounds   |        |      |          |         |       |          |
| 5) 2-Fluorophenol             | 5.469  | 112  | 1594920  | 102.678 | ng    | 0.00     |
| 7) Phenol-d6                  | 6.492  | 99   | 2187004  | 102.652 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 7.428  | 82   | 1947741  | 117.932 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 10.686 | 330  | 498494   | 119.497 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 9.216  | 172  | 2717293  | 100.417 | ng    | 0.00     |
| 79) Terphenyl-d14             | 12.963 | 244  | 2942618  | 96.752  | ng    | 0.00     |
|                               |        |      |          |         |       |          |
| Target Compounds              |        |      |          |         |       | Qvalue   |
| 2) 1,4-Dioxane                | 2.622  | 88   | 399158   | 52.590  | ng    | # 100    |
| 3) Pyridine                   | 3.369  | 79   | 1117030  | 55.582  | ng    | 90       |
| 4) n-Nitrosodimethylamine     | 3.334  | 42   | 571447   | 44.333  | ng    | # 54     |
| 6) Aniline                    | 6.528  | 93   | 1383326  | 55.124  | ng    | # 91     |
| 8) 2-Chlorophenol             | 6.645  | 128  | 810929   | 51.722  | ng    | # 77     |
| 9) Benzaldehyde               | 6.404  | 77   | 594346   | 46.621  | ng    | 93       |
| 10) Phenol                    | 6.504  | 94   | 1130541  | 55.035  | ng    | 82       |
| 11) bis(2-Chloroethyl)ether   | 6.598  | 93   | 895177   | 54.112  | ng    | # 80     |
| 12) 1,3-Dichlorobenzene       | 6.798  | 146  | 880038m  | 50.812  | ng    |          |
| 13) 1,4-Dichlorobenzene       | 6.875  | 146  | 881149   | 50.859  | ng    | 95       |
| 14) 1,2-Dichlorobenzene       | 7.034  | 146  | 790672   | 47.081  | ng    | 96       |
| 15) Benzyl Alcohol            | 6.998  | 79   | 833213   | 51.660  | ng    | 97       |
| 16) 2,2'-oxybis(1-Chloropr... | 7.134  | 45   | 1770455  | 43.700  | ng    | 99       |
| 17) 2-Methylphenol            | 7.110  | 107  | 738993   | 54.569  | ng    | 96       |
| 18) Hexachloroethane          | 7.369  | 117  | 340816   | 51.885  | ng    | 95       |
| 19) n-Nitroso-di-n-propyla... | 7.281  | 70   | 647759   | 53.216  | ng    | # 81     |
| 22) Acetophenone              | 7.275  | 105  | 1121663  | 50.925  | ng    | 92       |
| 24) Nitrobenzene              | 7.445  | 77   | 971800   | 56.989  | ng    | # 88     |
| 25) Isophorone                | 7.686  | 82   | 1795123  | 57.075  | ng    | # 87     |
| 26) 2-Nitrophenol             | 7.757  | 139  | 426197   | 77.085  | ng    | # 72     |
| 27) 2,4-Dimethylphenol        | 7.792  | 122  | 681746m  | 55.530  | ng    |          |
| 28) bis(2-Chloroethoxy)met... | 7.892  | 93   | 1098332  | 55.661  | ng    | # 97     |
| 29) 2,4-Dichlorophenol        | 7.998  | 162  | 669907   | 55.664  | ng    | 90       |
| 30) 1,2,4-Trichlorobenzene    | 8.081  | 180  | 749953   | 53.131  | ng    | 99       |
| 31) Naphthalene               | 8.163  | 128  | 2253065  | 51.941  | ng    | 99       |
| 32) Benzoic acid              | 7.933  | 122  | 620904   | 83.513  | ng    | # 84     |
| 33) 4-Chloroaniline           | 8.210  | 127  | 993835   | 55.598  | ng    | # 88     |
| 34) Hexachlorobutadiene       | 8.281  | 225  | 460359   | 51.894  | ng    | 99       |
| 35) Caprolactam               | 8.598  | 113  | 242041m  | 61.436  | ng    |          |
| 36) 4-Chloro-3-methylphenol   | 8.692  | 107  | 785849   | 55.795  | ng    | 90       |
| 37) 2-Methylnaphthalene       | 8.851  | 142  | 1474335  | 52.803  | ng    | 91       |
| 38) 1-Methylnaphthalene       | 8.951  | 142  | 1421707  | 52.757  | ng    | 90       |
| 40) 1,2,4,5-Tetrachloroben... | 9.022  | 216  | 688063   | 52.195  | ng    | 99       |
| 41) Hexachlorocyclopentadiene | 9.010  | 237  | 384945   | 54.067  | ng    | 97       |
| 43) 2,4,6-Trichlorophenol     | 9.128  | 196  | 500289   | 55.607  | ng    | 93       |
| 44) 2,4,5-Trichlorophenol     | 9.169  | 196  | 529068   | 55.739  | ng    | 97       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF102921\  
 Data File : BF126007.D  
 Acq On : 29 Oct 2021 16:08  
 Operator : JU/CG  
 Sample : SSTDICC060  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SSTDICC060

Manual Integrations  
 APPROVED

mohammad  
 11/1/2021 1:37:51 PM

Quant Time: Oct 29 17:20:42 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF102921.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Oct 29 17:15:51 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 46) 1,1'-Biphenyl             | 9.316  | 154  | 1760076  | 50.216 | ng    | 97       |
| 47) 2-Chloronaphthalene       | 9.339  | 162  | 1353018  | 51.200 | ng    | 94       |
| 48) 2-Nitroaniline            | 9.433  | 65   | 623385   | 66.044 | ng #  | 71       |
| 49) Acenaphthylene            | 9.757  | 152  | 2008976  | 50.327 | ng    | 96       |
| 50) Dimethylphthalate         | 9.622  | 163  | 1559198  | 52.600 | ng #  | 96       |
| 51) 2,6-Dinitrotoluene        | 9.680  | 165  | 368783   | 63.075 | ng #  | 83       |
| 52) Acenaphthene              | 9.927  | 154  | 1401263  | 54.015 | ng    | 98       |
| 53) 3-Nitroaniline            | 9.845  | 138  | 425649   | 63.506 | ng #  | 66       |
| 54) 2,4-Dinitrophenol         | 9.951  | 184  | 201179   | 99.362 | ng #  | 84       |
| 55) Dibenzofuran              | 10.098 | 168  | 1959434  | 50.569 | ng    | 97       |
| 56) 4-Nitrophenol             | 10.004 | 139  | 349174   | 67.039 | ng #  | 76       |
| 57) 2,4-Dinitrotoluene        | 10.080 | 165  | 484878   | 67.518 | ng #  | 89       |
| 58) Fluorene                  | 10.445 | 166  | 1374138  | 47.565 | ng    | 99       |
| 59) 2,3,4,6-Tetrachlorophenol | 10.216 | 232  | 445861   | 57.547 | ng #  | 86       |
| 60) Diethylphthalate          | 10.316 | 149  | 1554432  | 51.410 | ng    | 97       |
| 61) 4-Chlorophenyl-phenyle... | 10.433 | 204  | 666125   | 46.717 | ng    | 89       |
| 62) 4-Nitroaniline            | 10.463 | 138  | 411622   | 66.932 | ng    | 84       |
| 63) Azobenzene                | 10.598 | 77   | 1828229  | 52.424 | ng    | 93       |
| 65) 4,6-Dinitro-2-methylph... | 10.492 | 198  | 295111   | 88.747 | ng    | 88       |
| 66) n-Nitrosodiphenylamine    | 10.557 | 169  | 1321765  | 52.162 | ng    | 97       |
| 67) 4-Bromophenyl-phenylether | 10.922 | 248  | 485864   | 53.762 | ng    | 95       |
| 68) Hexachlorobenzene         | 10.992 | 284  | 515221   | 54.077 | ng #  | 83       |
| 69) Atrazine                  | 11.080 | 200  | 430980   | 53.703 | ng    | 91       |
| 70) Pentachlorophenol         | 11.180 | 266  | 326221   | 60.650 | ng    | 94       |
| 71) Phenanthrene              | 11.404 | 178  | 2182258  | 49.713 | ng    | 96       |
| 72) Anthracene                | 11.457 | 178  | 2186623  | 49.890 | ng    | 97       |
| 73) Carbazole                 | 11.610 | 167  | 2109614  | 50.633 | ng    | 98       |
| 74) Di-n-butylphthalate       | 11.939 | 149  | 2279146  | 51.345 | ng    | 98       |
| 75) Fluoranthene              | 12.586 | 202  | 2241392  | 50.650 | ng    | 96       |
| 78) Pyrene                    | 12.816 | 202  | 2311777  | 49.934 | ng    | 95       |
| 80) Butylbenzylphthalate      | 13.439 | 149  | 939941   | 62.413 | ng #  | 86       |
| 81) Benzo(a)anthracene        | 14.004 | 228  | 1708622  | 51.029 | ng    | 100      |
| 82) 3,3'-Dichlorobenzidine    | 13.968 | 252  | 602339   | 56.734 | ng #  | 99       |
| 83) Chrysene                  | 14.045 | 228  | 1747524  | 51.584 | ng    | 97       |
| 84) Bis(2-ethylhexyl)phtha... | 13.998 | 149  | 1016483  | 62.301 | ng    | 98       |
| 85) Di-n-octyl phthalate      | 14.610 | 149  | 1703440  | 67.393 | ng #  | 97       |
| 87) Indeno(1,2,3-cd)pyrene    | 16.939 | 276  | 1597143  | 53.386 | ng #  | 91       |
| 88) Benzo(b)fluoranthene      | 15.051 | 252  | 1473623  | 53.761 | ng #  | 96       |
| 89) Benzo(k)fluoranthene      | 15.080 | 252  | 1476052  | 56.061 | ng    | 99       |
| 90) Benzo(a)pyrene            | 15.415 | 252  | 1252865  | 56.618 | ng #  | 95       |
| 91) Dibenzo(a,h)anthracene    | 16.956 | 278  | 1284429  | 52.996 | ng #  | 95       |
| 92) Benzo(g,h,i)perylene      | 17.380 | 276  | 1356080  | 53.553 | ng #  | 90       |

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

