

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF110921\  
 Data File : BF126083.D  
 Acq On : 9 Nov 2021 15:42  
 Operator : JU/CG  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SSTDCCC040

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 11/09/2021  
 Supervised By :mohammad ahmed 11/11/2021

Quant Time: Nov 09 16:42:49 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF110321.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Nov 03 17:21:25 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 6.857  | 152  | 199483   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8            | 8.139  | 136  | 688975   | 20.000 | ng    | # 0.00   |
| 39) Acenaphthene-d10          | 9.892  | 164  | 421093   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10          | 11.374 | 188  | 771816   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12              | 14.015 | 240  | 563777   | 20.000 | ng    | # 0.00   |
| 86) Perylene-d12              | 15.474 | 264  | 380445   | 20.000 | ng    | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 5) 2-Fluorophenol             | 5.469  | 112  | 919221   | 79.133 | ng    | 0.00     |
| 7) Phenol-d6                  | 6.486  | 99   | 1261210  | 77.020 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 7.422  | 82   | 1274572  | 87.244 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 10.680 | 330  | 424225   | 87.909 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 9.216  | 172  | 2392902  | 76.792 | ng    | 0.00     |
| 79) Terphenyl-d14             | 12.963 | 244  | 2841095  | 81.263 | ng    | 0.00     |
| Target Compounds              |        |      |          |        |       |          |
|                               |        |      |          |        |       | Qvalue   |
| 2) 1,4-Dioxane                | 2.610  | 88   | 182398   | 37.763 | ng    | # 100    |
| 3) Pyridine                   | 3.357  | 79   | 487675   | 37.521 | ng    | 96       |
| 4) n-Nitrosodimethylamine     | 3.310  | 42   | 318557   | 34.951 | ng    | # 87     |
| 6) Aniline                    | 6.516  | 93   | 695531   | 38.520 | ng    | # 89     |
| 8) 2-Chlorophenol             | 6.639  | 128  | 496157   | 39.972 | ng    | 83       |
| 9) Benzaldehyde               | 6.404  | 77   | 351024   | 35.791 | ng    | 90       |
| 10) Phenol                    | 6.498  | 94   | 640345   | 38.244 | ng    | 81       |
| 11) bis(2-Chloroethyl)ether   | 6.592  | 93   | 471515   | 38.663 | ng    | 89       |
| 12) 1,3-Dichlorobenzene       | 6.798  | 146  | 560126m  | 39.622 | ng    |          |
| 13) 1,4-Dichlorobenzene       | 6.875  | 146  | 575500   | 39.923 | ng    | 97       |
| 14) 1,2-Dichlorobenzene       | 7.028  | 146  | 545911   | 39.308 | ng    | 96       |
| 15) Benzyl Alcohol            | 6.998  | 79   | 543359   | 38.854 | ng    | 93       |
| 16) 2,2'-oxybis(1-Chloropr... | 7.133  | 45   | 805633   | 35.926 | ng    | 93       |
| 17) 2-Methylphenol            | 7.104  | 107  | 412926   | 39.089 | ng    | # 89     |
| 18) Hexachloroethane          | 7.369  | 117  | 215460   | 39.784 | ng    | # 84     |
| 19) n-Nitroso-di-n-propyla... | 7.275  | 70   | 411362   | 38.106 | ng    | # 79     |
| 20) 3+4-Methylphenols         | 7.263  | 107  | 541151   | 37.891 | ng    | # 67     |
| 22) Acetophenone              | 7.269  | 105  | 767360   | 38.563 | ng    | 96       |
| 24) Nitrobenzene              | 7.439  | 77   | 599447   | 41.428 | ng    | # 86     |
| 25) Isophorone                | 7.680  | 82   | 1007190  | 38.292 | ng    | # 89     |
| 26) 2-Nitrophenol             | 7.751  | 139  | 247860   | 48.379 | ng    | # 60     |
| 27) 2,4-Dimethylphenol        | 7.786  | 122  | 370367   | 38.621 | ng    | # 75     |
| 28) bis(2-Chloroethoxy)met... | 7.886  | 93   | 599688   | 38.663 | ng    | # 96     |
| 29) 2,4-Dichlorophenol        | 7.992  | 162  | 449330   | 41.509 | ng    | 95       |
| 30) 1,2,4-Trichlorobenzene    | 8.080  | 180  | 543746   | 41.197 | ng    | 97       |
| 31) Naphthalene               | 8.163  | 128  | 1405701  | 39.391 | ng    | 99       |
| 32) Benzoic acid              | 7.916  | 122  | 275324   | 46.610 | ng    | # 72     |
| 33) 4-Chloroaniline           | 8.204  | 127  | 569846   | 39.955 | ng    | # 87     |
| 34) Hexachlorobutadiene       | 8.280  | 225  | 374937   | 41.267 | ng    | 99       |
| 35) Caprolactam               | 8.580  | 113  | 119436   | 39.622 | ng    | # 65     |
| 36) 4-Chloro-3-methylphenol   | 8.686  | 107  | 497454   | 39.852 | ng    | 89       |
| 37) 2-Methylnaphthalene       | 8.851  | 142  | 970566   | 39.379 | ng    | 98       |
| 38) 1-Methylnaphthalene       | 8.951  | 142  | 944267   | 39.553 | ng    | 98       |
| 40) 1,2,4,5-Tetrachloroben... | 9.016  | 216  | 560822   | 39.741 | ng    | 97       |
| 41) Hexachlorocyclopentadiene | 9.004  | 237  | 302999   | 41.324 | ng    | 98       |
| 43) 2,4,6-Trichlorophenol     | 9.127  | 196  | 368628   | 43.136 | ng    | 97       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF110921\  
 Data File : BF126083.D  
 Acq On : 9 Nov 2021 15:42  
 Operator : JU/CG  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SSTDCCC040

Quant Time: Nov 09 16:42:49 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF110321.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Nov 03 17:21:25 2021  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 11/09/2021  
 Supervised By :mohammad ahmed 11/11/2021

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 9.163  | 196  | 386291   | 41.843 | ng    | 93       |
| 46) 1,1'-Biphenyl             | 9.316  | 154  | 1261626  | 38.762 | ng    | 99       |
| 47) 2-Chloronaphthalene       | 9.339  | 162  | 996169   | 38.884 | ng    | 100      |
| 48) 2-Nitroaniline            | 9.433  | 65   | 345649   | 41.651 | ng    | # 79     |
| 49) Acenaphthylene            | 9.757  | 152  | 1497359  | 38.649 | ng    | 98       |
| 50) Dimethylphthalate         | 9.616  | 163  | 1179139  | 38.462 | ng    | # 98     |
| 51) 2,6-Dinitrotoluene        | 9.674  | 165  | 243655   | 46.368 | ng    | # 77     |
| 52) Acenaphthene              | 9.927  | 154  | 979963   | 40.103 | ng    | 96       |
| 53) 3-Nitroaniline            | 9.845  | 138  | 245624   | 45.596 | ng    | # 74     |
| 54) 2,4-Dinitrophenol         | 9.945  | 184  | 121532   | 52.790 | ng    | # 88     |
| 55) Dibenzofuran              | 10.098 | 168  | 1402698  | 38.131 | ng    | 98       |
| 56) 4-Nitrophenol             | 9.998  | 139  | 184269   | 43.342 | ng    | # 55     |
| 57) 2,4-Dinitrotoluene        | 10.074 | 165  | 331697   | 44.094 | ng    | # 85     |
| 58) Fluorene                  | 10.439 | 166  | 1108346  | 37.530 | ng    | 97       |
| 59) 2,3,4,6-Tetrachlorophenol | 10.216 | 232  | 331516   | 42.406 | ng    | # 87     |
| 60) Diethylphthalate          | 10.310 | 149  | 1178729  | 38.340 | ng    | 98       |
| 61) 4-Chlorophenyl-phenyle... | 10.433 | 204  | 614644   | 38.916 | ng    | 89       |
| 62) 4-Nitroaniline            | 10.457 | 138  | 247677   | 44.082 | ng    | 86       |
| 63) Azobenzene                | 10.592 | 77   | 1228653  | 36.805 | ng    | 92       |
| 65) 4,6-Dinitro-2-methylph... | 10.486 | 198  | 191682   | 50.960 | ng    | 91       |
| 66) n-Nitrosodiphenylamine    | 10.551 | 169  | 953862   | 38.809 | ng    | 96       |
| 67) 4-Bromophenyl-phenylether | 10.921 | 248  | 376742   | 40.300 | ng    | 97       |
| 68) Hexachlorobenzene         | 10.986 | 284  | 399574   | 40.916 | ng    | 92       |
| 69) Atrazine                  | 11.074 | 200  | 360471   | 41.463 | ng    | 93       |
| 70) Pentachlorophenol         | 11.180 | 266  | 232237   | 44.132 | ng    | 92       |
| 71) Phenanthrene              | 11.398 | 178  | 1635181  | 39.067 | ng    | 98       |
| 72) Anthracene                | 11.451 | 178  | 1641703  | 39.287 | ng    | 98       |
| 73) Carbazole                 | 11.604 | 167  | 1489451  | 38.430 | ng    | 99       |
| 74) Di-n-butylphthalate       | 11.939 | 149  | 1800701  | 39.874 | ng    | 99       |
| 75) Fluoranthene              | 12.586 | 202  | 1808746  | 38.806 | ng    | 96       |
| 77) Benzidine                 | 12.704 | 184  | 481066   | 25.526 | ng    | 96       |
| 78) Pyrene                    | 12.815 | 202  | 1811956  | 40.077 | ng    | 96       |
| 80) Butylbenzylphthalate      | 13.433 | 149  | 721553   | 43.814 | ng    | # 84     |
| 81) Benzo(a)anthracene        | 14.004 | 228  | 1507343  | 38.919 | ng    | 98       |
| 82) 3,3'-Dichlorobenzidine    | 13.962 | 252  | 452275   | 40.017 | ng    | 98       |
| 83) Chrysene                  | 14.039 | 228  | 1430293  | 39.689 | ng    | 98       |
| 84) Bis(2-ethylhexyl)phtha... | 13.998 | 149  | 956433   | 41.903 | ng    | # 99     |
| 85) Di-n-octyl phthalate      | 14.609 | 149  | 1323025  | 40.797 | ng    | # 99     |
| 87) Indeno(1,2,3-cd)pyrene    | 16.939 | 276  | 969190   | 43.040 | ng    | # 84     |
| 88) Benzo(b)fluoranthene      | 15.051 | 252  | 994855   | 40.202 | ng    | # 93     |
| 89) Benzo(k)fluoranthene      | 15.080 | 252  | 985624   | 41.059 | ng    | # 97     |
| 90) Benzo(a)pyrene            | 15.415 | 252  | 778860   | 39.950 | ng    | # 92     |
| 91) Dibenzo(a,h)anthracene    | 16.956 | 278  | 800625   | 43.288 | ng    | # 91     |
| 92) Benzo(g,h,i)perylene      | 17.380 | 276  | 779591   | 43.950 | ng    | # 82     |

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

