

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF010721\  
 Data File : BF122927.D  
 Acq On : 7 Jan 2021 16:13  
 Operator : JU/CG  
 Sample : SSTDICC010  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SSTDICC010

Quant Time: Jan 07 18:02:33 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF010721.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jan 07 17:57:24 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc  | Units | Dev(Min) |
|-------------------------------|--------|------|----------|-------|-------|----------|
| -----                         |        |      |          |       |       |          |
| Internal Standards            |        |      |          |       |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 6.916  | 152  | 173091   | 20.00 | ng    | 0.00     |
| 21) Naphthalene-d8            | 8.198  | 136  | 677279   | 20.00 | ng    | # 0.00   |
| 39) Acenaphthene-d10          | 9.963  | 164  | 375844   | 20.00 | ng    | 0.00     |
| 64) Phenanthrene-d10          | 11.451 | 188  | 718518   | 20.00 | ng    | 0.00     |
| 76) Chrysene-d12              | 14.098 | 240  | 588860   | 20.00 | ng    | # 0.00   |
| 86) Perylene-d12              | 15.592 | 264  | 556915   | 20.00 | ng    | 0.00     |
| System Monitoring Compounds   |        |      |          |       |       |          |
| 5) 2-Fluorophenol             | 5.522  | 112  | 280672   | 25.84 | ng    | -0.01    |
| 7) Phenol-d6                  | 6.522  | 99   | 380726   | 27.23 | ng    | -0.02    |
| 23) Nitrobenzene-d5           | 7.475  | 82   | 319449   | 24.25 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 10.745 | 330  | 97697    | 20.33 | ng    | -0.01    |
| 45) 2-Fluorobiphenyl          | 9.275  | 172  | 646879   | 24.02 | ng    | 0.00     |
| 79) Terphenyl-d14             | 13.039 | 244  | 791871   | 22.96 | ng    | -0.01    |
| Target Compounds              |        |      |          |       |       |          |
|                               |        |      |          |       |       | Qvalue   |
| 2) 1,4-Dioxane                | 2.722  | 88   | 62988    | 12.63 | ng    | 89       |
| 3) Pyridine                   | 3.475  | 79   | 174050   | 13.98 | ng    | 96       |
| 4) n-Nitrosodimethylamine     | 3.416  | 42   | 72100    | 14.74 | ng    | # 95     |
| 6) Aniline                    | 6.575  | 93   | 214167   | 13.40 | ng    | 98       |
| 8) 2-Chlorophenol             | 6.698  | 128  | 149853   | 12.71 | ng    | 99       |
| 9) Benzaldehyde               | 6.463  | 77   | 115430   | 15.46 | ng    | 97       |
| 10) Phenol                    | 6.539  | 94   | 185368   | 12.92 | ng    | 83       |
| 11) bis(2-Chloroethyl)ether   | 6.645  | 93   | 148772   | 13.80 | ng    | 98       |
| 12) 1,3-Dichlorobenzene       | 6.857  | 146  | 169743   | 12.04 | ng    | 98       |
| 13) 1,4-Dichlorobenzene       | 6.934  | 146  | 171875   | 12.15 | ng    | 98       |
| 14) 1,2-Dichlorobenzene       | 7.087  | 146  | 166038   | 12.55 | ng    | 98       |
| 15) Benzyl Alcohol            | 7.045  | 79   | 129029   | 13.79 | ng    | 99       |
| 16) 2,2'-oxybis(1-Chloropr... | 7.186  | 45   | 237111   | 17.21 | ng    | 94       |
| 17) 2-Methylphenol            | 7.151  | 107  | 123225   | 13.07 | ng    | 97       |
| 18) Hexachloroethane          | 7.434  | 117  | 63099    | 12.43 | ng    | 98       |
| 19) n-Nitroso-di-n-propyla... | 7.316  | 70   | 115883   | 14.97 | ng    | 94       |
| 20) 3+4-Methylphenols         | 7.304  | 107  | 163765   | 13.97 | ng    | # 66     |
| 22) Acetophenone              | 7.316  | 105  | 217096   | 12.99 | ng    | # 97     |
| 24) Nitrobenzene              | 7.492  | 77   | 159740   | 12.39 | ng    | 99       |
| 25) Isophorone                | 7.728  | 82   | 284828   | 13.08 | ng    | 97       |
| 26) 2-Nitrophenol             | 7.810  | 139  | 66603    | 10.12 | ng    | 98       |
| 27) 2,4-Dimethylphenol        | 7.839  | 122  | 125588   | 13.67 | ng    | 96       |
| 28) bis(2-Chloroethoxy)met... | 7.945  | 93   | 195071   | 13.09 | ng    | 99       |
| 29) 2,4-Dichlorophenol        | 8.045  | 162  | 123993   | 11.44 | ng    | 96       |
| 30) 1,2,4-Trichlorobenzene    | 8.139  | 180  | 143345   | 11.49 | ng    | 99       |
| 31) Naphthalene               | 8.222  | 128  | 438237   | 11.86 | ng    | 99       |
| 32) Benzoic acid              | 7.904  | 122  | 81879    | 13.95 | ng    | 98       |
| 33) 4-Chloroaniline           | 8.263  | 127  | 186017   | 12.15 | ng    | 97       |
| 34) Hexachlorobutadiene       | 8.345  | 225  | 83265    | 10.77 | ng    | 98       |
| 35) Caprolactam               | 8.598  | 113  | 41757    | 12.82 | ng    | # 84     |
| 36) 4-Chloro-3-methylphenol   | 8.733  | 107  | 129309   | 11.90 | ng    | 99       |
| 37) 2-Methylnaphthalene       | 8.916  | 142  | 300769   | 12.30 | ng    | 96       |
| 38) 1-Methylnaphthalene       | 9.016  | 142  | 285275   | 12.44 | ng    | 96       |
| 40) 1,2,4,5-Tetrachloroben... | 9.081  | 216  | 136321   | 11.07 | ng    | 99       |
| 41) Hexachlorocyclopentadiene | 9.075  | 237  | 71595    | 34.51 | ng    | 100      |
| 43) 2,4,6-Trichlorophenol     | 9.186  | 196  | 95027    | 11.74 | ng    | 94       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF010721\  
 Data File : BF122927.D  
 Acq On : 7 Jan 2021 16:13  
 Operator : JU/CG  
 Sample : SSTDICC010  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SSTDICC010

Quant Time: Jan 07 18:02:33 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF010721.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jan 07 17:57:24 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc  | Units | Dev(Min) |
|-------------------------------|--------|------|----------|-------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 9.222  | 196  | 92884    | 11.13 | ng    | 94       |
| 46) 1,1'-Biphenyl             | 9.380  | 154  | 364768   | 11.77 | ng    | 97       |
| 47) 2-Chloronaphthalene       | 9.404  | 162  | 290854   | 11.30 | ng    | 96       |
| 48) 2-Nitroaniline            | 9.492  | 65   | 82814    | 11.35 | ng    | 93       |
| 49) Acenaphthylene            | 9.822  | 152  | 431052   | 11.48 | ng    | 99       |
| 50) Dimethylphthalate         | 9.675  | 163  | 328848   | 11.06 | ng    | 99       |
| 51) 2,6-Dinitrotoluene        | 9.733  | 165  | 65384    | 10.26 | ng    | 94       |
| 52) Acenaphthene              | 9.992  | 154  | 284697   | 12.52 | ng    | 99       |
| 53) 3-Nitroaniline            | 9.898  | 138  | 79018    | 10.84 | ng    | # 87     |
| 54) 2,4-Dinitrophenol         | 10.004 | 184  | 20661    | 8.77  | ng    | # 14     |
| 55) Dibenzofuran              | 10.163 | 168  | 402725   | 11.91 | ng    | 97       |
| 56) 4-Nitrophenol             | 10.039 | 139  | 59046    | 14.57 | ng    | # 82     |
| 57) 2,4-Dinitrotoluene        | 10.133 | 165  | 83057    | 10.00 | ng    | 96       |
| 58) Fluorene                  | 10.510 | 166  | 325379   | 12.01 | ng    | 99       |
| 59) 2,3,4,6-Tetrachlorophenol | 10.275 | 232  | 84685    | 11.66 | ng    | 94       |
| 60) Diethylphthalate          | 10.380 | 149  | 331764   | 11.46 | ng    | 99       |
| 61) 4-Chlorophenyl-phenyle... | 10.504 | 204  | 160847   | 11.84 | ng    | 92       |
| 62) 4-Nitroaniline            | 10.510 | 138  | 76653    | 10.05 | ng    | 95       |
| 63) Azobenzene                | 10.663 | 77   | 333716   | 13.08 | ng    | 98       |
| 65) 4,6-Dinitro-2-methylph... | 10.539 | 198  | 28207    | 6.85  | ng    | 86       |
| 66) n-Nitrosodiphenylamine    | 10.616 | 169  | 283651   | 11.50 | ng    | 99       |
| 67) 4-Bromophenyl-phenylether | 10.992 | 248  | 99738    | 10.74 | ng    | 98       |
| 68) Hexachlorobenzene         | 11.063 | 284  | 105278   | 10.16 | ng    | 94       |
| 69) Atrazine                  | 11.139 | 200  | 90970    | 11.57 | ng    | 99       |
| 70) Pentachlorophenol         | 11.245 | 266  | 61014    | 14.93 | ng    | 98       |
| 71) Phenanthrene              | 11.474 | 178  | 489004   | 11.52 | ng    | 99       |
| 72) Anthracene                | 11.527 | 178  | 489631   | 11.56 | ng    | 98       |
| 73) Carbazole                 | 11.674 | 167  | 454380   | 11.85 | ng    | 98       |
| 74) Di-n-butylphthalate       | 12.016 | 149  | 549965   | 11.56 | ng    | 98       |
| 75) Fluoranthene              | 12.669 | 202  | 531373   | 11.77 | ng    | 98       |
| 77) Benzidine                 | 12.780 | 184  | 199013   | 12.22 | ng    | 99       |
| 78) Pyrene                    | 12.898 | 202  | 537128   | 12.53 | ng    | 99       |
| 80) Butylbenzylphthalate      | 13.521 | 149  | 228585   | 11.98 | ng    | 95       |
| 81) Benzo(a)anthracene        | 14.086 | 228  | 473634   | 11.86 | ng    | 98       |
| 82) 3,3'-Dichlorobenzidine    | 14.051 | 252  | 158797   | 11.62 | ng    | 98       |
| 83) Chrysene                  | 14.127 | 228  | 448432   | 11.61 | ng    | 99       |
| 84) Bis(2-ethylhexyl)phtha... | 14.086 | 149  | 315455   | 11.86 | ng    | 97       |
| 85) Di-n-octyl phthalate      | 14.698 | 149  | 525869   | 11.97 | ng    | 97       |
| 87) Indeno(1,2,3-cd)pyrene    | 17.086 | 276  | 415903   | 11.38 | ng    | 97       |
| 88) Benzo(b)fluoranthene      | 15.151 | 252  | 436327   | 11.56 | ng    | 99       |
| 89) Benzo(k)fluoranthene      | 15.180 | 252  | 426112   | 11.92 | ng    | 99       |
| 90) Benzo(a)pyrene            | 15.521 | 252  | 396752   | 11.71 | ng    | 97       |
| 91) Dibenzo(a,h)anthracene    | 17.109 | 278  | 345414   | 11.47 | ng    | 98       |
| 92) Benzo(g,h,i)perylene      | 17.539 | 276  | 329445   | 11.41 | ng    | 97       |

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

Quant Time: Jan 07 18:02:33 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF010721.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Thu Jan 07 17:57:24 2021  
Response via : Initial Calibration

