

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF013024\  
 Data File : BF137194.D  
 Acq On : 30 Jan 2024 12:55  
 Operator : CG\JU  
 Sample : SSTDICC020  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SSTDICC020

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 01/31/2024  
 Supervised By :Sohil Jodhani 01/31/2024

Quant Time: Jan 31 04:38:30 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF013024.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Jan 31 04:27:00 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 6.792  | 152  | 51095    | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8            | 8.069  | 136  | 206968   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10          | 9.828  | 164  | 122293   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10          | 11.322 | 188  | 240682   | 20.000 | ng    | # 0.00   |
| 76) Chrysene-d12              | 13.974 | 240  | 147372   | 20.000 | ng    | 0.00     |
| 86) Perylene-d12              | 15.451 | 264  | 114296   | 20.000 | ng    | 0.00     |
|                               |        |      |          |        |       |          |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 5) 2-Fluorophenol             | 5.416  | 112  | 133532   | 43.123 | ng    | 0.00     |
| 7) Phenol-d6                  | 6.422  | 99   | 191584   | 43.354 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 7.351  | 82   | 196550   | 42.518 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 10.616 | 330  | 58344    | 42.244 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 9.145  | 172  | 382283   | 42.972 | ng    | 0.00     |
| 79) Terphenyl-d14             | 12.916 | 244  | 469371   | 43.298 | ng    | 0.00     |
|                               |        |      |          |        |       |          |
| Target Compounds              |        |      |          |        |       | Qvalue   |
| 2) 1,4-Dioxane                | 2.546  | 88   | 35792    | 21.602 | ng    | 98       |
| 3) Pyridine                   | 3.299  | 79   | 59213m   | 19.650 | ng    |          |
| 4) n-Nitrosodimethylamine     | 3.252  | 42   | 28348m   | 20.758 | ng    |          |
| 6) Aniline                    | 6.457  | 93   | 103090   | 21.793 | ng    | 98       |
| 8) 2-Chlorophenol             | 6.575  | 128  | 73969    | 21.364 | ng    | 98       |
| 9) Benzaldehyde               | 6.345  | 77   | 42116    | 26.352 | ng    | 98       |
| 10) Phenol                    | 6.434  | 94   | 106364   | 21.563 | ng    | 99       |
| 11) bis(2-Chloroethyl)ether   | 6.528  | 93   | 66000    | 20.022 | ng    | 98       |
| 12) 1,3-Dichlorobenzene       | 6.734  | 146  | 87235    | 21.490 | ng    | 98       |
| 13) 1,4-Dichlorobenzene       | 6.810  | 146  | 87684    | 20.886 | ng    | 99       |
| 14) 1,2-Dichlorobenzene       | 6.963  | 146  | 84502    | 21.527 | ng    | 98       |
| 15) Benzyl Alcohol            | 6.934  | 79   | 65814    | 20.963 | ng    | 95       |
| 16) 2,2'-oxybis(1-Chloropr... | 7.063  | 45   | 129053   | 21.733 | ng    | 99       |
| 17) 2-Methylphenol            | 7.040  | 107  | 57394    | 19.842 | ng    | 98       |
| 18) Hexachloroethane          | 7.304  | 117  | 32909    | 21.734 | ng    | 98       |
| 19) n-Nitroso-di-n-propyla... | 7.198  | 70   | 66231    | 21.859 | ng    | 94       |
| 20) 3+4-Methylphenols         | 7.198  | 107  | 80469    | 21.976 | ng    | # 74     |
| 22) Acetophenone              | 7.198  | 105  | 122503   | 21.986 | ng    | # 99     |
| 24) Nitrobenzene              | 7.369  | 77   | 93074    | 20.744 | ng    | 98       |
| 25) Isophorone                | 7.610  | 82   | 179440   | 21.030 | ng    | 98       |
| 26) 2-Nitrophenol             | 7.687  | 139  | 31757    | 19.799 | ng    | # 88     |
| 27) 2,4-Dimethylphenol        | 7.722  | 122  | 50679    | 20.780 | ng    | 99       |
| 28) bis(2-Chloroethoxy)met... | 7.828  | 93   | 98610    | 21.762 | ng    | 99       |
| 29) 2,4-Dichlorophenol        | 7.928  | 162  | 65467    | 21.638 | ng    | 97       |
| 30) 1,2,4-Trichlorobenzene    | 8.010  | 180  | 78968    | 21.012 | ng    | 98       |
| 31) Naphthalene               | 8.092  | 128  | 235056   | 21.008 | ng    | 100      |
| 32) Benzoic acid              | 7.816  | 122  | 21169m   | 19.723 | ng    |          |
| 33) 4-Chloroaniline           | 8.145  | 127  | 81131    | 21.080 | ng    | 97       |
| 34) Hexachlorobutadiene       | 8.210  | 225  | 53481    | 21.402 | ng    | 98       |
| 35) Caprolactam               | 8.492  | 113  | 19808    | 21.055 | ng    | 97       |
| 36) 4-Chloro-3-methylphenol   | 8.622  | 107  | 76386    | 21.375 | ng    | 96       |
| 37) 2-Methylnaphthalene       | 8.786  | 142  | 159208   | 21.224 | ng    | 100      |
| 38) 1-Methylnaphthalene       | 8.881  | 142  | 145231   | 20.845 | ng    | 99       |
| 40) 1,2,4,5-Tetrachloroben... | 8.945  | 216  | 85673    | 20.937 | ng    | 99       |
| 41) Hexachlorocyclopentadiene | 8.934  | 237  | 39619    | 20.328 | ng    | 98       |
| 43) 2,4,6-Trichlorophenol     | 9.063  | 196  | 47481    | 20.457 | ng    | 95       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF013024\  
 Data File : BF137194.D  
 Acq On : 30 Jan 2024 12:55  
 Operator : CG\JU  
 Sample : SSTDICC020  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SSTDICC020

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 01/31/2024  
 Supervised By :Sohil Jodhani 01/31/2024

Quant Time: Jan 31 04:38:30 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF013024.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Jan 31 04:27:00 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 9.098  | 196  | 54987    | 21.674 | ng    | 98       |
| 46) 1,1'-Biphenyl             | 9.251  | 154  | 206798   | 21.254 | ng    | 98       |
| 47) 2-Chloronaphthalene       | 9.275  | 162  | 151354   | 21.019 | ng    | 97       |
| 48) 2-Nitroaniline            | 9.369  | 65   | 47804    | 20.511 | ng    | 90       |
| 49) Acenaphthylene            | 9.686  | 152  | 244847   | 21.240 | ng    | 100      |
| 50) Dimethylphthalate         | 9.551  | 163  | 176219   | 21.329 | ng    | 99       |
| 51) 2,6-Dinitrotoluene        | 9.610  | 165  | 37681    | 21.388 | ng    | 95       |
| 52) Acenaphthene              | 9.857  | 154  | 141011   | 20.810 | ng    | 98       |
| 53) 3-Nitroaniline            | 9.781  | 138  | 33627    | 21.210 | ng    | 88       |
| 54) 2,4-Dinitrophenol         | 9.886  | 184  | 6394     | 16.525 | ng #  | 84       |
| 55) Dibenzofuran              | 10.033 | 168  | 229820   | 21.321 | ng    | 99       |
| 56) 4-Nitrophenol             | 9.939  | 139  | 15904    | 18.004 | ng    | 98       |
| 57) 2,4-Dinitrotoluene        | 10.016 | 165  | 48651    | 21.283 | ng    | 84       |
| 58) Fluorene                  | 10.375 | 166  | 179606   | 21.255 | ng    | 99       |
| 59) 2,3,4,6-Tetrachlorophenol | 10.151 | 232  | 46971m   | 19.792 | ng    |          |
| 60) Diethylphthalate          | 10.257 | 149  | 172687   | 21.228 | ng    | 99       |
| 61) 4-Chlorophenyl-phenyle... | 10.375 | 204  | 90182    | 21.470 | ng    | 98       |
| 62) 4-Nitroaniline            | 10.398 | 138  | 31118    | 20.425 | ng    | 95       |
| 63) Azobenzene                | 10.533 | 77   | 189247   | 21.006 | ng    | 98       |
| 65) 4,6-Dinitro-2-methylph... | 10.416 | 198  | 19064    | 18.242 | ng    | 98       |
| 66) n-Nitrosodiphenylamine    | 10.492 | 169  | 155880   | 21.300 | ng    | 98       |
| 67) 4-Bromophenyl-phenylether | 10.863 | 248  | 55876    | 21.113 | ng #  | 92       |
| 68) Hexachlorobenzene         | 10.922 | 284  | 61981    | 20.968 | ng    | 95       |
| 69) Atrazine                  | 11.022 | 200  | 51052    | 21.132 | ng    | 99       |
| 70) Pentachlorophenol         | 11.122 | 266  | 33071    | 20.919 | ng    | 99       |
| 71) Phenanthrene              | 11.339 | 178  | 266372   | 21.141 | ng    | 99       |
| 72) Anthracene                | 11.392 | 178  | 277960   | 21.165 | ng    | 99       |
| 73) Carbazole                 | 11.551 | 167  | 233886   | 21.310 | ng    | 99       |
| 74) Di-n-butylphthalate       | 11.892 | 149  | 250723   | 21.133 | ng    | 99       |
| 75) Fluoranthene              | 12.533 | 202  | 289354   | 21.946 | ng    | 98       |
| 77) Benzidine                 | 12.669 | 184  | 56103    | 22.670 | ng    | 98       |
| 78) Pyrene                    | 12.763 | 202  | 299067   | 21.117 | ng    | 100      |
| 80) Butylbenzylphthalate      | 13.404 | 149  | 75245    | 20.769 | ng    | 97       |
| 81) Benzo(a)anthracene        | 13.963 | 228  | 221691   | 21.241 | ng    | 99       |
| 82) 3,3'-Dichlorobenzidine    | 13.933 | 252  | 59822    | 21.038 | ng    | 97       |
| 83) Chrysene                  | 13.998 | 228  | 214440   | 20.899 | ng    | 99       |
| 84) Bis(2-ethylhexyl)phtha... | 13.968 | 149  | 85501    | 20.285 | ng    | 99       |
| 85) Di-n-octyl phthalate      | 14.586 | 149  | 115873   | 18.261 | ng    | 99       |
| 87) Indeno(1,2,3-cd)pyrene    | 16.945 | 276  | 177849   | 22.148 | ng    | 99       |
| 88) Benzo(b)fluoranthene      | 15.015 | 252  | 150865m  | 21.463 | ng    |          |
| 89) Benzo(k)fluoranthene      | 15.045 | 252  | 156917   | 21.102 | ng    | 100      |
| 90) Benzo(a)pyrene            | 15.392 | 252  | 138461   | 20.932 | ng    | 98       |
| 91) Dibenzo(a,h)anthracene    | 16.974 | 278  | 142558   | 22.156 | ng    | 98       |
| 92) Benzo(g,h,i)perylene      | 17.398 | 276  | 140939   | 22.318 | ng    | 99       |

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

