

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP011123\  
 Data File : BP013439.D  
 Acq On : 11 Jan 2023 12:19  
 Operator : CG/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTDCCC040

Quant Time: Jan 12 02:23:15 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP121322.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jan 12 02:16:32 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.916  | 152  | 395258   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8            | 10.728 | 136  | 1523088  | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10          | 14.563 | 164  | 987069   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10          | 17.316 | 188  | 2222690  | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12              | 21.410 | 240  | 2228980  | 20.000 | ng    | 0.00     |
| 86) Perylene-d12              | 23.868 | 264  | 2434392  | 20.000 | ng    | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 5) 2-Fluorophenol             | 5.469  | 112  | 1785838  | 82.553 | ng    | 0.00     |
| 7) Phenol-d6                  | 7.075  | 99   | 2406426  | 85.181 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 9.087  | 82   | 2412671  | 86.842 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 16.057 | 330  | 1090111  | 91.599 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 13.181 | 172  | 5760593  | 79.241 | ng    | 0.00     |
| 79) Terphenyl-d14             | 19.874 | 244  | 9251762  | 81.888 | ng    | 0.00     |
| Target Compounds              |        |      |          |        |       |          |
|                               |        |      |          |        |       | Qvalue   |
| 2) 1,4-Dioxane                | 3.340  | 88   | 391150   | 40.808 | ng    | 92       |
| 3) Pyridine                   | 3.752  | 79   | 1163898  | 42.758 | ng    | 92       |
| 4) n-Nitrosodimethylamine     | 3.664  | 42   | 564152   | 46.246 | ng    | 93       |
| 6) Aniline                    | 7.240  | 93   | 1458220  | 42.625 | ng    | 98       |
| 8) 2-Chlorophenol             | 7.475  | 128  | 1031054  | 40.430 | ng    | 97       |
| 9) Benzaldehyde               | 7.052  | 77   | 635472   | 37.489 | ng    | 95       |
| 10) Phenol                    | 7.105  | 94   | 1337053  | 42.282 | ng    | 96       |
| 11) bis(2-Chloroethyl)ether   | 7.340  | 93   | 1056772  | 41.510 | ng    | 96       |
| 12) 1,3-Dichlorobenzene       | 7.805  | 146  | 1189840  | 40.101 | ng    | 98       |
| 13) 1,4-Dichlorobenzene       | 7.952  | 146  | 1213598  | 40.161 | ng    | 99       |
| 14) 1,2-Dichlorobenzene       | 8.269  | 146  | 1157624  | 40.363 | ng    | 99       |
| 15) Benzyl Alcohol            | 8.157  | 79   | 894799   | 43.662 | ng    | 97       |
| 16) 2,2'-oxybis(1-Chloropr... | 8.452  | 45   | 1690091  | 45.835 | ng    | 97       |
| 17) 2-Methylphenol            | 8.363  | 107  | 888756   | 42.521 | ng    | 99       |
| 18) Hexachloroethane          | 8.999  | 117  | 416148   | 40.666 | ng    | 97       |
| 19) n-Nitroso-di-n-propyla... | 8.728  | 70   | 834364   | 45.475 | ng    | 95       |
| 20) 3+4-Methylphenols         | 8.693  | 107  | 1230522  | 43.303 | ng    | 98       |
| 22) Acetophenone              | 8.746  | 105  | 1577880  | 41.632 | ng    | # 96     |
| 24) Nitrobenzene              | 9.128  | 77   | 1240004  | 42.825 | ng    | 99       |
| 25) Isophorone                | 9.657  | 82   | 2096660  | 44.434 | ng    | 99       |
| 26) 2-Nitrophenol             | 9.840  | 139  | 553565   | 38.218 | ng    | 94       |
| 27) 2,4-Dimethylphenol        | 9.899  | 122  | 832151   | 41.679 | ng    | 98       |
| 28) bis(2-Chloroethoxy)met... | 10.134 | 93   | 1276877  | 42.201 | ng    | 99       |
| 29) 2,4-Dichlorophenol        | 10.375 | 162  | 1031584  | 41.865 | ng    | 99       |
| 30) 1,2,4-Trichlorobenzene    | 10.587 | 180  | 1152142  | 39.615 | ng    | 99       |
| 31) Naphthalene               | 10.775 | 128  | 3382042  | 40.553 | ng    | 100      |
| 32) Benzoic acid              | 10.028 | 122  | 711739   | 42.052 | ng    | 97       |
| 33) 4-Chloroaniline           | 10.893 | 127  | 1343895  | 43.208 | ng    | 98       |
| 34) Hexachlorobutadiene       | 11.057 | 225  | 732388   | 39.475 | ng    | 99       |
| 35) Caprolactam               | 11.687 | 113  | 317860   | 49.961 | ng    | # 83     |
| 36) 4-Chloro-3-methylphenol   | 12.010 | 107  | 1055851  | 45.876 | ng    | 97       |
| 37) 2-Methylnaphthalene       | 12.387 | 142  | 2378718  | 42.412 | ng    | 99       |
| 38) 1-Methylnaphthalene       | 12.604 | 142  | 2327481  | 42.615 | ng    | 99       |
| 40) 1,2,4,5-Tetrachloroben... | 12.751 | 216  | 1329037  | 37.742 | ng    | 100      |
| 41) Hexachlorocyclopentadiene | 12.728 | 237  | 614438   | 34.605 | ng    | 99       |
| 43) 2,4,6-Trichlorophenol     | 12.992 | 196  | 904630   | 40.738 | ng    | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP011123\  
 Data File : BP013439.D  
 Acq On : 11 Jan 2023 12:19  
 Operator : CG/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTDCCC040

Quant Time: Jan 12 02:23:15 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP121322.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jan 12 02:16:32 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 13.063 | 196  | 998554   | 41.622 | ng    | 98       |
| 46) 1,1'-Biphenyl             | 13.392 | 154  | 3059621  | 39.203 | ng    | 99       |
| 47) 2-Chloronaphthalene       | 13.440 | 162  | 2424252  | 39.081 | ng    | 100      |
| 48) 2-Nitroaniline            | 13.645 | 65   | 750653   | 45.013 | ng    | 92       |
| 49) Acenaphthylene            | 14.287 | 152  | 3828617  | 41.959 | ng    | 100      |
| 50) Dimethylphthalate         | 14.016 | 163  | 3180645  | 44.054 | ng    | 99       |
| 51) 2,6-Dinitrotoluene        | 14.139 | 165  | 681680   | 46.140 | ng    | 94       |
| 52) Acenaphthene              | 14.628 | 154  | 2329537  | 41.067 | ng    | 100      |
| 53) 3-Nitroaniline            | 14.475 | 138  | 718298   | 44.277 | ng    | 97       |
| 54) 2,4-Dinitrophenol         | 14.681 | 184  | 395534   | 42.176 | ng    | 95       |
| 55) Dibenzofuran              | 14.963 | 168  | 3818355  | 42.062 | ng    | 99       |
| 56) 4-Nitrophenol             | 14.781 | 139  | 594034   | 48.113 | ng    | 95       |
| 57) 2,4-Dinitrotoluene        | 14.928 | 165  | 913610   | 45.639 | ng    | 91       |
| 58) Fluorene                  | 15.610 | 166  | 3099394  | 44.112 | ng    | 100      |
| 59) 2,3,4,6-Tetrachlorophenol | 15.186 | 232  | 925462   | 45.127 | ng    | 99       |
| 60) Diethylphthalate          | 15.381 | 149  | 3108948  | 46.840 | ng    | 99       |
| 61) 4-Chlorophenyl-phenyle... | 15.604 | 204  | 1577818  | 43.391 | ng    | 98       |
| 62) 4-Nitroaniline            | 15.639 | 138  | 736862   | 47.423 | ng    | 96       |
| 63) Azobenzene                | 15.898 | 77   | 2950057  | 47.826 | ng    | 96       |
| 65) 4,6-Dinitro-2-methylph... | 15.692 | 198  | 566938   | 39.705 | ng    | 89       |
| 66) n-Nitrosodiphenylamine    | 15.822 | 169  | 2677420  | 38.581 | ng    | 99       |
| 67) 4-Bromophenyl-phenylether | 16.504 | 248  | 1000063  | 38.051 | ng    | 97       |
| 68) Hexachlorobenzene         | 16.616 | 284  | 1162156  | 37.765 | ng    | 97       |
| 69) Atrazine                  | 16.775 | 200  | 865638   | 44.458 | ng    | 99       |
| 70) Pentachlorophenol         | 16.969 | 266  | 746870   | 41.211 | ng    | 98       |
| 71) Phenanthrene              | 17.363 | 178  | 5069417  | 40.082 | ng    | 100      |
| 72) Anthracene                | 17.451 | 178  | 4857188  | 40.907 | ng    | 99       |
| 73) Carbazole                 | 17.727 | 167  | 4482953  | 43.067 | ng    | 100      |
| 74) Di-n-butylphthalate       | 18.269 | 149  | 5273361  | 46.059 | ng    | 100      |
| 75) Fluoranthene              | 19.327 | 202  | 6139876  | 45.102 | ng    | 100      |
| 77) Benzidine                 | 19.510 | 184  | 689666   | 21.783 | ng    | 99       |
| 78) Pyrene                    | 19.680 | 202  | 6338638  | 39.576 | ng    | 100      |
| 80) Butylbenzylphthalate      | 20.545 | 149  | 2418346  | 42.417 | ng    | 97       |
| 81) Benzo(a)anthracene        | 21.392 | 228  | 6501881  | 42.676 | ng    | 99       |
| 82) 3,3'-Dichlorobenzidine    | 21.321 | 252  | 1963034  | 43.269 | ng    | 99       |
| 83) Chrysene                  | 21.445 | 228  | 6203933  | 41.746 | ng    | 100      |
| 84) Bis(2-ethylhexyl)phtha... | 21.304 | 149  | 3498699  | 44.520 | ng    | 99       |
| 85) Di-n-octyl phthalate      | 22.245 | 149  | 5854952  | 45.045 | ng    | 97       |
| 87) Indeno(1,2,3-cd)pyrene    | 26.445 | 276  | 7714961  | 38.786 | ng    | 99       |
| 88) Benzo(b)fluoranthene      | 23.115 | 252  | 6295894  | 40.168 | ng    | 100      |
| 89) Benzo(k)fluoranthene      | 23.163 | 252  | 6473386  | 40.826 | ng    | 99       |
| 90) Benzo(a)pyrene            | 23.757 | 252  | 5313280  | 40.394 | ng    | 99       |
| 91) Dibenzo(a,h)anthracene    | 26.456 | 278  | 6456228  | 39.055 | ng    | 99       |
| 92) Benzo(g,h,i)perylene      | 27.233 | 276  | 6363371  | 38.851 | ng    | 100      |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP011123\  
 Data File : BP013439.D  
 Acq On : 11 Jan 2023 12:19  
 Operator : CG/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTDCCC040

Quant Time: Jan 12 02:23:15 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP121322.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Jan 12 02:16:32 2023  
 Response via : Initial Calibration

