

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP052423\  
 Data File : BP015276.D  
 Acq On : 24 May 2023 17:12  
 Operator : CG/JU  
 Sample : SSTDICC020  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTDICC020

Manual Integrations  
 APPROVED

Reviewed By : Christian Giraldo 05/25/2023  
 Supervised By : Jagrut Upadhyay 05/25/2023

Quant Time: May 25 01:28:32 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP052423.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu May 25 01:19:28 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.110  | 152  | 155349   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8            | 10.940 | 136  | 582156   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10          | 14.751 | 164  | 334654   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10          | 17.504 | 188  | 707402   | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12              | 21.586 | 240  | 652510   | 20.000 | ng    | 0.00     |
| 86) Perylene-d12              | 24.139 | 264  | 748572   | 20.000 | ng    | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 5) 2-Fluorophenol             | 5.634  | 112  | 392463   | 41.863 | ng    | 0.00     |
| 7) Phenol-d6                  | 7.258  | 99   | 509653   | 41.482 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 9.287  | 82   | 496568   | 41.648 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 16.239 | 330  | 184855   | 40.648 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 13.375 | 172  | 1063416  | 42.662 | ng    | 0.00     |
| 79) Terphenyl-d14             | 20.039 | 244  | 1488575  | 42.910 | ng    | 0.00     |
| Target Compounds              |        |      |          |        |       |          |
|                               |        |      |          |        |       | Qvalue   |
| 2) 1,4-Dioxane                | 3.452  | 88   | 89729    | 21.609 | ng    | 98       |
| 3) Pyridine                   | 3.875  | 79   | 220892   | 20.814 | ng    | 96       |
| 4) n-Nitrosodimethylamine     | 3.781  | 42   | 102746   | 20.746 | ng    | # 94     |
| 6) Aniline                    | 7.428  | 93   | 319459   | 20.572 | ng    | 100      |
| 8) 2-Chlorophenol             | 7.669  | 128  | 205927   | 21.020 | ng    | 99       |
| 9) Benzaldehyde               | 7.240  | 77   | 141337   | 21.159 | ng    | 99       |
| 10) Phenol                    | 7.287  | 94   | 253951   | 20.639 | ng    | 99       |
| 11) bis(2-Chloroethyl)ether   | 7.528  | 93   | 215775   | 20.922 | ng    | 99       |
| 12) 1,3-Dichlorobenzene       | 7.999  | 146  | 229202   | 21.025 | ng    | 99       |
| 13) 1,4-Dichlorobenzene       | 8.146  | 146  | 230764   | 21.030 | ng    | 99       |
| 14) 1,2-Dichlorobenzene       | 8.469  | 146  | 220743   | 21.085 | ng    | 98       |
| 15) Benzyl Alcohol            | 8.352  | 79   | 169326   | 19.745 | ng    | 99       |
| 16) 2,2'-oxybis(1-Chloropr... | 8.646  | 45   | 280224m  | 21.066 | ng    |          |
| 17) 2-Methylphenol            | 8.557  | 107  | 181831   | 20.765 | ng    | 99       |
| 18) Hexachloroethane          | 9.199  | 117  | 86909    | 21.015 | ng    | 93       |
| 19) n-Nitroso-di-n-propyla... | 8.922  | 70   | 163536   | 20.798 | ng    | 98       |
| 20) 3+4-Methylphenols         | 8.887  | 107  | 240932   | 20.367 | ng    | 97       |
| 22) Acetophenone              | 8.946  | 105  | 313208   | 20.714 | ng    | # 99     |
| 24) Nitrobenzene              | 9.328  | 77   | 245366   | 20.667 | ng    | 95       |
| 25) Isophorone                | 9.857  | 82   | 429445   | 20.071 | ng    | 99       |
| 26) 2-Nitrophenol             | 10.046 | 139  | 103362   | 19.836 | ng    | 96       |
| 27) 2,4-Dimethylphenol        | 10.099 | 122  | 178034   | 20.379 | ng    | 95       |
| 28) bis(2-Chloroethoxy)met... | 10.340 | 93   | 270570   | 20.518 | ng    | 100      |
| 29) 2,4-Dichlorophenol        | 10.581 | 162  | 182693   | 20.169 | ng    | 98       |
| 30) 1,2,4-Trichlorobenzene    | 10.799 | 180  | 208460   | 20.631 | ng    | 98       |
| 31) Naphthalene               | 10.993 | 128  | 632367   | 20.626 | ng    | 99       |
| 32) Benzoic acid              | 10.204 | 122  | 92362    | 17.829 | ng    | 93       |
| 33) 4-Chloroaniline           | 11.104 | 127  | 254334   | 20.065 | ng    | 98       |
| 34) Hexachlorobutadiene       | 11.269 | 225  | 132145   | 20.642 | ng    | 100      |
| 35) Caprolactam               | 11.875 | 113  | 50261    | 18.073 | ng    | 96       |
| 36) 4-Chloro-3-methylphenol   | 12.210 | 107  | 197097   | 20.013 | ng    | 98       |
| 37) 2-Methylnaphthalene       | 12.587 | 142  | 447608   | 20.542 | ng    | 98       |
| 38) 1-Methylnaphthalene       | 12.804 | 142  | 414942   | 20.409 | ng    | 100      |
| 40) 1,2,4,5-Tetrachloroben... | 12.951 | 216  | 227119   | 20.683 | ng    | 99       |
| 41) Hexachlorocyclopentadiene | 12.928 | 237  | 117842   | 20.261 | ng    | 98       |
| 43) 2,4,6-Trichlorophenol     | 13.193 | 196  | 144739   | 20.307 | ng    | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP052423\  
 Data File : BP015276.D  
 Acq On : 24 May 2023 17:12  
 Operator : CG/JU  
 Sample : SSTDICC020  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTDICC020

Manual Integrations  
 APPROVED

Reviewed By : Christian Giraldo 05/25/2023  
 Supervised By : Jagrut Upadhyay 05/25/2023

Quant Time: May 25 01:28:32 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP052423.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu May 25 01:19:28 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 13.263 | 196  | 168670   | 20.130 | ng    | 95       |
| 46) 1,1'-Biphenyl             | 13.587 | 154  | 556824   | 21.000 | ng    | 100      |
| 47) 2-Chloronaphthalene       | 13.634 | 162  | 417701   | 20.883 | ng    | 100      |
| 48) 2-Nitroaniline            | 13.840 | 65   | 122975   | 20.126 | ng    | 98       |
| 49) Acenaphthylene            | 14.475 | 152  | 668584   | 20.890 | ng    | 100      |
| 50) Dimethylphthalate         | 14.204 | 163  | 546936   | 20.890 | ng    | 99       |
| 51) 2,6-Dinitrotoluene        | 14.328 | 165  | 112500   | 20.352 | ng    | 97       |
| 52) Acenaphthene              | 14.816 | 154  | 395236   | 20.734 | ng    | 99       |
| 53) 3-Nitroaniline            | 14.657 | 138  | 114279   | 20.228 | ng    | 97       |
| 54) 2,4-Dinitrophenol         | 14.869 | 184  | 56706    | 19.443 | ng    | 96       |
| 55) Dibenzofuran              | 15.151 | 168  | 653636   | 20.961 | ng    | 99       |
| 56) 4-Nitrophenol             | 14.957 | 139  | 83992    | 19.625 | ng    | 97       |
| 57) 2,4-Dinitrotoluene        | 15.116 | 165  | 145447   | 20.012 | ng    | 97       |
| 58) Fluorene                  | 15.798 | 166  | 504661   | 20.954 | ng    | 99       |
| 59) 2,3,4,6-Tetrachlorophenol | 15.375 | 232  | 136263   | 20.333 | ng    | 100      |
| 60) Diethylphthalate          | 15.557 | 149  | 547009   | 20.973 | ng    | 99       |
| 61) 4-Chlorophenyl-phenyle... | 15.786 | 204  | 258763   | 20.983 | ng    | 100      |
| 62) 4-Nitroaniline            | 15.816 | 138  | 110766   | 19.642 | ng    | 98       |
| 63) Azobenzene                | 16.081 | 77   | 527421   | 21.198 | ng    | 99       |
| 65) 4,6-Dinitro-2-methylph... | 15.881 | 198  | 86777    | 20.196 | ng    | 99       |
| 66) n-Nitrosodiphenylamine    | 16.004 | 169  | 444631   | 21.093 | ng    | 100      |
| 67) 4-Bromophenyl-phenylether | 16.686 | 248  | 156891   | 20.511 | ng    | 98       |
| 68) Hexachlorobenzene         | 16.804 | 284  | 174522   | 20.718 | ng    | 97       |
| 69) Atrazine                  | 16.951 | 200  | 141134   | 20.969 | ng    | 98       |
| 70) Pentachlorophenol         | 17.151 | 266  | 115287   | 19.971 | ng    | 97       |
| 71) Phenanthrene              | 17.545 | 178  | 794293   | 21.067 | ng    | 99       |
| 72) Anthracene                | 17.639 | 178  | 797170   | 21.112 | ng    | 99       |
| 73) Carbazole                 | 17.904 | 167  | 706462   | 21.129 | ng    | 100      |
| 74) Di-n-butylphthalate       | 18.439 | 149  | 922373   | 21.481 | ng    | 100      |
| 75) Fluoranthene              | 19.498 | 202  | 938789   | 21.314 | ng    | 100      |
| 77) Benzidine                 | 19.680 | 184  | 224643   | 21.649 | ng    | 99       |
| 78) Pyrene                    | 19.851 | 202  | 969381   | 20.763 | ng    | 99       |
| 80) Butylbenzylphthalate      | 20.710 | 149  | 409492   | 20.697 | ng    | 98       |
| 81) Benzo(a)anthracene        | 21.569 | 228  | 950083   | 20.956 | ng    | 100      |
| 82) 3,3'-Dichlorobenzidine    | 21.492 | 252  | 294118   | 21.004 | ng    | 100      |
| 83) Chrysene                  | 21.621 | 228  | 904634   | 20.809 | ng    | 100      |
| 84) Bis(2-ethylhexyl)phtha... | 21.469 | 149  | 612487   | 21.019 | ng    | 100      |
| 85) Di-n-octyl phthalate      | 22.439 | 149  | 1018154  | 20.654 | ng    | 100      |
| 87) Indeno(1,2,3-cd)pyrene    | 26.845 | 276  | 1138565  | 20.984 | ng    | 99       |
| 88) Benzo(b)fluoranthene      | 23.351 | 252  | 974338   | 21.408 | ng    | 97       |
| 89) Benzo(k)fluoranthene      | 23.404 | 252  | 926567   | 20.217 | ng    | 99       |
| 90) Benzo(a)pyrene            | 24.021 | 252  | 919037   | 20.915 | ng    | 100      |
| 91) Dibenzo(a,h)anthracene    | 26.862 | 278  | 937969   | 21.051 | ng    | 100      |
| 92) Benzo(g,h,i)perylene      | 27.674 | 276  | 946916   | 20.915 | ng    | 99       |

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

## Manual Integrations

Quant Time: May 25 01:28:32 2023  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP052423.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Thu May 25 01:19:28 2023  
Response via : Initial Calibration

