

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP110920\  
 Data File : BP003866.D  
 Acq On : 09 Nov 2020 13:55  
 Operator : CG/JU  
 Sample : PB132858BS  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 PB132858BS

Manual Integrations  
 APPROVED

mohammad  
 11/10/2020 12:00:22 PM

Quant Time: Nov 09 14:25:57 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\8270-BP110320.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Nov 06 11:33:14 2020  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.304  | 152  | 165468   | 20.00  | ng    | 0.00     |
| 21) Naphthalene-d8            | 11.134 | 136  | 636704   | 20.00  | ng    | # 0.00   |
| 39) Acenaphthene-d10          | 14.922 | 164  | 379210   | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10          | 17.645 | 188  | 775775   | 20.00  | ng    | # 0.00   |
| 76) Chrysene-d12              | 21.674 | 240  | 723007   | 20.00  | ng    | 0.00     |
| 86) Perylene-d12              | 24.151 | 264  | 706112   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 5) 2-Fluorophenol             | 5.822  | 112  | 1334699  | 134.62 | ng    | 0.00     |
| 7) Phenol-d6                  | 7.469  | 99   | 1687900  | 118.60 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 9.469  | 82   | 1102578m | 84.81  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 16.398 | 330  | 626558   | 132.10 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 13.545 | 172  | 2295719  | 84.64  | ng    | 0.00     |
| 79) Terphenyl-d14             | 20.139 | 244  | 3008439  | 80.39  | ng    | 0.00     |
| Target Compounds              |        |      |          |        |       |          |
|                               |        |      |          |        |       | Qvalue   |
| 2) 1,4-Dioxane                | 3.511  | 88   | 138591   | 32.40  | ng    | # 63     |
| 3) Pyridine                   | 3.946  | 79   | 424242   | 33.82  | ng    | # 62     |
| 4) n-Nitrosodimethylamine     | 3.846  | 42   | 242447   | 41.98  | ng    | # 54     |
| 6) Aniline                    | 7.604  | 93   | 527216   | 32.62  | ng    | # 85     |
| 8) 2-Chlorophenol             | 7.869  | 128  | 494185   | 46.95  | ng    | # 81     |
| 9) Benzaldehyde               | 7.404  | 77   | 135583   | 17.92  | ng    | # 81     |
| 10) Phenol                    | 7.499  | 94   | 614734   | 44.76  | ng    | 83       |
| 11) bis(2-Chloroethyl)ether   | 7.693  | 93   | 489141   | 44.31  | ng    | # 83     |
| 12) 1,3-Dichlorobenzene       | 8.199  | 146  | 549285   | 42.58  | ng    | # 95     |
| 13) 1,4-Dichlorobenzene       | 8.340  | 146  | 560532   | 43.10  | ng    | 96       |
| 14) 1,2-Dichlorobenzene       | 8.669  | 146  | 533831   | 42.97  | ng    | 97       |
| 15) Benzyl Alcohol            | 8.534  | 79   | 452842   | 45.94  | ng    | # 87     |
| 16) 2,2'-oxybis(1-Chloropr... | 8.828  | 45   | 731009   | 39.32  | ng    | 81       |
| 17) 2-Methylphenol            | 8.763  | 107  | 408036   | 45.25  | ng    | # 94     |
| 18) Hexachloroethane          | 9.416  | 117  | 204585   | 44.26  | ng    | 97       |
| 19) n-Nitroso-di-n-propyla... | 9.104  | 70   | 372047   | 43.74  | ng    | # 77     |
| 20) 3+4-Methylphenols         | 9.087  | 107  | 547787   | 45.51  | ng    | 95       |
| 22) Acetophenone              | 9.122  | 105  | 729109   | 42.66  | ng    | # 86     |
| 24) Nitrobenzene              | 9.510  | 77   | 556007   | 43.04  | ng    | # 82     |
| 25) Isophorone                | 10.034 | 82   | 996481   | 45.13  | ng    | # 91     |
| 26) 2-Nitrophenol             | 10.234 | 139  | 260419   | 51.62  | ng    | # 73     |
| 27) 2,4-Dimethylphenol        | 10.298 | 122  | 439027   | 52.17  | ng    | 92       |
| 28) bis(2-Chloroethoxy)met... | 10.504 | 93   | 609133   | 40.20  | ng    | 98       |
| 29) 2,4-Dichlorophenol        | 10.787 | 162  | 459507   | 45.32  | ng    | 97       |
| 30) 1,2,4-Trichlorobenzene    | 11.004 | 180  | 521721   | 42.23  | ng    | 99       |
| 31) Naphthalene               | 11.187 | 128  | 1444834  | 42.29  | ng    | 100      |
| 32) Benzoic acid              | 10.457 | 122  | 195108   | 35.62  | ng    | # 81     |
| 33) 4-Chloroaniline           | 11.275 | 127  | 363712   | 25.05  | ng    | # 90     |
| 34) Hexachlorobutadiene       | 11.498 | 225  | 336098   | 41.71  | ng    | 100      |
| 35) Caprolactam               | 12.010 | 113  | 138410   | 44.15  | ng    | # 22     |
| 36) 4-Chloro-3-methylphenol   | 12.398 | 107  | 476935   | 43.57  | ng    | 91       |
| 37) 2-Methylnaphthalene       | 12.769 | 142  | 1041786  | 42.61  | ng    | 97       |
| 38) 1-Methylnaphthalene       | 12.987 | 142  | 960839   | 41.20  | ng    | 97       |
| 40) 1,2,4,5-Tetrachloroben... | 13.145 | 216  | 569602   | 45.39  | ng    | 98       |
| 41) Hexachlorocyclopentadiene | 13.139 | 237  | 726210   | 114.70 | ng    | 99       |
| 43) 2,4,6-Trichlorophenol     | 13.369 | 196  | 353918   | 45.62  | ng    | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP110920\  
 Data File : BP003866.D  
 Acq On : 09 Nov 2020 13:55  
 Operator : CG/JU  
 Sample : PB132858BS  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 PB132858BS

Manual Integrations  
 APPROVED

mohammad  
 11/10/2020 12:00:22 PM

Quant Time: Nov 09 14:25:57 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\8270-BP110320.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Nov 06 11:33:14 2020  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc  | Units | Dev(Min) |
|-------------------------------|--------|------|----------|-------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 13.451 | 196  | 373289   | 45.70 | ng    | 96       |
| 46) 1,1'-Biphenyl             | 13.751 | 154  | 1286331  | 43.70 | ng    | 98       |
| 47) 2-Chloronaphthalene       | 13.804 | 162  | 1013248  | 41.95 | ng    | 99       |
| 48) 2-Nitroaniline            | 13.981 | 65   | 313085   | 43.44 | ng    | # 60     |
| 49) Acenaphthylene            | 14.645 | 152  | 1543271  | 43.51 | ng    | 99       |
| 50) Dimethylphthalate         | 14.351 | 163  | 1230727  | 41.52 | ng    | 99       |
| 51) 2,6-Dinitrotoluene        | 14.463 | 165  | 275876   | 45.36 | ng    | # 73     |
| 52) Acenaphthene              | 14.986 | 154  | 1094535  | 46.33 | ng    | 93       |
| 53) 3-Nitroaniline            | 14.792 | 138  | 194701   | 28.92 | ng    | # 75     |
| 54) 2,4-Dinitrophenol         | 15.004 | 184  | 257348   | 76.80 | ng    | # 60     |
| 55) Dibenzofuran              | 15.310 | 168  | 1492834  | 42.53 | ng    | 99       |
| 56) 4-Nitrophenol             | 15.116 | 139  | 433442   | 89.33 | ng    | # 62     |
| 57) 2,4-Dinitrotoluene        | 15.251 | 165  | 371330   | 45.47 | ng    | # 75     |
| 58) Fluorene                  | 15.957 | 166  | 1191681  | 42.40 | ng    | 98       |
| 59) 2,3,4,6-Tetrachlorophenol | 15.545 | 232  | 323613   | 43.95 | ng    | 97       |
| 60) Diethylphthalate          | 15.704 | 149  | 1210568  | 41.71 | ng    | 98       |
| 61) 4-Chlorophenyl-phenyle... | 15.933 | 204  | 643015   | 41.86 | ng    | 98       |
| 62) 4-Nitroaniline            | 15.951 | 138  | 271003   | 38.61 | ng    | 86       |
| 63) Azobenzene                | 16.228 | 77   | 1139722  | 41.59 | ng    | 86       |
| 65) 4,6-Dinitro-2-methylph... | 16.028 | 198  | 201475   | 48.20 | ng    | 86       |
| 66) n-Nitrosodiphenylamine    | 16.145 | 169  | 1050563  | 44.24 | ng    | 97       |
| 67) 4-Bromophenyl-phenylether | 16.828 | 248  | 396410   | 43.67 | ng    | 98       |
| 68) Hexachlorobenzene         | 16.986 | 284  | 437591   | 42.25 | ng    | 99       |
| 69) Atrazine                  | 17.075 | 200  | 320777   | 41.14 | ng    | 99       |
| 70) Pentachlorophenol         | 17.316 | 266  | 454873   | 91.81 | ng    | 99       |
| 71) Phenanthrene              | 17.686 | 178  | 1836826  | 42.19 | ng    | 99       |
| 72) Anthracene                | 17.780 | 178  | 1874187  | 44.57 | ng    | 98       |
| 73) Carbazole                 | 18.027 | 167  | 1670584  | 42.82 | ng    | 99       |
| 74) Di-n-butylphthalate       | 18.545 | 149  | 2019451  | 44.72 | ng    | 99       |
| 75) Fluoranthene              | 19.616 | 202  | 2150316  | 42.80 | ng    | 99       |
| 77) Benzidine                 | 19.763 | 184  | 698566   | 40.96 | ng    | 99       |
| 78) Pyrene                    | 19.969 | 202  | 2186465  | 44.56 | ng    | 98       |
| 80) Butylbenzylphthalate      | 20.786 | 149  | 893019   | 48.85 | ng    | # 87     |
| 81) Benzo(a)anthracene        | 21.657 | 228  | 2069878  | 43.41 | ng    | 99       |
| 82) 3,3'-Dichlorobenzidine    | 21.568 | 252  | 567539   | 34.51 | ng    | 100      |
| 83) Chrysene                  | 21.715 | 228  | 1979385  | 43.21 | ng    | 98       |
| 84) Bis(2-ethylhexyl)phtha... | 21.533 | 149  | 1307063  | 48.50 | ng    | 100      |
| 85) Di-n-octyl phthalate      | 22.468 | 149  | 2228747  | 49.64 | ng    | # 87     |
| 87) Indeno(1,2,3-cd)pyrene    | 26.715 | 276  | 2219829  | 43.58 | ng    | 97       |
| 88) Benzo(b)fluoranthene      | 23.398 | 252  | 2064004  | 44.67 | ng    | 99       |
| 89) Benzo(k)fluoranthene      | 23.445 | 252  | 1996763  | 43.76 | ng    | 99       |
| 90) Benzo(a)pyrene            | 24.045 | 252  | 1836884  | 42.94 | ng    | 99       |
| 91) Dibenzo(a,h)anthracene    | 26.715 | 278  | 1900978  | 44.77 | ng    | 98       |
| 92) Benzo(g,h,i)perylene      | 27.503 | 276  | 1857205  | 45.19 | ng    | 97       |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP110920\  
 Data File : BP003866.D  
 Acq On : 09 Nov 2020 13:55  
 Operator : CG/JU  
 Sample : PB132858BS  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampled :  
 PB132858BS

Manual Integrations  
 APPROVED

mohammad  
 11/10/2020 12:00:22 PM

Quant Time: Nov 09 14:25:57 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\8270-BP110320.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Nov 06 11:33:14 2020  
 Response via : Initial Calibration

