

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP043025\  
 Data File : BP024479.D  
 Acq On : 30 Apr 2025 18:50  
 Operator : RC/JU  
 Sample : Q1900-08  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 WC-2

Quant Time: May 01 01:06:37 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP041425.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Apr 18 12:04:48 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-------------------------------|--------|------|----------|---------|-------|----------|
| Internal Standards            |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.711  | 152  | 135426   | 20.000  | ng    | -0.01    |
| 21) Naphthalene-d8            | 10.487 | 136  | 515132   | 20.000  | ng    | #-0.01   |
| 39) Acenaphthene-d10          | 14.340 | 164  | 312586   | 20.000  | ng    | -0.01    |
| 64) Phenanthrene-d10          | 17.139 | 188  | 637680   | 20.000  | ng    | # 0.00   |
| 76) Chrysene-d12              | 21.598 | 240  | 798028   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12              | 24.945 | 264  | 927368   | 20.000  | ng    | 0.01     |
| System Monitoring Compounds   |        |      |          |         |       |          |
| 5) 2-Fluorophenol             | 5.334  | 112  | 1103106  | 134.681 | ng    | -0.01    |
| 7) Phenol-d6                  | 6.899  | 99   | 1275077  | 113.693 | ng    | -0.01    |
| 23) Nitrobenzene-d5           | 8.858  | 82   | 845070   | 93.543  | ng    | -0.01    |
| 42) 2,4,6-Tribromophenol      | 15.857 | 330  | 710051   | 164.181 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 12.951 | 172  | 1794317  | 87.889  | ng    | -0.01    |
| 79) Terphenyl-d14             | 19.875 | 244  | 3556078  | 89.818  | ng    | 0.00     |
| Target Compounds              |        |      |          |         |       |          |
| 2) 1,4-Dioxane                | 3.287  | 88   | 306      | 0.088   | ng    | # 1      |
| 3) Pyridine                   | 3.670  | 79   | 189      | 0.021   | ng    | # 1      |
| 4) n-Nitrosodimethylamine     | 3.587  | 42   | 483      | 0.159   | ng    | # 1      |
| 6) Aniline                    | 7.058  | 93   | 49       | 0.005   | ng    | # 1      |
| 8) 2-Chlorophenol             | 7.163  | 128  | 228      | 0.025   | ng    | 79       |
| 9) Benzaldehyde               | 6.899  | 77   | 2424     | 0.437   | ng    | # 3      |
| 10) Phenol                    | 6.740  | 94   | 35       | 0.003   | ng    | 81       |
| 11) bis(2-Chloroethyl)ether   | 7.146  | 93   | 58       | 0.006   | ng    | # 47     |
| 12) 1,3-Dichlorobenzene       | 7.616  | 146  | 7        | 0.001   | ng    | # 1      |
| 13) 1,4-Dichlorobenzene       | 7.728  | 146  | 25       | 0.003   | ng    | # 1      |
| 14) 1,2-Dichlorobenzene       | 8.040  | 146  | 57       | 0.006   | ng    | # 58     |
| 15) Benzyl Alcohol            | 7.940  | 79   | 255      | 0.035   | ng    | # 43     |
| 16) 2,2'-oxybis(1-Chloropr... | 8.252  | 45   | 219      | 0.022   | ng    | 86       |
| 17) 2-Methylphenol            | 8.175  | 107  | 58       | 0.008   | ng    | # 1      |
| 18) Hexachloroethane          | 8.810  | 117  | 81       | 0.022   | ng    | # 42     |
| 19) n-Nitroso-di-n-propyla... | 8.505  | 70   | 50       | 0.007   | ng    | # 1      |
| 20) 3+4-Methylphenols         | 8.481  | 107  | 35       | 0.003   | ng    | # 1      |
| 22) Acetophenone              | 8.528  | 105  | 1034     | 0.081   | ng    | # 40     |
| 24) Nitrobenzene              | 8.928  | 77   | 123      | 0.014   | ng    | # 66     |
| 25) Isophorone                | 9.410  | 82   | 88       | 0.005   | ng    | # 63     |
| 26) 2-Nitrophenol             | 9.610  | 139  | 17       | 0.004   | ng    | # 1      |
| 28) bis(2-Chloroethoxy)met... | 9.922  | 93   | 149      | 0.013   | ng    | # 1      |
| 29) 2,4-Dichlorophenol        | 10.128 | 162  | 24       | 0.003   | ng    | # 61     |
| 30) 1,2,4-Trichlorobenzene    | 10.469 | 180  | 15       | 0.002   | ng    | 94       |
| 31) Naphthalene               | 10.534 | 128  | 2088     | 0.077   | ng    | # 84     |
| 32) Benzoic acid              | 9.781  | 122  | 52679    | 8.233   | ng    | 99       |
| 33) 4-Chloroaniline           | 10.663 | 127  | 20       | 0.002   | ng    | # 1      |
| 34) Hexachlorobutadiene       | 10.816 | 225  | 12       | 0.003   | ng    | # 1      |
| 35) Caprolactam               | 11.428 | 113  | 70       | 0.025   | ng    | # 16     |
| 36) 4-Chloro-3-methylphenol   | 11.804 | 107  | 96       | 0.011   | ng    | 86       |
| 37) 2-Methylnaphthalene       | 12.157 | 142  | 1288     | 0.068   | ng    | # 70     |
| 38) 1-Methylnaphthalene       | 12.369 | 142  | 954      | 0.052   | ng    | # 87     |
| 40) 1,2,4,5-Tetrachloroben... | 12.516 | 216  | 28       | 0.003   | ng    | # 50     |
| 41) Hexachlorocyclopentadiene | 12.557 | 237  | 13       | 0.004   | ng    | 85       |
| 43) 2,4,6-Trichlorophenol     | 12.775 | 196  | 16       | 0.003   | ng    | # 67     |
| 44) 2,4,5-Trichlorophenol     | 12.863 | 196  | 8        | 0.001   | ng    | # 1      |

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Quant Time: May 01 01:06:37 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP041425.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Apr 18 12:04:48 2025  
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| Compound                      | R.T.   | QIon | Response | Conc  | Units | Dev(Min) |
|-------------------------------|--------|------|----------|-------|-------|----------|
| 46) 1,1'-Biphenyl             | 13.169 | 154  | 212      | 0.009 | ng    | # 60     |
| 47) 2-Chloronaphthalene       | 13.475 | 162  | 22       | 0.001 | ng    | 96       |
| 48) 2-Nitroaniline            | 13.428 | 65   | 32       | 0.007 | ng    | # 17     |
| 49) Acenaphthylene            | 14.063 | 152  | 353      | 0.013 | ng    | # 83     |
| 50) Dimethylphthalate         | 13.798 | 163  | 109743   | 4.828 | ng    | 98       |
| 51) 2,6-Dinitrotoluene        | 13.945 | 165  | 156      | 0.032 | ng    | # 60     |
| 52) Acenaphthene              | 14.410 | 154  | 1113     | 0.065 | ng    | 97       |
| 53) 3-Nitroaniline            | 14.245 | 138  | 41       | 0.008 | ng    | # 1      |
| 54) 2,4-Dinitrophenol         | 14.434 | 184  | 28       | 0.010 | ng    | # 1      |
| 55) Dibenzofuran              | 14.751 | 168  | 707      | 0.025 | ng    | # 73     |
| 56) 4-Nitrophenol             | 14.581 | 139  | 78       | 0.018 | ng    | # 10     |
| 57) 2,4-Dinitrotoluene        | 14.710 | 165  | 107      | 0.016 | ng    | # 58     |
| 58) Fluorene                  | 15.416 | 166  | 868      | 0.040 | ng    | # 73     |
| 59) 2,3,4,6-Tetrachlorophenol | 14.998 | 232  | 33       | 0.006 | ng    | # 1      |
| 60) Diethylphthalate          | 15.181 | 149  | 3385     | 0.144 | ng    | 97       |
| 61) 4-Chlorophenyl-phenyle... | 15.398 | 204  | 14       | 0.001 | ng    | # 1      |
| 62) 4-Nitroaniline            | 15.369 | 138  | 82       | 0.015 | ng    | 95       |
| 63) Azobenzene                | 15.687 | 77   | 152      | 0.007 | ng    | # 70     |
| 65) 4,6-Dinitro-2-methylph... | 15.516 | 198  | 37       | 0.009 | ng    | # 1      |
| 66) n-Nitrosodiphenylamine    | 15.857 | 169  | 20023    | 1.052 | ng    | # 51     |
| 67) 4-Bromophenyl-phenylether | 16.322 | 248  | 15       | 0.002 | ng    | # 1      |
| 68) Hexachlorobenzene         | 16.439 | 284  | 18       | 0.002 | ng    | # 1      |
| 69) Atrazine                  | 16.628 | 200  | 24       | 0.005 | ng    | 89       |
| 70) Pentachlorophenol         | 16.787 | 266  | 8        | 0.001 | ng    | # 1      |
| 71) Phenanthrene              | 17.181 | 178  | 3101     | 0.089 | ng    | # 91     |
| 72) Anthracene                | 17.286 | 178  | 636      | 0.019 | ng    | # 20     |
| 73) Carbazole                 | 17.586 | 167  | 585      | 0.018 | ng    | # 1      |
| 74) Di-n-butylphthalate       | 18.151 | 149  | 15226    | 0.373 | ng    | 98       |
| 75) Fluoranthene              | 19.280 | 202  | 1488     | 0.036 | ng    | 86       |
| 77) Benzidine                 | 19.510 | 184  | 24       | 0.002 | ng    | # 1      |
| 78) Pyrene                    | 19.657 | 202  | 2203     | 0.043 | ng    | 95       |
| 80) Butylbenzylphthalate      | 20.604 | 149  | 1470     | 0.068 | ng    | # 78     |
| 81) Benzo(a)anthracene        | 21.604 | 228  | 3051     | 0.062 | ng    | # 72     |
| 82) 3,3'-Dichlorobenzidine    | 21.469 | 252  | 101      | 0.006 | ng    | # 37     |
| 83) Chrysene                  | 21.651 | 228  | 816      | 0.017 | ng    | # 84     |
| 84) Bis(2-ethylhexyl)phtha... | 21.510 | 149  | 9389     | 0.295 | ng    | 95       |
| 85) Di-n-octyl phthalate      | 22.798 | 149  | 608      | 0.012 | ng    | # 87     |
| 87) Indeno(1,2,3-cd)pyrene    | 28.768 | 276  | 62       | 0.001 | ng    | # 63     |
| 88) Benzo(b)fluoranthene      | 23.904 | 252  | 889      | 0.016 | ng    | # 1      |
| 89) Benzo(k)fluoranthene      | 23.957 | 252  | 191      | 0.004 | ng    | # 1      |
| 90) Benzo(a)pyrene            | 24.786 | 252  | 316      | 0.007 | ng    | # 1      |
| 91) Dibenzo(a,h)anthracene    | 28.821 | 278  | 55       | 0.001 | ng    | # 1      |
| 92) Benzo(g,h,i)perylene      | 30.162 | 276  | 195      | 0.004 | ng    | 92       |

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

BNA P

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WC-2

Quant Time: May 01 01:06:37 2025

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP041425.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Apr 18 12:04:48 2025

Response via : Initial Calibration

