

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080525\  
 Data File : BP025302.D  
 Acq On : 05 Aug 2025 14:47  
 Operator : CG/JU  
 Sample : SSTDICCC040  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTDICCC040

Manual Integrations  
 APPROVED

Reviewed By :Rahul Chavli 08/06/2025  
 Supervised By :Jagrut Upadhyay 08/06/2025

Quant Time: Aug 06 06:01:49 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP080525.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Aug 06 05:56:10 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.802  | 152  | 201963   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8            | 10.572 | 136  | 880642   | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10          | 14.419 | 164  | 554286   | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10          | 17.225 | 188  | 1096367  | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12              | 21.660 | 240  | 1098659  | 20.000 | ng    | 0.00     |
| 86) Perylene-d12              | 25.036 | 264  | 1264736  | 20.000 | ng    | 0.00     |
|                               |        |      |          |        |       |          |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 5) 2-Fluorophenol             | 5.408  | 112  | 996403   | 78.838 | ng    | 0.00     |
| 7) Phenol-d6                  | 6.966  | 99   | 1318375  | 79.077 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 8.943  | 82   | 1316251  | 80.869 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 15.925 | 330  | 470979   | 82.296 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 13.031 | 172  | 3200514  | 80.179 | ng    | 0.00     |
| 79) Terphenyl-d14             | 19.942 | 244  | 4595144  | 79.015 | ng    | 0.00     |
|                               |        |      |          |        |       |          |
| Target Compounds              |        |      |          |        |       | Qvalue   |
| 2) 1,4-Dioxane                | 3.337  | 88   | 194354   | 39.221 | ng    | 100      |
| 3) Pyridine                   | 3.725  | 79   | 541403   | 39.325 | ng    | 100      |
| 4) n-Nitrosodimethylamine     | 3.631  | 42   | 246918   | 38.631 | ng    | 100      |
| 6) Aniline                    | 7.131  | 93   | 895891   | 39.490 | ng    | 100      |
| 8) 2-Chlorophenol             | 7.372  | 128  | 545256   | 39.670 | ng    | 100      |
| 9) Benzaldehyde               | 6.949  | 77   | 566415   | 48.256 | ng    | 100      |
| 10) Phenol                    | 6.990  | 94   | 699120   | 39.538 | ng    | 100      |
| 11) bis(2-Chloroethyl)ether   | 7.225  | 93   | 539982   | 38.975 | ng    | 100      |
| 12) 1,3-Dichlorobenzene       | 7.696  | 146  | 595110   | 38.533 | ng    | 100      |
| 13) 1,4-Dichlorobenzene       | 7.837  | 146  | 609114   | 38.983 | ng    | 100      |
| 14) 1,2-Dichlorobenzene       | 8.149  | 146  | 587683   | 38.826 | ng    | 100      |
| 15) Benzyl Alcohol            | 8.031  | 79   | 502383   | 39.218 | ng    | 100      |
| 16) 2,2'-oxybis(1-Chloropr... | 8.325  | 45   | 850917   | 38.879 | ng    | 100      |
| 17) 2-Methylphenol            | 8.225  | 107  | 474988   | 40.146 | ng    | 100      |
| 18) Hexachloroethane          | 8.878  | 117  | 218135   | 39.037 | ng    | 100      |
| 19) n-Nitroso-di-n-propyla... | 8.596  | 70   | 436056   | 40.160 | ng    | 100      |
| 20) 3+4-Methylphenols         | 8.549  | 107  | 640969   | 39.342 | ng    | 100      |
| 22) Acetophenone              | 8.607  | 105  | 867835   | 39.255 | ng    | 100      |
| 24) Nitrobenzene              | 8.984  | 77   | 606097   | 40.280 | ng    | 100      |
| 25) Isophorone                | 9.502  | 82   | 1255611  | 39.682 | ng    | 100      |
| 26) 2-Nitrophenol             | 9.684  | 139  | 225091   | 35.376 | ng    | 100      |
| 27) 2,4-Dimethylphenol        | 9.749  | 122  | 570930   | 39.137 | ng    | 100      |
| 28) bis(2-Chloroethoxy)met... | 9.984  | 93   | 759841   | 39.476 | ng    | 100      |
| 29) 2,4-Dichlorophenol        | 10.213 | 162  | 522153   | 40.884 | ng    | 100      |
| 30) 1,2,4-Trichlorobenzene    | 10.437 | 180  | 548893   | 38.956 | ng    | 100      |
| 31) Naphthalene               | 10.625 | 128  | 1791087  | 39.178 | ng    | 100      |
| 32) Benzoic acid              | 9.849  | 122  | 297213   | 38.952 | ng    | 100      |
| 33) 4-Chloroaniline           | 10.725 | 127  | 808713   | 39.774 | ng    | 100      |
| 34) Hexachlorobutadiene       | 10.913 | 225  | 316426   | 39.122 | ng    | 100      |
| 35) Caprolactam               | 11.472 | 113  | 202605   | 40.802 | ng    | 100      |
| 36) 4-Chloro-3-methylphenol   | 11.831 | 107  | 600241   | 40.617 | ng    | 100      |
| 37) 2-Methylnaphthalene       | 12.225 | 142  | 1154835  | 39.028 | ng    | 100      |
| 38) 1-Methylnaphthalene       | 12.454 | 142  | 1205124  | 39.032 | ng    | 100      |
| 40) 1,2,4,5-Tetrachloroben... | 12.595 | 216  | 606083   | 40.051 | ng    | 100      |
| 41) Hexachlorocyclopentadiene | 12.590 | 237  | 366850   | 39.278 | ng    | 100      |
| 43) 2,4,6-Trichlorophenol     | 12.837 | 196  | 406436   | 41.576 | ng    | 100      |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080525\  
 Data File : BP025302.D  
 Acq On : 05 Aug 2025 14:47  
 Operator : CG/JU  
 Sample : SSTDICCC040  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTDICCC040

Manual Integrations  
 APPROVED

Reviewed By :Rahul Chavli 08/06/2025  
 Supervised By :Jagrut Upadhyay 08/06/2025

Quant Time: Aug 06 06:01:49 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP080525.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Aug 06 05:56:10 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 12.901 | 196  | 456928   | 41.667 | ng    | 100      |
| 46) 1,1'-Biphenyl             | 13.243 | 154  | 1605905  | 39.314 | ng    | 100      |
| 47) 2-Chloronaphthalene       | 13.284 | 162  | 1225849  | 39.335 | ng    | 100      |
| 48) 2-Nitroaniline            | 13.484 | 65   | 341536   | 38.066 | ng    | 100      |
| 49) Acenaphthylene            | 14.142 | 152  | 2070346  | 39.828 | ng    | 100      |
| 50) Dimethylphthalate         | 13.872 | 163  | 1586291  | 39.258 | ng    | 100      |
| 51) 2,6-Dinitrotoluene        | 13.978 | 165  | 318700   | 42.760 | ng    | 100      |
| 52) Acenaphthene              | 14.490 | 154  | 1234915  | 38.947 | ng    | 100      |
| 53) 3-Nitroaniline            | 14.307 | 138  | 369122   | 42.324 | ng    | 100      |
| 54) 2,4-Dinitrophenol         | 14.513 | 184  | 105656   | 38.143 | ng    | 100      |
| 55) Dibenzofuran              | 14.831 | 168  | 1863874  | 39.336 | ng    | 100      |
| 56) 4-Nitrophenol             | 14.607 | 139  | 314409   | 40.848 | ng    | 100      |
| 57) 2,4-Dinitrotoluene        | 14.778 | 165  | 431918   | 38.103 | ng    | 100      |
| 58) Fluorene                  | 15.489 | 166  | 1458854  | 39.200 | ng    | 100      |
| 59) 2,3,4,6-Tetrachlorophenol | 15.054 | 232  | 381224   | 41.372 | ng    | 100      |
| 60) Diethylphthalate          | 15.260 | 149  | 1596506  | 39.190 | ng    | 100      |
| 61) 4-Chlorophenyl-phenyle... | 15.484 | 204  | 699128   | 39.626 | ng    | 100      |
| 62) 4-Nitroaniline            | 15.489 | 138  | 374812   | 43.373 | ng    | 100      |
| 63) Azobenzene                | 15.784 | 77   | 1497305  | 39.565 | ng    | 100      |
| 65) 4,6-Dinitro-2-methylph... | 15.560 | 198  | 165146   | 37.407 | ng    | 100      |
| 66) n-Nitrosodiphenylamine    | 15.695 | 169  | 1327098  | 39.754 | ng    | 100      |
| 67) 4-Bromophenyl-phenylether | 16.395 | 248  | 436238   | 39.545 | ng    | 100      |
| 68) Hexachlorobenzene         | 16.513 | 284  | 482334   | 38.967 | ng    | 100      |
| 69) Atrazine                  | 16.666 | 200  | 465752   | 39.673 | ng    | 100      |
| 70) Pentachlorophenol         | 16.860 | 266  | 315442   | 40.000 | ng    | 100      |
| 71) Phenanthrene              | 17.266 | 178  | 2351974  | 39.167 | ng    | 100      |
| 72) Anthracene                | 17.354 | 178  | 2401769  | 39.667 | ng    | 100      |
| 73) Carbazole                 | 17.631 | 167  | 2307398  | 39.700 | ng    | 100      |
| 74) Di-n-butylphthalate       | 18.236 | 149  | 2828001  | 40.220 | ng    | 100      |
| 75) Fluoranthene              | 19.342 | 202  | 2745884  | 39.501 | ng    | 100      |
| 77) Benzidine                 | 19.530 | 184  | 2131359  | 53.598 | ng    | 100      |
| 78) Pyrene                    | 19.719 | 202  | 2831811  | 38.761 | ng    | 100      |
| 80) Butylbenzylphthalate      | 20.672 | 149  | 1241891  | 41.913 | ng    | 100      |
| 81) Benzo(a)anthracene        | 21.642 | 228  | 2859970  | 39.173 | ng    | 100      |
| 82) 3,3'-Dichlorobenzidine    | 21.554 | 252  | 1039593  | 40.358 | ng    | 100      |
| 83) Chrysene                  | 21.701 | 228  | 2674978  | 39.503 | ng    | 100      |
| 84) Bis(2-ethylhexyl)phtha... | 21.607 | 149  | 1959652  | 41.420 | ng    | 100      |
| 85) Di-n-octyl phthalate      | 22.901 | 149  | 3300305  | 41.675 | ng    | 100      |
| 87) Indeno(1,2,3-cd)pyrene    | 28.901 | 276  | 3649652m | 40.200 | ng    |          |
| 88) Benzo(b)fluoranthene      | 23.971 | 252  | 2981039  | 39.375 | ng    | 100      |
| 89) Benzo(k)fluoranthene      | 24.048 | 252  | 2961242  | 39.456 | ng    | 100      |
| 90) Benzo(a)pyrene            | 24.883 | 252  | 2911834  | 39.938 | ng    | 100      |
| 91) Dibenzo(a,h)anthracene    | 28.971 | 278  | 2998158  | 40.383 | ng    | 100      |
| 92) Benzo(g,h,i)perylene      | 30.053 | 276  | 2917156  | 40.025 | ng    | 100      |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080525\  
Data File : BP025302.D  
Acq On : 05 Aug 2025 14:47  
Operator : CG/JU  
Sample : SSTDICCC040  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SSTDICCC040

Quant Time: Aug 06 06:01:49 2025  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP080525.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Wed Aug 06 05:56:10 2025  
Response via : Initial Calibration

Manual Integrations  
APPROVED

Reviewed By :Rahul Chavli 08/06/2025  
Supervised By :Jagrut Upadhyay 08/06/2025

