

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM050825\  
 Data File : BM050143.D  
 Acq On : 08 May 2025 12:39  
 Operator : RC/JU  
 Sample : PB167889BSD  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 PB167889BSD

Manual Integrations  
 APPROVED

Reviewed By :Rahul Chavli 05/09/2025  
 Supervised By :Jagrut Upadhyay 05/09/2025

Quant Time: May 08 13:14:53 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM042825.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu May 08 12:03:21 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-------------------------------|--------|------|----------|---------|-------|----------|
| -----                         |        |      |          |         |       |          |
| Internal Standards            |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.740  | 152  | 264346   | 20.000  | ng    | 0.00     |
| 21) Naphthalene-d8            | 10.534 | 136  | 911076   | 20.000  | ng    | 0.00     |
| 39) Acenaphthene-d10          | 14.392 | 164  | 548910   | 20.000  | ng    | 0.00     |
| 64) Phenanthrene-d10          | 17.139 | 188  | 982803   | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12              | 21.386 | 240  | 868474   | 20.000  | ng    | 0.00     |
| 86) Perylene-d12              | 24.380 | 264  | 893841   | 20.000  | ng    | 0.00     |
| System Monitoring Compounds   |        |      |          |         |       |          |
| 5) 2-Fluorophenol             | 5.340  | 112  | 2021970  | 133.939 | ng    | 0.00     |
| 7) Phenol-d6                  | 6.934  | 99   | 2452721  | 129.267 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 8.904  | 82   | 1448557  | 78.347  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 15.886 | 330  | 1063477  | 131.029 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 13.004 | 172  | 3723241  | 80.242  | ng    | 0.00     |
| 79) Terphenyl-d14             | 19.768 | 244  | 5036934  | 85.821  | ng    | 0.00     |
| Target Compounds              |        |      |          |         |       |          |
|                               |        |      |          |         |       | Qvalue   |
| 2) 1,4-Dioxane                | 3.234  | 88   | 230236   | 35.570  | ng    | 99       |
| 3) Pyridine                   | 3.634  | 79   | 631457   | 37.969  | ng    | 97       |
| 4) n-Nitrosodimethylamine     | 3.546  | 42   | 272864   | 41.833  | ng    | 99       |
| 6) Aniline                    | 7.075  | 93   | 646417   | 26.980  | ng    | 98       |
| 8) 2-Chlorophenol             | 7.316  | 128  | 713271   | 43.699  | ng    | 99       |
| 9) Benzaldehyde               | 6.887  | 77   | 294210   | 23.171  | ng    | 99       |
| 10) Phenol                    | 6.957  | 94   | 838666   | 43.662  | ng    | 99       |
| 11) bis(2-Chloroethyl)ether   | 7.169  | 93   | 669133   | 41.719  | ng    | 99       |
| 12) 1,3-Dichlorobenzene       | 7.628  | 146  | 816752   | 41.424  | ng    | 100      |
| 13) 1,4-Dichlorobenzene       | 7.775  | 146  | 821437   | 41.470  | ng    | 99       |
| 14) 1,2-Dichlorobenzene       | 8.093  | 146  | 788459   | 41.208  | ng    | 99       |
| 15) Benzyl Alcohol            | 7.993  | 79   | 559284   | 42.504  | ng    | 96       |
| 16) 2,2'-oxybis(1-Chloropr... | 8.263  | 45   | 810918   | 41.593  | ng    | 99       |
| 17) 2-Methylphenol            | 8.204  | 107  | 540207   | 42.691  | ng    | 99       |
| 18) Hexachloroethane          | 8.816  | 117  | 293460   | 41.677  | ng    | 99       |
| 19) n-Nitroso-di-n-propyla... | 8.551  | 70   | 488851   | 40.172  | ng    | 100      |
| 20) 3+4-Methylphenols         | 8.534  | 107  | 717828   | 42.007  | ng    | 98       |
| 22) Acetophenone              | 8.569  | 105  | 969487   | 42.801  | ng    | # 98     |
| 24) Nitrobenzene              | 8.945  | 77   | 695574   | 43.509  | ng    | 98       |
| 25) Isophorone                | 9.475  | 82   | 1290030  | 40.932  | ng    | 99       |
| 26) 2-Nitrophenol             | 9.657  | 139  | 367311   | 44.974  | ng    | 98       |
| 27) 2,4-Dimethylphenol        | 9.728  | 122  | 595475   | 43.965  | ng    | 99       |
| 28) bis(2-Chloroethoxy)met... | 9.945  | 93   | 817207   | 41.952  | ng    | 99       |
| 29) 2,4-Dichlorophenol        | 10.210 | 162  | 674561   | 44.457  | ng    | 99       |
| 30) 1,2,4-Trichlorobenzene    | 10.398 | 180  | 792832   | 42.886  | ng    | 99       |
| 31) Naphthalene               | 10.587 | 128  | 1954746  | 42.366  | ng    | 100      |
| 32) Benzoic acid              | 9.910  | 122  | 394151m  | 45.501  | ng    |          |
| 33) 4-Chloroaniline           | 10.716 | 127  | 306640   | 16.214  | ng    | 98       |
| 34) Hexachlorobutadiene       | 10.869 | 225  | 513285   | 43.740  | ng    | 99       |
| 35) Caprolactam               | 11.510 | 113  | 167178   | 37.787  | ng    | 96       |
| 36) 4-Chloro-3-methylphenol   | 11.857 | 107  | 576231   | 41.105  | ng    | 99       |
| 37) 2-Methylnaphthalene       | 12.204 | 142  | 1271852  | 41.115  | ng    | 100      |
| 38) 1-Methylnaphthalene       | 12.422 | 142  | 1327396  | 40.548  | ng    | 99       |
| 40) 1,2,4,5-Tetrachloroben... | 12.581 | 216  | 870924   | 45.834  | ng    | 99       |
| 41) Hexachlorocyclopentadiene | 12.551 | 237  | 1147740  | 98.838  | ng    | 100      |
| 43) 2,4,6-Trichlorophenol     | 12.833 | 196  | 549841   | 45.622  | ng    | 97       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM050825\  
 Data File : BM050143.D  
 Acq On : 08 May 2025 12:39  
 Operator : RC/JU  
 Sample : PB167889BSD  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 PB167889BSD

Manual Integrations  
 APPROVED

Reviewed By :Rahul Chavli 05/09/2025  
 Supervised By :Jagrut Upadhyay 05/09/2025

Quant Time: May 08 13:14:53 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM042825.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu May 08 12:03:21 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 12.922 | 196  | 594889   | 45.773 | ng    | 98       |
| 46) 1,1'-Biphenyl             | 13.216 | 154  | 1813471  | 44.444 | ng    | 99       |
| 47) 2-Chloronaphthalene       | 13.263 | 162  | 1442235  | 44.629 | ng    | 100      |
| 48) 2-Nitroaniline            | 13.480 | 65   | 346667   | 45.707 | ng    | 99       |
| 49) Acenaphthylene            | 14.110 | 152  | 2217366  | 43.518 | ng    | 100      |
| 50) Dimethylphthalate         | 13.851 | 163  | 1687386  | 42.064 | ng    | 99       |
| 51) 2,6-Dinitrotoluene        | 13.975 | 165  | 364981   | 43.944 | ng    | 99       |
| 52) Acenaphthene              | 14.451 | 154  | 1280948  | 43.432 | ng    | 99       |
| 53) 3-Nitroaniline            | 14.316 | 138  | 187182   | 23.606 | ng    | 98       |
| 54) 2,4-Dinitrophenol         | 14.533 | 184  | 430670   | 80.951 | ng    | 98       |
| 55) Dibenzofuran              | 14.792 | 168  | 2092337  | 42.639 | ng    | 99       |
| 56) 4-Nitrophenol             | 14.663 | 139  | 485201   | 75.484 | ng    | 99       |
| 57) 2,4-Dinitrotoluene        | 14.775 | 165  | 497539   | 43.352 | ng    | 97       |
| 58) Fluorene                  | 15.439 | 166  | 1702697  | 42.392 | ng    | 100      |
| 59) 2,3,4,6-Tetrachlorophenol | 15.033 | 232  | 482259   | 42.868 | ng    | 99       |
| 60) Diethylphthalate          | 15.216 | 149  | 1537382  | 40.672 | ng    | 100      |
| 61) 4-Chlorophenyl-phenyle... | 15.433 | 204  | 942516   | 42.723 | ng    | 99       |
| 62) 4-Nitroaniline            | 15.480 | 138  | 285268   | 37.288 | ng    | 97       |
| 63) Azobenzene                | 15.727 | 77   | 1345417  | 42.471 | ng    | 99       |
| 65) 4,6-Dinitro-2-methylph... | 15.545 | 198  | 279844   | 45.488 | ng    | 97       |
| 66) n-Nitrosodiphenylamine    | 15.651 | 169  | 1392965  | 46.156 | ng    | 100      |
| 67) 4-Bromophenyl-phenylether | 16.327 | 248  | 542393   | 45.638 | ng    | 98       |
| 68) Hexachlorobenzene         | 16.451 | 284  | 625015   | 45.623 | ng    | 97       |
| 69) Atrazine                  | 16.610 | 200  | 478867   | 43.989 | ng    | 100      |
| 70) Pentachlorophenol         | 16.804 | 266  | 748706   | 97.091 | ng    | 99       |
| 71) Phenanthrene              | 17.180 | 178  | 2456477  | 44.316 | ng    | 100      |
| 72) Anthracene                | 17.274 | 178  | 2500364  | 44.573 | ng    | 100      |
| 73) Carbazole                 | 17.551 | 167  | 2142721  | 44.019 | ng    | 99       |
| 74) Di-n-butylphthalate       | 18.104 | 149  | 2500884  | 43.142 | ng    | 100      |
| 75) Fluoranthene              | 19.204 | 202  | 2767529  | 42.525 | ng    | 100      |
| 77) Benzidine                 | 19.392 | 184  | 926472   | 30.033 | ng    | 99       |
| 78) Pyrene                    | 19.568 | 202  | 2838617  | 46.545 | ng    | 100      |
| 80) Butylbenzylphthalate      | 20.462 | 149  | 1034547  | 45.291 | ng    | 97       |
| 81) Benzo(a)anthracene        | 21.368 | 228  | 2693046  | 45.030 | ng    | 100      |
| 82) 3,3'-Dichlorobenzidine    | 21.292 | 252  | 571243   | 24.952 | ng    | 99       |
| 83) Chrysene                  | 21.427 | 228  | 2451903  | 44.297 | ng    | 99       |
| 84) Bis(2-ethylhexyl)phtha... | 21.280 | 149  | 1507874  | 44.196 | ng    | 99       |
| 85) Di-n-octyl phthalate      | 22.403 | 149  | 2495544  | 44.353 | ng    | 100      |
| 87) Indeno(1,2,3-cd)pyrene    | 27.768 | 276  | 3344168  | 50.211 | ng    | 100      |
| 88) Benzo(b)fluoranthene      | 23.433 | 252  | 2565912  | 44.142 | ng    | 100      |
| 89) Benzo(k)fluoranthene      | 23.492 | 252  | 2534162  | 43.098 | ng    | 100      |
| 90) Benzo(a)pyrene            | 24.239 | 252  | 2428116  | 44.598 | ng    | 99       |
| 91) Dibenzo(a,h)anthracene    | 27.809 | 278  | 2734690  | 50.370 | ng    | 99       |
| 92) Benzo(g,h,i)perylene      | 28.821 | 276  | 2620801  | 50.422 | ng    | 99       |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM050825\  
Data File : BM050143.D  
Acq On : 08 May 2025 12:39  
Operator : RC/JU  
Sample : PB167889BSD  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

## Instrument :

BNA\_M

## ClientSampleId :

PB167889BSD

## Manual Integrations

## APPROVED

Reviewed By :Rahul Chavli 05/09/2025

Supervised By :Jagrut Upadhyay 05/09/2025

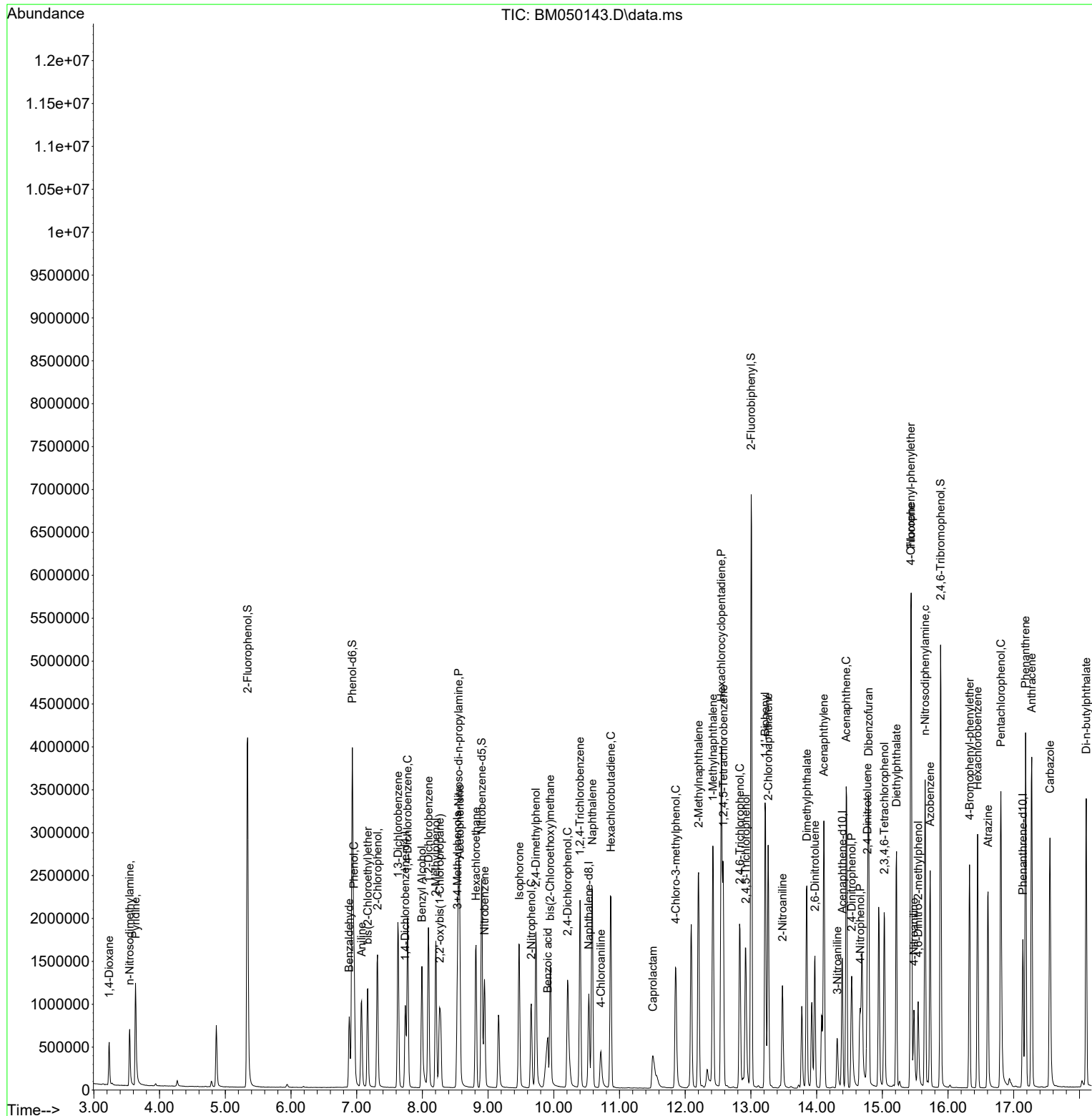
Quant Time: May 08 13:14:53 2025

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM042825.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Thu May 08 12:03:21 2025

Response via : Initial Calibration



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM050825\  
Data File : BM050143.D  
Acq On : 08 May 2025 12:39  
Operator : RC/JU  
Sample : PB167889BSD  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
PB167889BSD

Manual Integrations  
APPROVED

Reviewed By :Rahul Chavli 05/09/2025  
Supervised By :Jagrut Upadhyay 05/09/2025

Quant Time: May 08 13:14:53 2025  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM042825.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Thu May 08 12:03:21 2025  
Response via : Initial Calibration

