

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM042125\  
 Data File : BM049980.D  
 Acq On : 21 Apr 2025 11:15  
 Operator : RC/JU  
 Sample : PB167645BS  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 PB167645BS

Quant Time: Apr 21 11:59:36 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM040825.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Apr 09 04:00:55 2025  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units | Dev(Min) |
|-------------------------------|--------|------|----------|---------|-------|----------|
| -----                         |        |      |          |         |       |          |
| Internal Standards            |        |      |          |         |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.769  | 152  | 405465   | 20.000  | ng    | -0.01    |
| 21) Naphthalene-d8            | 10.563 | 136  | 1378196  | 20.000  | ng    | -0.01    |
| 39) Acenaphthene-d10          | 14.410 | 164  | 810401   | 20.000  | ng    | -0.01    |
| 64) Phenanthrene-d10          | 17.157 | 188  | 1404843  | 20.000  | ng    | 0.00     |
| 76) Chrysene-d12              | 21.398 | 240  | 1154135  | 20.000  | ng    | 0.00     |
| 86) Perylene-d12              | 24.397 | 264  | 1149601  | 20.000  | ng    | -0.01    |
| System Monitoring Compounds   |        |      |          |         |       |          |
| 5) 2-Fluorophenol             | 5.357  | 112  | 2892461  | 121.135 | ng    | 0.00     |
| 7) Phenol-d6                  | 6.957  | 99   | 3494435  | 117.624 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 8.928  | 82   | 2042556  | 82.528  | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 15.910 | 330  | 1446565  | 122.719 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 13.033 | 172  | 5148618  | 86.150  | ng    | -0.01    |
| 79) Terphenyl-d14             | 19.786 | 244  | 6229008  | 101.145 | ng    | 0.00     |
| Target Compounds              |        |      |          |         |       |          |
|                               |        |      |          |         |       | Qvalue   |
| 2) 1,4-Dioxane                | 3.252  | 88   | 306579   | 31.176  | ng    | 99       |
| 3) Pyridine                   | 3.652  | 79   | 873359   | 33.803  | ng    | 99       |
| 4) n-Nitrosodimethylamine     | 3.563  | 42   | 394723   | 40.145  | ng    | 98       |
| 6) Aniline                    | 7.099  | 93   | 1051356  | 41.300  | ng    | 100      |
| 8) 2-Chlorophenol             | 7.340  | 128  | 1059061  | 43.280  | ng    | 99       |
| 9) Benzaldehyde               | 6.910  | 77   | 469566   | 30.800  | ng    | 99       |
| 10) Phenol                    | 6.981  | 94   | 1238080  | 42.705  | ng    | 99       |
| 11) bis(2-Chloroethyl)ether   | 7.193  | 93   | 972046   | 42.756  | ng    | 99       |
| 12) 1,3-Dichlorobenzene       | 7.657  | 146  | 1209916  | 41.821  | ng    | 99       |
| 13) 1,4-Dichlorobenzene       | 7.804  | 146  | 1217909  | 41.838  | ng    | 99       |
| 14) 1,2-Dichlorobenzene       | 8.116  | 146  | 1173072  | 42.784  | ng    | 99       |
| 15) Benzyl Alcohol            | 8.010  | 79   | 807273   | 43.152  | ng    | 99       |
| 16) 2,2'-oxybis(1-Chloropr... | 8.293  | 45   | 1141979  | 45.066  | ng    | 99       |
| 17) 2-Methylphenol            | 8.228  | 107  | 791650   | 44.311  | ng    | 99       |
| 18) Hexachloroethane          | 8.845  | 117  | 419803   | 41.451  | ng    | 95       |
| 19) n-Nitroso-di-n-propyla... | 8.575  | 70   | 693522   | 42.149  | ng    | 100      |
| 20) 3+4-Methylphenols         | 8.551  | 107  | 1057066  | 43.399  | ng    | 99       |
| 22) Acetophenone              | 8.593  | 105  | 1378862  | 41.167  | ng    | # 99     |
| 24) Nitrobenzene              | 8.969  | 77   | 993197   | 41.676  | ng    | 97       |
| 25) Isophorone                | 9.498  | 82   | 1842352  | 44.436  | ng    | 99       |
| 26) 2-Nitrophenol             | 9.681  | 139  | 546955   | 43.302  | ng    | 99       |
| 27) 2,4-Dimethylphenol        | 9.751  | 122  | 870924   | 60.841  | ng    | 99       |
| 28) bis(2-Chloroethoxy)met... | 9.975  | 93   | 1162126  | 41.868  | ng    | 99       |
| 29) 2,4-Dichlorophenol        | 10.222 | 162  | 990222   | 41.615  | ng    | 100      |
| 30) 1,2,4-Trichlorobenzene    | 10.428 | 180  | 1161491  | 42.194  | ng    | 99       |
| 31) Naphthalene               | 10.610 | 128  | 2846837  | 40.200  | ng    | 99       |
| 32) Benzoic acid              | 9.916  | 122  | 547361m  | 33.900  | ng    |          |
| 33) 4-Chloroaniline           | 10.728 | 127  | 575602   | 23.018  | ng    | 99       |
| 34) Hexachlorobutadiene       | 10.898 | 225  | 743966   | 43.693  | ng    | 99       |
| 35) Caprolactam               | 11.522 | 113  | 243129   | 35.625  | ng    | 94       |
| 36) 4-Chloro-3-methylphenol   | 11.869 | 107  | 827174   | 39.686  | ng    | 99       |
| 37) 2-Methylnaphthalene       | 12.228 | 142  | 1836048  | 36.798  | ng    | 99       |
| 38) 1-Methylnaphthalene       | 12.445 | 142  | 1933779  | 39.782  | ng    | 100      |
| 40) 1,2,4,5-Tetrachloroben... | 12.604 | 216  | 1290451  | 45.516  | ng    | 99       |
| 41) Hexachlorocyclopentadiene | 12.581 | 237  | 1664984  | 167.595 | ng    | 98       |
| 43) 2,4,6-Trichlorophenol     | 12.851 | 196  | 801684   | 44.436  | ng    | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM042125\  
 Data File : BM049980.D  
 Acq On : 21 Apr 2025 11:15  
 Operator : RC/JU  
 Sample : PB167645BS  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 PB167645BS

Quant Time: Apr 21 11:59:36 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM040825.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Apr 09 04:00:55 2025  
 Response via : Initial Calibration

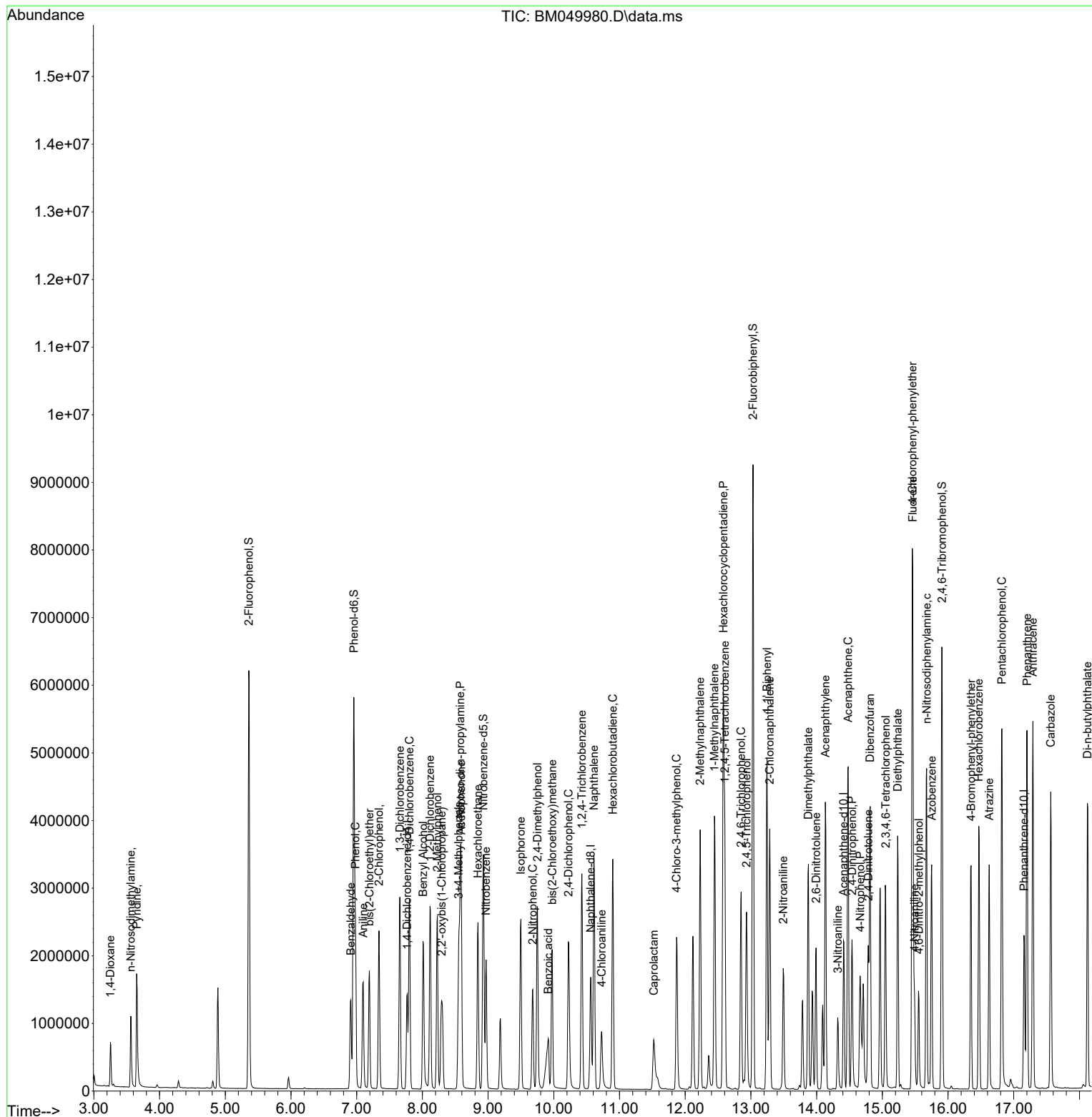
| Compound                       | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|--------|------|----------|--------|-------|----------|
| 44) 2,4,5-Trichlorophenol      | 12.933 | 196  | 852296   | 43.472 | ng    | 99       |
| 46) 1,1'-Biphenyl              | 13.245 | 154  | 2574209  | 42.731 | ng    | 99       |
| 47) 2-Chloronaphthalene        | 13.286 | 162  | 2039547  | 43.074 | ng    | 99       |
| 48) 2-Nitroaniline             | 13.492 | 65   | 468922   | 42.808 | ng    | 98       |
| 49) Acenaphthylene             | 14.133 | 152  | 3135276  | 45.457 | ng    | 100      |
| 50) Dimethylphthalate          | 13.875 | 163  | 2323401  | 40.638 | ng    | 100      |
| 51) 2,6-Dinitrotoluene         | 13.992 | 165  | 506087   | 42.250 | ng    | 97       |
| 52) Acenaphthene               | 14.480 | 154  | 1796540  | 40.950 | ng    | 99       |
| 53) 3-Nitroaniline             | 14.322 | 138  | 311855   | 26.937 | ng    | 94       |
| 54) 2,4-Dinitrophenol          | 14.539 | 184  | 620782   | 72.325 | ng    | 96       |
| 55) Dibenzofuran               | 14.816 | 168  | 2916639  | 39.768 | ng    | 98       |
| 56) 4-Nitrophenol              | 14.663 | 139  | 761858   | 73.846 | ng    | 99       |
| 57) 2,4-Dinitrotoluene         | 14.786 | 165  | 688368   | 43.076 | ng    | 97       |
| 58) Fluorene                   | 15.463 | 166  | 2364711  | 40.563 | ng    | 99       |
| 59) 2,3,4,6-Tetrachlorophenol  | 15.051 | 232  | 682403   | 39.712 | ng    | 97       |
| 60) Diethylphthalate           | 15.233 | 149  | 2076932  | 37.881 | ng    | 99       |
| 61) 4-Chlorophenyl-phenylether | 15.457 | 204  | 1297022  | 41.498 | ng    | 98       |
| 62) 4-Nitroaniline             | 15.492 | 138  | 454257   | 37.942 | ng    | 99       |
| 63) Azobenzene                 | 15.751 | 77   | 1784083  | 38.007 | ng    | 96       |
| 65) 4,6-Dinitro-2-methylph...  | 15.557 | 198  | 389517   | 44.000 | ng    | 96       |
| 66) n-Nitrosodiphenylamine     | 15.674 | 169  | 1955452  | 46.512 | ng    | 99       |
| 67) 4-Bromophenyl-phenylether  | 16.351 | 248  | 756840   | 46.408 | ng    | 94       |
| 68) Hexachlorobenzene          | 16.468 | 284  | 872280   | 46.630 | ng    | 99       |
| 69) Atrazine                   | 16.627 | 200  | 657210   | 60.291 | ng    | 97       |
| 70) Pentachlorophenol          | 16.821 | 266  | 1088541  | 79.040 | ng    | 99       |
| 71) Phenanthrene               | 17.204 | 178  | 3394421  | 44.072 | ng    | 100      |
| 72) Anthracene                 | 17.292 | 178  | 3404218  | 44.851 | ng    | 100      |
| 73) Carbazole                  | 17.563 | 167  | 2965804  | 41.486 | ng    | 99       |
| 74) Di-n-butylphthalate        | 18.121 | 149  | 3247632  | 39.441 | ng    | 99       |
| 75) Fluoranthene               | 19.221 | 202  | 3793074  | 40.054 | ng    | 99       |
| 77) Benzidine                  | 19.404 | 184  | 1414967  | 92.414 | ng    | 100      |
| 78) Pyrene                     | 19.580 | 202  | 3840802  | 48.287 | ng    | 99       |
| 80) Butylbenzylphthalate       | 20.480 | 149  | 1310426  | 43.845 | ng    | 96       |
| 81) Benzo(a)anthracene         | 21.380 | 228  | 3470021  | 45.239 | ng    | 100      |
| 82) 3,3'-Dichlorobenzidine     | 21.303 | 252  | 910767   | 31.442 | ng    | 99       |
| 83) Chrysene                   | 21.445 | 228  | 3150721  | 43.206 | ng    | 100      |
| 84) Bis(2-ethylhexyl)phtha...  | 21.298 | 149  | 1796205  | 41.249 | ng    | 99       |
| 85) Di-n-octyl phthalate       | 22.433 | 149  | 2905764  | 38.144 | ng    | 100      |
| 87) Indeno(1,2,3-cd)pyrene     | 27.803 | 276  | 3835487  | 44.759 | ng    | 99       |
| 88) Benzo(b)fluoranthene       | 23.456 | 252  | 3132543  | 42.666 | ng    | 99       |
| 89) Benzo(k)fluoranthene       | 23.515 | 252  | 3120017  | 43.851 | ng    | 100      |
| 90) Benzo(a)pyrene             | 24.262 | 252  | 3005590  | 46.586 | ng    | 99       |
| 91) Dibenzo(a,h)anthracene     | 27.844 | 278  | 3122653  | 44.240 | ng    | 100      |
| 92) Benzo(g,h,i)perylene       | 28.856 | 276  | 3029590  | 42.003 | ng    | 99       |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM042125\  
Data File : BM049980.D  
Acq On : 21 Apr 2025 11:15  
Operator : RC/JU  
Sample : PB167645BS  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
PB167645BS

Quant Time: Apr 21 11:59:36 2025  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM040825.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Wed Apr 09 04:00:55 2025  
Response via : Initial Calibration



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM042125\  
Data File : BM049980.D  
Acq On : 21 Apr 2025 11:15  
Operator : RC/JU  
Sample : PB167645BS  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
PB167645BS

Quant Time: Apr 21 11:59:36 2025  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM040825.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Wed Apr 09 04:00:55 2025  
Response via : Initial Calibration

