

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM020924\  
 Data File : BM044047.D  
 Acq On : 09 Feb 2024 20:19  
 Operator : MA/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTDCCC040EC

Quant Time: Feb 09 22:58:38 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM020824.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Feb 09 02:42:01 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.910  | 152  | 484814   | 20.000 | ng    | 0.00     |
| 21) Naphthalene-d8            | 10.710 | 136  | 2018789  | 20.000 | ng    | 0.00     |
| 39) Acenaphthene-d10          | 14.551 | 164  | 1121571  | 20.000 | ng    | 0.00     |
| 64) Phenanthrene-d10          | 17.304 | 188  | 2101339  | 20.000 | ng    | 0.00     |
| 76) Chrysene-d12              | 21.503 | 240  | 1621850  | 20.000 | ng    | 0.00     |
| 86) Perylene-d12              | 23.897 | 264  | 1859389  | 20.000 | ng    | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 5) 2-Fluorophenol             | 5.475  | 112  | 2570673  | 77.120 | ng    | 0.00     |
| 7) Phenol-d6                  | 7.075  | 99   | 3353327  | 79.066 | ng    | 0.00     |
| 23) Nitrobenzene-d5           | 9.081  | 82   | 3059386  | 80.416 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol      | 16.045 | 330  | 955206   | 78.660 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl          | 13.180 | 172  | 6188525  | 77.873 | ng    | 0.00     |
| 79) Terphenyl-d14             | 19.945 | 244  | 7216577  | 78.003 | ng    | 0.00     |
| Target Compounds              |        |      |          |        |       |          |
|                               |        |      |          |        |       | Qvalue   |
| 2) 1,4-Dioxane                | 3.387  | 88   | 492245   | 37.210 | ng    | 99       |
| 3) Pyridine                   | 3.781  | 79   | 1345340m | 38.520 | ng    |          |
| 4) n-Nitrosodimethylamine     | 3.704  | 42   | 542557   | 40.943 | ng    | 98       |
| 6) Aniline                    | 7.240  | 93   | 1665493  | 39.939 | ng    | 98       |
| 8) 2-Chlorophenol             | 7.469  | 128  | 1351782  | 38.834 | ng    | 100      |
| 9) Benzaldehyde               | 7.057  | 77   | 853351   | 39.153 | ng    | 99       |
| 10) Phenol                    | 7.104  | 94   | 1706005  | 39.857 | ng    | 99       |
| 11) bis(2-Chloroethyl)ether   | 7.345  | 93   | 1423282  | 40.154 | ng    | 98       |
| 12) 1,3-Dichlorobenzene       | 7.792  | 146  | 1532931  | 38.787 | ng    | 99       |
| 13) 1,4-Dichlorobenzene       | 7.945  | 146  | 1548341  | 38.866 | ng    | 99       |
| 14) 1,2-Dichlorobenzene       | 8.257  | 146  | 1463738  | 38.561 | ng    | 100      |
| 15) Benzyl Alcohol            | 8.157  | 79   | 1106504  | 40.069 | ng    | 99       |
| 16) 2,2'-oxybis(1-Chloropr... | 8.451  | 45   | 1761590m | 39.204 | ng    |          |
| 17) 2-Methylphenol            | 8.351  | 107  | 1112750  | 39.603 | ng    | 98       |
| 18) Hexachloroethane          | 8.981  | 117  | 584691   | 39.199 | ng    | 98       |
| 19) n-Nitroso-di-n-propyla... | 8.728  | 70   | 1036227  | 39.789 | ng    | 99       |
| 20) 3+4-Methylphenols         | 8.686  | 107  | 1540340  | 39.831 | ng    | 98       |
| 22) Acetophenone              | 8.739  | 105  | 2028588  | 39.588 | ng    | 100      |
| 24) Nitrobenzene              | 9.122  | 77   | 1557476  | 39.977 | ng    | 99       |
| 25) Isophorone                | 9.645  | 82   | 2918070  | 39.833 | ng    | 99       |
| 26) 2-Nitrophenol             | 9.828  | 139  | 729503   | 40.387 | ng    | 100      |
| 27) 2,4-Dimethylphenol        | 9.886  | 122  | 909872   | 39.395 | ng    | 97       |
| 28) bis(2-Chloroethoxy)met... | 10.139 | 93   | 1856451  | 39.184 | ng    | 99       |
| 29) 2,4-Dichlorophenol        | 10.357 | 162  | 1205562  | 39.646 | ng    | 97       |
| 30) 1,2,4-Trichlorobenzene    | 10.569 | 180  | 1345379  | 39.347 | ng    | 99       |
| 31) Naphthalene               | 10.763 | 128  | 4338371  | 38.937 | ng    | 99       |
| 32) Benzoic acid              | 10.039 | 122  | 877358   | 37.711 | ng    | 99       |
| 33) 4-Chloroaniline           | 10.880 | 127  | 1642806  | 38.938 | ng    | 99       |
| 34) Hexachlorobutadiene       | 11.033 | 225  | 770122   | 39.349 | ng    | 99       |
| 35) Caprolactam               | 11.675 | 113  | 377009   | 38.529 | ng    | 98       |
| 36) 4-Chloro-3-methylphenol   | 12.004 | 107  | 1303222  | 39.579 | ng    | 100      |
| 37) 2-Methylnaphthalene       | 12.374 | 142  | 2869985  | 38.670 | ng    | 97       |
| 38) 1-Methylnaphthalene       | 12.592 | 142  | 2827058  | 38.765 | ng    | 100      |
| 40) 1,2,4,5-Tetrachloroben... | 12.733 | 216  | 1295899  | 39.751 | ng    | 100      |
| 41) Hexachlorocyclopentadiene | 12.710 | 237  | 568911   | 39.988 | ng    | 99       |
| 43) 2,4,6-Trichlorophenol     | 12.980 | 196  | 891684   | 40.191 | ng    | 100      |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM020924\  
 Data File : BM044047.D  
 Acq On : 09 Feb 2024 20:19  
 Operator : MA/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTDCCC040EC

Quant Time: Feb 09 22:58:38 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM020824.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Feb 09 02:42:01 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 13.051 | 196  | 979961   | 40.249 | ng    | 98       |
| 46) 1,1'-Biphenyl             | 13.386 | 154  | 3490408  | 39.439 | ng    | 99       |
| 47) 2-Chloronaphthalene       | 13.427 | 162  | 2779179  | 39.643 | ng    | 99       |
| 48) 2-Nitroaniline            | 13.639 | 65   | 815036   | 41.326 | ng    | 94       |
| 49) Acenaphthylene            | 14.274 | 152  | 4131907  | 39.506 | ng    | 100      |
| 50) Dimethylphthalate         | 14.027 | 163  | 3200163  | 38.483 | ng    | 99       |
| 51) 2,6-Dinitrotoluene        | 14.145 | 165  | 724889   | 39.510 | ng    | 98       |
| 52) Acenaphthene              | 14.616 | 154  | 2524567  | 39.001 | ng    | 100      |
| 53) 3-Nitroaniline            | 14.468 | 138  | 785149   | 39.820 | ng    | 97       |
| 54) 2,4-Dinitrophenol         | 14.674 | 184  | 373597   | 37.138 | ng    | 98       |
| 55) Dibenzofuran              | 14.951 | 168  | 3870877  | 38.636 | ng    | 98       |
| 56) 4-Nitrophenol             | 14.768 | 139  | 626670   | 39.097 | ng    | 95       |
| 57) 2,4-Dinitrotoluene        | 14.927 | 165  | 957026   | 39.910 | ng    | 99       |
| 58) Fluorene                  | 15.604 | 166  | 2932442  | 37.878 | ng    | 99       |
| 59) 2,3,4,6-Tetrachlorophenol | 15.174 | 232  | 741134   | 38.793 | ng    | 98       |
| 60) Diethylphthalate          | 15.392 | 149  | 3261297  | 37.948 | ng    | 99       |
| 61) 4-Chlorophenyl-phenyle... | 15.604 | 204  | 1435446  | 38.266 | ng    | 99       |
| 62) 4-Nitroaniline            | 15.633 | 138  | 784779   | 39.540 | ng    | 99       |
| 63) Azobenzene                | 15.898 | 77   | 3183253  | 39.479 | ng    | 99       |
| 65) 4,6-Dinitro-2-methylph... | 15.686 | 198  | 509245   | 40.669 | ng    | 98       |
| 66) n-Nitrosodiphenylamine    | 15.821 | 169  | 2547473  | 39.903 | ng    | 100      |
| 67) 4-Bromophenyl-phenylether | 16.498 | 248  | 809297   | 38.954 | ng    | 99       |
| 68) Hexachlorobenzene         | 16.598 | 284  | 999030   | 39.255 | ng    | 99       |
| 69) Atrazine                  | 16.780 | 200  | 699001   | 40.436 | ng    | 99       |
| 70) Pentachlorophenol         | 16.945 | 266  | 676335   | 40.852 | ng    | 99       |
| 71) Phenanthrene              | 17.345 | 178  | 4327186  | 38.996 | ng    | 99       |
| 72) Anthracene                | 17.439 | 178  | 4320646  | 39.331 | ng    | 99       |
| 73) Carbazole                 | 17.715 | 167  | 4184530  | 39.168 | ng    | 99       |
| 74) Di-n-butylphthalate       | 18.292 | 149  | 5411236  | 38.967 | ng    | 100      |
| 75) Fluoranthene              | 19.368 | 202  | 4792511  | 38.449 | ng    | 99       |
| 77) Benzidine                 | 19.562 | 184  | 1187314  | 47.108 | ng    | 99       |
| 78) Pyrene                    | 19.727 | 202  | 4957718  | 39.691 | ng    | 99       |
| 80) Butylbenzylphthalate      | 20.650 | 149  | 2220251  | 39.988 | ng    | 99       |
| 81) Benzo(a)anthracene        | 21.492 | 228  | 4406996  | 39.064 | ng    | 99       |
| 82) 3,3'-Dichlorobenzidine    | 21.427 | 252  | 1464668  | 39.569 | ng    | 99       |
| 83) Chrysene                  | 21.544 | 228  | 4179189  | 39.210 | ng    | 100      |
| 84) Bis(2-ethylhexyl)phtha... | 21.439 | 149  | 3210236  | 39.187 | ng    | 100      |
| 85) Di-n-octyl phthalate      | 22.374 | 149  | 5575197  | 38.968 | ng    | 100      |
| 87) Indeno(1,2,3-cd)pyrene    | 26.391 | 276  | 5137573  | 39.039 | ng    | 100      |
| 88) Benzo(b)fluoranthene      | 23.174 | 252  | 4586070  | 39.456 | ng    | 100      |
| 89) Benzo(k)fluoranthene      | 23.221 | 252  | 4355540  | 39.436 | ng    | 100      |
| 90) Benzo(a)pyrene            | 23.791 | 252  | 4011646  | 39.514 | ng    | 100      |
| 91) Dibenzo(a,h)anthracene    | 26.415 | 278  | 4306873  | 39.125 | ng    | 100      |
| 92) Benzo(g,h,i)perylene      | 27.156 | 276  | 4338135  | 38.858 | ng    | 100      |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM020924\  
Data File : BM044047.D  
Acq On : 09 Feb 2024 20:19  
Operator : MA/JU  
Sample : SSTDCCC040  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTDCCC040EC

Quant Time: Feb 09 22:58:38 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM020824.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Fri Feb 09 02:42:01 2024  
Response via : Initial Calibration

