

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP062821\  
 Data File : BP006074.D  
 Acq On : 28 Jun 2021 13:15  
 Operator : CG/JU  
 Sample : M2824-02MSD  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 ABN-1MSD

Manual Integrations  
 APPROVED

mohammad  
 6/29/2021 2:29:54 PM

Quant Time: Jun 28 14:04:00 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP060821.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jun 28 11:34:30 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.904  | 152  | 70816    | 20.000 | ng     | 0.00     |
| 21) Naphthalene-d8            | 10.704 | 136  | 272936   | 20.000 | ng     | 0.00     |
| 39) Acenaphthene-d10          | 14.533 | 164  | 167905   | 20.000 | ng     | 0.00     |
| 64) Phenanthrene-d10          | 17.280 | 188  | 325755   | 20.000 | ng     | 0.00     |
| 76) Chrysene-d12              | 21.357 | 240  | 374583   | 20.000 | ng     | 0.00     |
| 86) Perylene-d12              | 23.756 | 264  | 435095   | 20.000 | ng     | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 5) 2-Fluorophenol             | 5.487  | 112  | 19679    | 4.459  | ng     | 0.01     |
| 7) Phenol-d6                  | 7.087  | 99   | 110171   | 19.260 | ng     | 0.00     |
| 23) Nitrobenzene-d5           | 9.069  | 82   | 363756   | 65.341 | ng     | 0.00     |
| 42) 2,4,6-Tribromophenol      | 16.027 | 330  | 6010     | 2.086  | ng     | 0.00     |
| 45) 2-Fluorobiphenyl          | 13.163 | 172  | 803254   | 62.820 | ng     | 0.00     |
| 79) Terphenyl-d14             | 19.827 | 244  | 1256242  | 62.798 | ng     | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.340  | 88   | 68486    | 34.349 | ng     | 97       |
| 3) Pyridine                   | 3.757  | 79   | 172225   | 36.752 | ng     | 98       |
| 4) n-Nitrosodimethylamine     | 3.663  | 42   | 81511    | 37.421 | ng     | 87       |
| 6) Aniline                    | 7.234  | 93   | 274639   | 43.005 | ng     | 99       |
| 8) 2-Chlorophenol             | 7.469  | 128  | 236196   | 46.678 | ng     | 99       |
| 9) Benzaldehyde               | 7.045  | 77   | 131642   | 53.984 | ng     | 98       |
| 10) Phenol                    | 7.110  | 94   | 277806   | 48.616 | ng     | 96       |
| 11) bis(2-Chloroethyl)ether   | 7.328  | 93   | 205652   | 47.493 | ng     | 99       |
| 12) 1,3-Dichlorobenzene       | 7.793  | 146  | 249457   | 44.089 | ng     | 98       |
| 13) 1,4-Dichlorobenzene       | 7.940  | 146  | 255450   | 44.271 | ng     | 99       |
| 14) 1,2-Dichlorobenzene       | 8.257  | 146  | 244019   | 44.243 | ng     | 99       |
| 15) Benzyl Alcohol            | 8.151  | 79   | 201075   | 46.669 | ng     | 97       |
| 16) 2,2'-oxybis(1-Chloropr... | 8.440  | 45   | 174137   | 43.504 | ng     | 97       |
| 17) 2-Methylphenol            | 8.363  | 107  | 189458   | 48.661 | ng     | 97       |
| 18) Hexachloroethane          | 8.981  | 117  | 90159    | 43.197 | ng     | 97       |
| 19) n-Nitroso-di-n-propyla... | 8.716  | 70   | 155413   | 44.164 | ng     | 95       |
| 20) 3+4-Methylphenols         | 8.692  | 107  | 251698   | 47.859 | ng     | 95       |
| 22) Acetophenone              | 8.734  | 105  | 324743   | 45.204 | ng     | 97       |
| 24) Nitrobenzene              | 9.116  | 77   | 238030   | 41.175 | ng     | 99       |
| 25) Isophorone                | 9.645  | 82   | 410888   | 39.967 | ng     | 99       |
| 26) 2-Nitrophenol             | 9.822  | 139  | 123415   | 44.744 | ng     | 94       |
| 27) 2,4-Dimethylphenol        | 9.892  | 122  | 206939   | 50.360 | ng     | 99       |
| 28) bis(2-Chloroethoxy)met... | 10.122 | 93   | 254857   | 47.499 | ng     | 99       |
| 29) 2,4-Dichlorophenol        | 10.369 | 162  | 215886   | 43.717 | ng     | 99       |
| 30) 1,2,4-Trichlorobenzene    | 10.569 | 180  | 230128   | 41.474 | ng     | 97       |
| 31) Naphthalene               | 10.757 | 128  | 674879   | 42.736 | ng     | 99       |
| 32) Benzoic acid              | 10.081 | 122  | 131836   | 43.503 | ng     | 97       |
| 33) 4-Chloroaniline           | 10.875 | 127  | 169979   | 26.644 | ng     | 99       |
| 34) Hexachlorobutadiene       | 11.039 | 225  | 149320   | 39.786 | ng     | 97       |
| 35) Caprolactam               | 11.686 | 113  | 66991    | 47.097 | ng     | 96       |
| 36) 4-Chloro-3-methylphenol   | 12.010 | 107  | 226848   | 45.241 | ng     | 97       |
| 37) 2-Methylnaphthalene       | 12.363 | 142  | 476422   | 42.373 | ng     | 98       |
| 38) 1-Methylnaphthalene       | 12.580 | 142  | 442878   | 41.817 | ng     | 100      |
| 40) 1,2,4,5-Tetrachloroben... | 12.733 | 216  | 254339   | 43.034 | ng     | 98       |
| 41) Hexachlorocyclopentadiene | 12.710 | 237  | 217567   | 73.730 | ng     | 98       |
| 43) 2,4,6-Trichlorophenol     | 12.980 | 196  | 170267   | 42.002 | ng     | 97       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP062821\  
 Data File : BP006074.D  
 Acq On : 28 Jun 2021 13:15  
 Operator : CG/JU  
 Sample : M2824-02MSD  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 ABN-1MSD

Manual Integrations  
 APPROVED

mohammad  
 6/29/2021 2:29:54 PM

Quant Time: Jun 28 14:04:00 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP060821.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jun 28 11:34:30 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 44) 2,4,5-Trichlorophenol     | 13.063 | 196  | 183648   | 42.066 | ng    | 96       |
| 46) 1,1'-Biphenyl             | 13.369 | 154  | 588393   | 43.463 | ng    | 99       |
| 47) 2-Chloronaphthalene       | 13.416 | 162  | 467331   | 42.274 | ng    | 99       |
| 48) 2-Nitroaniline            | 13.627 | 65   | 129319   | 44.491 | ng    | 97       |
| 49) Acenaphthylene            | 14.257 | 152  | 745108   | 43.933 | ng    | 99       |
| 50) Dimethylphthalate         | 13.998 | 163  | 598497   | 43.353 | ng    | 100      |
| 51) 2,6-Dinitrotoluene        | 14.122 | 165  | 133483   | 42.540 | ng    | 96       |
| 52) Acenaphthene              | 14.598 | 154  | 437490   | 42.477 | ng    | 98       |
| 53) 3-Nitroaniline            | 14.451 | 138  | 98906    | 32.522 | ng    | 95       |
| 54) 2,4-Dinitrophenol         | 14.669 | 184  | 140426   | 75.535 | ng    | 97       |
| 55) Dibenzofuran              | 14.933 | 168  | 702297   | 41.463 | ng    | 99       |
| 56) 4-Nitrophenol             | 14.786 | 139  | 204245   | 82.857 | ng    | 89       |
| 57) 2,4-Dinitrotoluene        | 14.910 | 165  | 183399   | 43.700 | ng    | 99       |
| 58) Fluorene                  | 15.580 | 166  | 566704   | 42.942 | ng    | 98       |
| 59) 2,3,4,6-Tetrachlorophenol | 15.169 | 232  | 163120m  | 42.394 | ng    |          |
| 60) Diethylphthalate          | 15.357 | 149  | 593822   | 42.873 | ng    | 99       |
| 61) 4-Chlorophenyl-phenyle... | 15.574 | 204  | 288580   | 42.170 | ng    | 99       |
| 62) 4-Nitroaniline            | 15.616 | 138  | 137213   | 43.835 | ng    | 94       |
| 63) Azobenzene                | 15.869 | 77   | 515931   | 41.172 | ng    | 94       |
| 65) 4,6-Dinitro-2-methylph... | 15.680 | 198  | 96920    | 40.487 | ng    | 96       |
| 66) n-Nitrosodiphenylamine    | 15.792 | 169  | 498154   | 46.115 | ng    | 98       |
| 67) 4-Bromophenyl-phenylether | 16.469 | 248  | 186413   | 44.584 | ng    | 98       |
| 68) Hexachlorobenzene         | 16.586 | 284  | 223624   | 44.621 | ng    | 98       |
| 69) Atrazine                  | 16.745 | 200  | 184791   | 54.983 | ng    | 98       |
| 70) Pentachlorophenol         | 16.939 | 266  | 252908   | 90.757 | ng    | 100      |
| 71) Phenanthrene              | 17.321 | 178  | 886281   | 45.304 | ng    | 100      |
| 72) Anthracene                | 17.416 | 178  | 894924   | 46.481 | ng    | 100      |
| 73) Carbazole                 | 17.692 | 167  | 810940   | 43.869 | ng    | 100      |
| 74) Di-n-butylphthalate       | 18.233 | 149  | 1009589  | 47.139 | ng    | 99       |
| 75) Fluoranthene              | 19.286 | 202  | 1060813  | 44.651 | ng    | 98       |
| 77) Benzidine                 | 19.468 | 184  | 662086   | 84.945 | ng    | 99       |
| 78) Pyrene                    | 19.639 | 202  | 1108104  | 42.375 | ng    | 100      |
| 80) Butylbenzylphthalate      | 20.504 | 149  | 482183   | 44.819 | ng    | 96       |
| 81) Benzo(a)anthracene        | 21.345 | 228  | 1146049  | 42.384 | ng    | 99       |
| 82) 3,3'-Dichlorobenzidine    | 21.274 | 252  | 393771   | 42.122 | ng    | 98       |
| 83) Chrysene                  | 21.398 | 228  | 1128048  | 43.072 | ng    | 99       |
| 84) Bis(2-ethylhexyl)phtha... | 21.256 | 149  | 715602   | 45.353 | ng    | 99       |
| 85) Di-n-octyl phthalate      | 22.174 | 149  | 1286881  | 45.591 | ng    | 99       |
| 87) Indeno(1,2,3-cd)pyrene    | 26.286 | 276  | 1721161  | 46.259 | ng    | 98       |
| 88) Benzo(b)fluoranthene      | 23.021 | 252  | 1297765  | 43.549 | ng    | 98       |
| 89) Benzo(k)fluoranthene      | 23.074 | 252  | 1307336  | 45.242 | ng    | 99       |
| 90) Benzo(a)pyrene            | 23.650 | 252  | 1314099  | 46.904 | ng    | 98       |
| 91) Dibenzo(a,h)anthracene    | 26.297 | 278  | 1462282  | 45.662 | ng    | 98       |
| 92) Benzo(g,h,i)perylene      | 27.062 | 276  | 1471375  | 46.511 | ng    | 98       |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP062821\  
Data File : BP006074.D  
Acq On : 28 Jun 2021 13:15  
Operator : CG/JU  
Sample : M2824-02MSD  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
ABN-1MSD

Manual Integrations  
APPROVED

mohammad  
6/29/2021 2:29:54 PM

Quant Time: Jun 28 14:04:00 2021  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP060821.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Mon Jun 28 11:34:30 2021  
Response via : Initial Calibration

