

Data Path : Z:\HPCHEM1\BNA G\DATA\BG031418\  
 Data File : BG033163.D  
 Acq On : 15 Mar 2018 3:15  
 Operator : SJ/JU  
 Sample : J1749-04MS  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BU-02-031318MS

Manual Integrations  
 APPROVED

Sohil  
 3/15/2018 7:00:15 PM

Quant Time: Mar 15 11:19:58 2018  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Mar 14 13:59:34 2018  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.90  | 152  | 38648    | 20.00 | ng    | 0.00     |
| 21) Naphthalene-d8        | 11.78 | 136  | 176432   | 20.00 | ng    | 0.00     |
| 38) Acenaphthene-d10      | 15.52 | 164  | 111139   | 20.00 | ng    | 0.00     |
| 63) Phenanthrene-d10      | 18.24 | 188  | 265164   | 20.00 | ng    | 0.00     |
| 75) Chrysene-d12          | 22.60 | 240  | 241561   | 20.00 | ng    | 0.00     |
| 86) Perylene-d12          | 26.40 | 264  | 261715   | 20.00 | ng    | -0.01    |

## System Monitoring Compounds

|                          |       |     |         |        |    |      |
|--------------------------|-------|-----|---------|--------|----|------|
| 5) 2-Fluorophenol        | 6.32  | 112 | 350833  | 145.23 | ng | 0.00 |
| 7) Phenol-d6             | 8.01  | 99  | 445582  | 132.33 | ng | 0.00 |
| 23) Nitrobenzene-d5      | 10.10 | 82  | 317523  | 120.24 | ng | 0.00 |
| 41) 2,4,6-Tribromophenol | 16.99 | 330 | 260917  | 223.28 | ng | 0.00 |
| 44) 2-Fluorobiphenyl     | 14.14 | 172 | 719984  | 93.43  | ng | 0.00 |
| 78) Terphenyl-d14        | 20.76 | 244 | 1255036 | 116.73 | ng | 0.00 |

## Target Compounds

|                                |       |     |        |        |    | Qvalue |
|--------------------------------|-------|-----|--------|--------|----|--------|
| 2) 1,4-Dioxane                 | 3.96  | 88  | 31219  | 29.777 | ng | # 92   |
| 3) Pyridine                    | 4.43  | 79  | 78177  | 27.054 | ng | 93     |
| 4) n-Nitrosodimethylamine      | 4.32  | 42  | 57954  | 50.024 | ng | # 82   |
| 6) Aniline                     | 8.19  | 93  | 60189  | 13.206 | ng | 96     |
| 8) 2-Chlorophenol              | 8.45  | 128 | 120967 | 45.749 | ng | 98     |
| 9) Benzaldehyde                | 8.00  | 77  | 32798  | 16.901 | ng | 96     |
| 10) Phenol                     | 8.03  | 94  | 137631 | 40.548 | ng | 95     |
| 11) bis(2-Chloroethyl)ether    | 8.28  | 93  | 108581 | 39.121 | ng | 99     |
| 12) 1,3-Dichlorobenzene        | 8.78  | 146 | 117687 | 38.001 | ng | 99     |
| 13) 1,4-Dichlorobenzene        | 8.93  | 146 | 120781 | 38.744 | ng | 96     |
| 14) 1,2-Dichlorobenzene        | 9.27  | 146 | 117260 | 39.253 | ng | 99     |
| 15) Benzyl Alcohol             | 9.13  | 79  | 109684 | 44.310 | ng | 95     |
| 16) 2,2'-oxybis(1-Chloropropan | 9.42  | 45  | 210642 | 44.147 | ng | 99     |
| 17) 2-Methylphenol             | 9.33  | 107 | 105476 | 42.140 | ng | 96     |
| 18) Hexachloroethane           | 10.02 | 117 | 43908  | 41.368 | ng | 90     |
| 19) n-Nitroso-di-n-propylamine | 9.71  | 70  | 96680  | 40.671 | ng | # 89   |
| 20) 3+4-Methylphenols          | 9.67  | 107 | 146869 | 42.201 | ng | 95     |
| 22) Acetophenone               | 9.74  | 105 | 181994 | 40.845 | ng | # 98   |
| 24) Nitrobenzene               | 10.14 | 77  | 141095 | 49.861 | ng | 94     |
| 25) Isophorone                 | 10.67 | 82  | 280294 | 45.262 | ng | # 94   |
| 26) 2-Nitrophenol              | 10.87 | 139 | 67683  | 61.756 | ng | 98     |
| 27) 2,4-Dimethylphenol         | 10.90 | 122 | 133005 | 49.441 | ng | 92     |
| 28) bis(2-Chloroethoxy)methane | 11.14 | 93  | 155234 | 43.030 | ng | 96     |
| 29) 2,4-Dichlorophenol         | 11.41 | 162 | 127094 | 52.054 | ng | 94     |
| 30) 1,2,4-Trichlorobenzene     | 11.64 | 180 | 116518 | 40.712 | ng | 96     |
| 31) Naphthalene                | 11.84 | 128 | 393509 | 42.684 | ng | 99     |
| 32) Benzoic acid               | 11.04 | 122 | 71383m | 44.221 | ng |        |
| 33) 4-Chloroaniline            | 11.93 | 127 | 24525  | 5.950  | ng | 98     |
| 34) Hexachlorobutadiene        | 12.09 | 225 | 68398  | 42.606 | ng | 97     |
| 35) Caprolactam                | 12.68 | 113 | 39667  | 40.574 | ng | 93     |
| 36) 4-Chloro-3-methylphenol    | 13.00 | 107 | 158567 | 55.808 | ng | 95     |
| 37) 2-Methylnaphthalene        | 13.39 | 142 | 294369 | 44.134 | ng | 94     |
| 39) 1,2,4,5-Tetrachlorobenzene | 13.74 | 216 | 136917 | 41.500 | ng | # 97   |
| 40) Hexachlorocyclopentadiene  | 13.71 | 237 | 165199 | 98.252 | ng | 93     |

Data Path : Z:\HPCHEM1\BNA G\DATA\BG031418\  
 Data File : BG033163.D  
 Acq On : 15 Mar 2018 3:15  
 Operator : SJ/JU  
 Sample : J1749-04MS  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BU-02-031318MS

Manual Integrations  
 APPROVED

Sohil  
 3/15/2018 7:00:15 PM

Quant Time: Mar 15 11:19:58 2018  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Mar 14 13:59:34 2018  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc    | Units | Dev(Min) |
|--------------------------------|-------|------|----------|---------|-------|----------|
| 42) 2,4,6-Trichlorophenol      | 13.97 | 196  | 104056   | 50.013  | nq    | 99       |
| 43) 2,4,5-Trichlorophenol      | 14.04 | 196  | 119844   | 49.768  | nq    | 96       |
| 45) 1,1'-Biphenyl              | 14.36 | 154  | 382401   | 39.375  | nq    | 97       |
| 46) 2-Chloronaphthalene        | 14.41 | 162  | 289009   | 39.837  | nq    | 97       |
| 47) 2-Nitroaniline             | 14.60 | 65   | 110139   | 57.699  | nq    | 93       |
| 48) Acenaphthylene             | 15.25 | 152  | 488327   | 41.114  | nq    | 99       |
| 49) Dimethylphthalate          | 14.94 | 163  | 411197   | 43.574  | nq    | 100      |
| 50) 2,6-Dinitrotoluene         | 15.07 | 165  | 89568    | 55.612  | nq    | 97       |
| 51) Acenaphthene               | 15.58 | 154  | 345099   | 46.653  | nq    | 99       |
| 52) 3-Nitroaniline             | 15.41 | 138  | 27933    | 19.727  | nq    | # 86     |
| 53) 2,4-Dinitrophenol          | 15.61 | 184  | 72358    | 122.188 | nq    | # 80     |
| 54) Dibenzofuran               | 15.91 | 168  | 472103   | 44.823  | nq    | 91       |
| 55) 4-Nitrophenol              | 15.69 | 139  | 148135   | 94.456  | nq    | 94       |
| 56) 2,4-Dinitrotoluene         | 15.85 | 165  | 130061   | 64.808  | nq    | # 89     |
| 57) Fluorene                   | 16.55 | 166  | 384014   | 44.174  | nq    | 99       |
| 58) 2,3,4,6-Tetrachlorophenol  | 16.12 | 232  | 105668m  | 56.519  | nq    |          |
| 59) Diethylphthalate           | 16.28 | 149  | 451309   | 45.343  | nq    | 98       |
| 60) 4-Chlorophenyl-phenylether | 16.52 | 204  | 197302   | 46.395  | nq    | 90       |
| 61) 4-Nitroaniline             | 16.56 | 138  | 67245    | 34.582  | nq    | 96       |
| 62) Azobenzene                 | 16.82 | 77   | 393057   | 44.142  | nq    | 96       |
| 64) 4,6-Dinitro-2-methylphenol | 16.61 | 198  | 57989    | 65.492  | nq    | 98       |
| 65) n-Nitrosodiphenylamine     | 16.73 | 169  | 354389   | 41.083  | nq    | 100      |
| 66) 4-Bromophenyl-phenylether  | 17.42 | 248  | 130156   | 46.421  | nq    | # 88     |
| 67) Hexachlorobenzene          | 17.55 | 284  | 138988   | 45.871  | nq    | 96       |
| 68) Atrazine                   | 17.66 | 200  | 110674   | 38.978  | nq    | 97       |
| 69) Pentachlorophenol          | 17.89 | 266  | 166854   | 87.426  | nq    | 98       |
| 70) Phenanthrene               | 18.29 | 178  | 644083   | 43.538  | nq    | 99       |
| 71) Anthracene                 | 18.38 | 178  | 653780   | 44.020  | nq    | 99       |
| 72) Carbazole                  | 18.64 | 167  | 598660   | 40.895  | nq    | 97       |
| 73) Di-n-butylphthalate        | 19.13 | 149  | 760790   | 42.362  | nq    | 99       |
| 74) Fluoranthene               | 20.24 | 202  | 734107   | 44.923  | nq    | 97       |
| 76) Benzidine                  | 20.40 | 184  | 107447m  | 13.049  | nq    |          |
| 77) Pyrene                     | 20.60 | 202  | 732715   | 43.012  | nq    | 99       |
| 79) Butylbenzylphthalate       | 21.46 | 149  | 338712   | 44.846  | nq    | 94       |
| 80) Benzo(a)anthracene         | 22.58 | 228  | 691231   | 44.624  | nq    | 100      |
| 81) 3,3'-Dichlorobenzidine     | 22.47 | 252  | 79610    | 13.877  | nq    | 97       |
| 82) Chrysene                   | 22.66 | 228  | 651738   | 44.045  | nq    | 98       |
| 83) Bis(2-ethylhexyl)phthalate | 22.40 | 149  | 465232   | 41.613  | nq    | 95       |
| 84) Di-n-octyl phthalate       | 23.81 | 149  | 783127   | 45.956  | nq    | 99       |
| 85) Indeno(1,2,3-cd)pyrene     | 30.83 | 276  | 814695   | 47.210  | nq    | # 89     |
| 87) Benzo(b)fluoranthene       | 25.18 | 252  | 700773   | 45.286  | nq    | 99       |
| 88) Benzo(k)fluoranthene       | 25.26 | 252  | 676442   | 43.985  | nq    | 99       |
| 89) Benzo(a)pyrene             | 26.23 | 252  | 668578   | 45.425  | nq    | 98       |
| 90) Dibenzo(a,h)anthracene     | 30.90 | 278  | 681413   | 45.995  | nq    | 96       |
| 91) Benzo(g,h,i)perylene       | 32.23 | 276  | 674716   | 46.200  | nq    | 99       |

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

