

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010419\  
 Data File : BG038962.D  
 Acq On : 5 Jan 2019 13:30  
 Operator : JU/SJ  
 Sample : IDOC-01  
 Misc : IDOC-BN-FE  
 ALS Vial : 33 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 IDOC-01

Manual Integrations  
 APPROVED

Sohil  
 1/7/2019 12:48:27 PM

Quant Time: Jan 06 23:15:33 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Dec 12 15:26:27 2018  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.05  | 152  | 27768    | 20.00 | ng    | -0.01    |
| 21) Naphthalene-d8        | 10.87 | 136  | 113533   | 20.00 | ng    | -0.01    |
| 39) Acenaphthene-d10      | 14.69 | 164  | 80281    | 20.00 | ng    | 0.00     |
| 64) Phenanthrene-d10      | 17.44 | 188  | 187285   | 20.00 | ng    | 0.00     |
| 76) Chrysene-d12          | 21.72 | 240  | 180675   | 20.00 | ng    | 0.00     |
| 87) Perylene-d12          | 25.02 | 264  | 197124   | 20.00 | ng    | 0.00     |

|                             |       |     |        |        |    |       |
|-----------------------------|-------|-----|--------|--------|----|-------|
| System Monitoring Compounds |       |     |        |        |    |       |
| 5) 2-Fluorophenol           | 5.60  | 112 | 191954 | 122.61 | ng | 0.00  |
| 7) Phenol-d6                | 7.21  | 99  | 269260 | 125.28 | ng | 0.00  |
| 23) Nitrobenzene-d5         | 9.23  | 82  | 164049 | 73.82  | ng | -0.01 |
| 42) 2,4,6-Tribromophenol    | 16.18 | 330 | 149954 | 135.53 | ng | 0.00  |
| 45) 2-Fluorobiphenyl        | 13.31 | 172 | 452926 | 77.40  | ng | -0.01 |
| 79) Terphenyl-d14           | 20.04 | 244 | 762826 | 91.89  | ng | 0.00  |

|                                |       |     |        |        |    |      |
|--------------------------------|-------|-----|--------|--------|----|------|
| Target Compounds               |       |     |        | Qvalue |    |      |
| 2) 1,4-Dioxane                 | 3.48  | 88  | 17392  | 23.869 | ng | 97   |
| 3) Pyridine                    | 3.90  | 79  | 942    | 0.470  | ng | # 71 |
| 4) n-Nitrosodimethylamine      | 3.81  | 42  | 34890  | 35.495 | ng | # 88 |
| 6) Aniline                     | 7.38  | 93  | 5138   | 1.873  | ng | 97   |
| 8) 2-Chlorophenol              | 7.61  | 128 | 73256  | 40.434 | ng | 98   |
| 9) Benzaldehyde                | 7.19  | 77  | 31553  | 22.631 | ng | 94   |
| 10) Phenol                     | 7.24  | 94  | 92806  | 40.645 | ng | 98   |
| 11) bis(2-Chloroethyl)ether    | 7.47  | 93  | 58968  | 32.437 | ng | 95   |
| 12) 1,3-Dichlorobenzene        | 7.94  | 146 | 78148  | 36.481 | ng | 95   |
| 13) 1,4-Dichlorobenzene        | 8.08  | 146 | 79667  | 36.312 | ng | 99   |
| 14) 1,2-Dichlorobenzene        | 8.40  | 146 | 76909  | 37.184 | ng | 96   |
| 15) Benzyl Alcohol             | 8.29  | 79  | 66558  | 39.675 | ng | 98   |
| 16) 2,2'-oxybis(1-Chloropropan | 8.57  | 45  | 122466 | 29.408 | ng | 96   |
| 17) 2-Methylphenol             | 8.49  | 107 | 62221  | 40.672 | ng | 98   |
| 18) Hexachloroethane           | 9.12  | 117 | 27802  | 34.679 | ng | 97   |
| 19) n-Nitroso-di-n-propylamine | 8.85  | 70  | 51245  | 33.948 | ng | 97   |
| 20) 3+4-Methylphenols          | 8.82  | 107 | 89128  | 42.186 | ng | 99   |
| 22) Acetophenone               | 8.88  | 105 | 104937 | 37.004 | ng | # 96 |
| 24) Nitrobenzene               | 9.27  | 77  | 81657  | 36.321 | ng | 99   |
| 25) Isophorone                 | 9.78  | 82  | 147139 | 36.850 | ng | # 97 |
| 26) 2-Nitrophenol              | 9.98  | 139 | 45351  | 39.413 | ng | 93   |
| 27) 2,4-Dimethylphenol         | 10.03 | 122 | 73631  | 44.649 | ng | 97   |
| 28) bis(2-Chloroethoxy)methane | 10.26 | 93  | 84557  | 37.525 | ng | 99   |
| 29) 2,4-Dichlorophenol         | 10.51 | 162 | 85924  | 46.835 | ng | 97   |
| 30) 1,2,4-Trichlorobenzene     | 10.73 | 180 | 92442  | 41.715 | ng | 95   |
| 31) Naphthalene                | 10.92 | 128 | 236063 | 42.200 | ng | 99   |
| 32) Benzoic acid               | 10.14 | 122 | 1177m  | 11.050 | ng |      |
| 33) 4-Chloroaniline            | 11.04 | 127 | 11429  | 4.706  | ng | 93   |
| 34) Hexachlorobutadiene        | 11.18 | 225 | 60409  | 40.287 | ng | 97   |
| 35) Caprolactam                | 11.82 | 113 | 9817m  | 12.930 | ng |      |
| 36) 4-Chloro-3-methylphenol    | 12.15 | 107 | 86461  | 41.298 | ng | 97   |
| 37) 2-Methylnaphthalene        | 12.52 | 142 | 182346 | 45.692 | ng | 100  |
| 38) 1-Methylnaphthalene        | 12.74 | 142 | 173922 | 44.912 | ng | 98   |
| 40) 1,2,4,5-Tetrachlorobenzene | 12.88 | 216 | 115318 | 38.428 | ng | 98   |

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 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Dec 12 15:26:27 2018  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 41) Hexachlorocyclopentadiene  | 12.84 | 237  | 120054   | 72.819 | nq    | 94       |
| 43) 2,4,6-Trichlorophenol      | 13.12 | 196  | 80161    | 43.445 | nq    | 94       |
| 44) 2,4,5-Trichlorophenol      | 13.20 | 196  | 86405    | 44.229 | nq    | 95       |
| 46) 1,1'-Biphenyl              | 13.52 | 154  | 236689   | 35.169 | nq    | 99       |
| 47) 2-Chloronaphthalene        | 13.57 | 162  | 195823   | 37.668 | nq    | 99       |
| 48) 2-Nitroaniline             | 13.78 | 65   | 59683    | 36.043 | nq    | 92       |
| 49) Acenaphthylene             | 14.42 | 152  | 312449   | 40.203 | nq    | 99       |
| 50) Dimethylphthalate          | 14.14 | 163  | 261810   | 39.934 | nq    | 99       |
| 51) 2,6-Dinitrotoluene         | 14.28 | 165  | 55483    | 37.345 | nq #  | 88       |
| 52) Acenaphthene               | 14.75 | 154  | 180945   | 38.936 | nq    | 98       |
| 53) 3-Nitroaniline             | 14.61 | 138  | 43914    | 28.996 | nq    | 97       |
| 54) 2,4-Dinitrophenol          | 14.82 | 184  | 42037    | 49.655 | nq    | 98       |
| 55) Dibenzofuran               | 15.09 | 168  | 306101   | 38.425 | nq    | 97       |
| 56) 4-Nitrophenol              | 14.92 | 139  | 81517    | 70.950 | nq    | 98       |
| 57) 2,4-Dinitrotoluene         | 15.06 | 165  | 81682    | 39.035 | nq #  | 88       |
| 58) Fluorene                   | 15.73 | 166  | 244430   | 41.415 | nq    | 98       |
| 59) 2,3,4,6-Tetrachlorophenol  | 15.32 | 232  | 77528    | 44.110 | nq    | 97       |
| 60) Diethylphthalate           | 15.49 | 149  | 255311   | 35.474 | nq    | 98       |
| 61) 4-Chlorophenyl-phenylether | 15.72 | 204  | 141609   | 38.455 | nq    | 99       |
| 62) 4-Nitroaniline             | 15.77 | 138  | 55122    | 35.353 | nq    | 97       |
| 63) Azobenzene                 | 16.01 | 77   | 205494   | 34.179 | nq    | 96       |
| 65) 4,6-Dinitro-2-methylphenol | 15.82 | 198  | 45221    | 33.848 | nq    | 91       |
| 66) n-Nitrosodiphenylamine     | 15.94 | 169  | 221518   | 40.255 | nq    | 97       |
| 67) 4-Bromophenyl-phenylether  | 16.62 | 248  | 92157    | 40.291 | nq    | 94       |
| 68) Hexachlorobenzene          | 16.73 | 284  | 98136    | 41.390 | nq    | 96       |
| 69) Atrazine                   | 16.88 | 200  | 80707    | 39.520 | nq    | 98       |
| 70) Pentachlorophenol          | 17.08 | 266  | 102590   | 73.151 | nq    | 99       |
| 71) Phenanthrene               | 17.48 | 178  | 405089   | 41.710 | nq    | 99       |
| 72) Anthracene                 | 17.57 | 178  | 409294   | 42.689 | nq    | 99       |
| 73) Carbazole                  | 17.85 | 167  | 357678   | 37.721 | nq    | 99       |
| 74) Di-n-butylphthalate        | 18.38 | 149  | 423494   | 38.923 | nq    | 99       |
| 75) Fluoranthene               | 19.49 | 202  | 484868   | 40.351 | nq    | 97       |
| 77) Benzidine                  | 19.68 | 184  | 2383m    | 0.446  | nq    |          |
| 78) Pyrene                     | 19.85 | 202  | 485127   | 40.644 | nq    | 99       |
| 80) Butylbenzylphthalate       | 20.71 | 149  | 183404   | 39.330 | nq    | 99       |
| 81) Benzo(a)anthracene         | 21.70 | 228  | 467534   | 42.467 | nq    | 99       |
| 82) 3,3'-Dichlorobenzidine     | 21.61 | 252  | 86603    | 19.834 | nq    | 99       |
| 83) Chrysene                   | 21.77 | 228  | 433234   | 40.855 | nq    | 98       |
| 84) Bis(2-ethylhexyl)phthalate | 21.56 | 149  | 252744   | 38.215 | nq    | 100      |
| 85) Di-n-octyl phthalate       | 22.79 | 149  | 432101   | 39.011 | nq    | 99       |
| 86) Indeno(1,2,3-cd)pyrene     | 28.80 | 276  | 531962   | 42.670 | nq #  | 78       |
| 88) Benzo(b)fluoranthene       | 23.96 | 252  | 453457m  | 41.898 | nq    |          |
| 89) Benzo(k)fluoranthene       | 24.03 | 252  | 455515   | 42.266 | nq    | 98       |
| 90) Benzo(a)pyrene             | 24.86 | 252  | 450774   | 43.172 | nq #  | 96       |
| 91) Dibenzo(a,h)anthracene     | 28.86 | 278  | 441167   | 42.436 | nq    | 97       |
| 92) Benzo(g,h,i)perylene       | 29.99 | 276  | 433103   | 42.273 | nq    | 96       |

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

