

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM050919\  
 Data File : BM020230.D  
 Acq On : 10 May 2019 06:36  
 Operator : JU/SJ  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTDCCC040

Quant Time: May 10 07:08:59 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM043019.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed May 01 03:58:37 2019  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 25% Max. Rel. Area : 150%

|      | Compound                     | Amount | Calc.  | %Dev   | Area% | Dev(min) |
|------|------------------------------|--------|--------|--------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4       | 20.000 | 20.000 | 0.0    | 167   | -0.02    |
| 2    | 1,4-Dioxane                  | 40.000 | 43.337 | -8.3   | 195   | -0.01    |
| 3    | Pyridine                     | 40.000 | 47.197 | -18.0  | 193   | 0.00     |
| 4    | n-Nitrosodimethylamine       | 40.000 | 40.812 | -2.0   | 179   | 0.00     |
| 5 S  | 2-Fluorophenol               | 80.000 | 91.060 | -13.8  | 198   | -0.01    |
| 6    | Aniline                      | 40.000 | 44.205 | -10.5  | 193   | -0.02    |
| 7 S  | Phenol-d6                    | 80.000 | 87.967 | -10.0  | 191   | 0.00     |
| 8    | 2-Chlorophenol               | 40.000 | 44.906 | -12.3  | 193   | -0.01    |
| 9    | Benzaldehyde                 | 40.000 | 39.992 | 0.0    | 169   | -0.01    |
| 10 C | Phenol                       | 40.000 | 44.183 | -10.5  | 189   | 0.00     |
| 11   | bis(2-Chloroethyl)ether      | 40.000 | 44.945 | -12.4  | 190   | -0.02    |
| 12   | 1,3-Dichlorobenzene          | 40.000 | 44.521 | -11.3  | 192   | -0.02    |
| 13 C | 1,4-Dichlorobenzene          | 40.000 | 44.033 | -10.1  | 190   | -0.02    |
| 14   | 1,2-Dichlorobenzene          | 40.000 | 43.686 | -9.2   | 190   | -0.02    |
| 15   | Benzyl Alcohol               | 40.000 | 39.995 | 0.0    | 170   | -0.01    |
| 16   | 2,2'-oxybis(1-Chloropropane) | 40.000 | 41.773 | -4.4   | 173   | -0.02    |
| 17   | 2-Methylphenol               | 40.000 | 43.165 | -7.9   | 184   | 0.00     |
| 18   | Hexachloroethane             | 40.000 | 39.513 | 1.2    | 173   | -0.02    |
| 19 P | n-Nitroso-di-n-propylamine   | 40.000 | 39.790 | 0.5    | 167   | -0.02    |
| 20   | 3+4-Methylphenols            | 40.000 | 42.899 | -7.2   | 182   | 0.00     |
| 21 I | Naphthalene-d8               | 20.000 | 20.000 | 0.0    | 153   | -0.02    |
| 22   | Acetophenone                 | 40.000 | 43.199 | -8.0   | 172   | -0.01    |
| 23 S | Nitrobenzene-d5              | 80.000 | 87.379 | -9.2   | 173   | -0.01    |
| 24   | Nitrobenzene                 | 40.000 | 42.709 | -6.8   | 170   | -0.01    |
| 25   | Isophorone                   | 40.000 | 43.651 | -9.1   | 169   | -0.01    |
| 26 C | 2-Nitrophenol                | 40.000 | 54.291 | -35.7# | 212   | -0.02    |
| 27   | 2,4-Dimethylphenol           | 40.000 | 43.855 | -9.6   | 175   | -0.02    |
| 28   | bis(2-Chloroethoxy)methane   | 40.000 | 44.747 | -11.9  | 174   | -0.02    |
| 29 C | 2,4-Dichlorophenol           | 40.000 | 45.816 | -14.5  | 181   | -0.01    |
| 30   | 1,2,4-Trichlorobenzene       | 40.000 | 45.631 | -14.1  | 182   | -0.02    |
| 31   | Naphthalene                  | 40.000 | 45.240 | -13.1  | 180   | -0.01    |
| 32   | Benzoic acid                 | 40.000 | 43.280 | -8.2   | 177   | 0.02     |
| 33   | 4-Chloroaniline              | 40.000 | 43.215 | -8.0   | 168   | -0.01    |
| 34 C | Hexachlorobutadiene          | 40.000 | 43.490 | -8.7   | 172   | -0.03    |
| 35   | Caprolactam                  | 40.000 | 46.581 | -16.5  | 175   | 0.02     |
| 36 C | 4-Chloro-3-methylphenol      | 40.000 | 42.369 | -5.9   | 162   | 0.00     |
| 37   | 2-Methylnaphthalene          | 40.000 | 43.896 | -9.7   | 171   | -0.02    |
| 38   | 1-Methylnaphthalene          | 40.000 | 43.858 | -9.6   | 172   | -0.02    |
| 39 I | Acenaphthene-d10             | 20.000 | 20.000 | 0.0    | 138   | -0.01    |
| 40   | 1,2,4,5-Tetrachlorobenzene   | 40.000 | 45.493 | -13.7  | 166   | -0.02    |
| 41 P | Hexachlorocyclopentadiene    | 40.000 | 10.975 | 72.6#  | 40    | -0.02    |
| 42 S | 2,4,6-Tribromophenol         | 80.000 | 89.639 | -12.0  | 156   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol        | 40.000 | 48.307 | -20.8# | 171   | -0.01    |
| 44   | 2,4,5-Trichlorophenol        | 40.000 | 45.793 | -14.5  | 164   | -0.01    |

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 Max. RRF Dev : 25% Max. Rel. Area : 150%

|      | Compound                   | Amount | Calc.  | %Dev   | Area% | Dev(min) |
|------|----------------------------|--------|--------|--------|-------|----------|
| 45 S | 2-Fluorobiphenyl           | 80.000 | 90.240 | -12.8  | 164   | -0.02    |
| 46   | 1,1'-Biphenyl              | 40.000 | 45.450 | -13.6  | 163   | -0.02    |
| 47   | 2-Chloronaphthalene        | 40.000 | 45.123 | -12.8  | 165   | -0.01    |
| 48   | 2-Nitroaniline             | 40.000 | 43.544 | -8.9   | 152   | -0.01    |
| 49   | Acenaphthylene             | 40.000 | 43.891 | -9.7   | 157   | -0.01    |
| 50   | Dimethylphthalate          | 40.000 | 42.765 | -6.9   | 150   | -0.02    |
| 51   | 2,6-Dinitrotoluene         | 40.000 | 44.179 | -10.4  | 155   | -0.01    |
| 52 C | Acenaphthene               | 40.000 | 44.159 | -10.4  | 157   | -0.01    |
| 53   | 3-Nitroaniline             | 40.000 | 44.713 | -11.8  | 157   | -0.01    |
| 54 P | 2,4-Dinitrophenol          | 40.000 | 25.017 | 37.5#  | 80    | 0.00     |
| 55   | Dibenzofuran               | 40.000 | 42.899 | -7.2   | 154   | -0.01    |
| 56 P | 4-Nitrophenol              | 40.000 | 41.005 | -2.5   | 141   | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 40.000 | 44.164 | -10.4  | 153   | 0.00     |
| 58   | Fluorene                   | 40.000 | 42.173 | -5.4   | 150   | -0.02    |
| 59   | 2,3,4,6-Tetrachlorophenol  | 40.000 | 43.598 | -9.0   | 155   | 0.00     |
| 60   | Diethylphthalate           | 40.000 | 41.345 | -3.4   | 143   | -0.02    |
| 61   | 4-Chlorophenyl-phenylether | 40.000 | 42.353 | -5.9   | 151   | -0.02    |
| 62   | 4-Nitroaniline             | 40.000 | 42.985 | -7.5   | 148   | 0.00     |
| 63   | Azobenzene                 | 40.000 | 38.426 | 3.9    | 135   | -0.02    |
| 64 I | Phenanthrene-d10           | 20.000 | 20.000 | 0.0    | 123   | -0.01    |
| 65   | 4,6-Dinitro-2-methylphenol | 40.000 | 30.922 | 22.7   | 99    | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 40.000 | 46.420 | -16.1  | 147   | -0.02    |
| 67   | 4-Bromophenyl-phenylether  | 40.000 | 45.381 | -13.5  | 145   | -0.02    |
| 68   | Hexachlorobenzene          | 40.000 | 45.127 | -12.8  | 143   | -0.01    |
| 69   | Atrazine                   | 40.000 | 42.589 | -6.5   | 133   | -0.01    |
| 70 C | Pentachlorophenol          | 40.000 | 47.768 | -19.4  | 150   | 0.00     |
| 71   | Phenanthrene               | 40.000 | 44.486 | -11.2  | 142   | -0.01    |
| 72   | Anthracene                 | 40.000 | 44.571 | -11.4  | 142   | -0.01    |
| 73   | Carbazole                  | 40.000 | 43.860 | -9.6   | 139   | 0.00     |
| 74   | Di-n-butylphthalate        | 40.000 | 43.901 | -9.8   | 136   | -0.02    |
| 75 C | Fluoranthene               | 40.000 | 43.096 | -7.7   | 137   | -0.01    |
| 76 I | Chrysene-d12               | 20.000 | 20.000 | 0.0    | 121   | -0.01    |
| 77   | Benzidine                  | 40.000 | 42.450 | -6.1   | 131   | 0.00     |
| 78   | Pyrene                     | 40.000 | 43.994 | -10.0  | 138   | -0.01    |
| 79 S | Terphenyl-d14              | 80.000 | 85.381 | -6.7   | 133   | -0.01    |
| 80   | Butylbenzylphthalate       | 40.000 | 47.233 | -18.1  | 143   | -0.02    |
| 81   | Benzo(a)anthracene         | 40.000 | 44.244 | -10.6  | 139   | -0.01    |
| 82   | 3,3'-Dichlorobenzidine     | 40.000 | 45.733 | -14.3  | 141   | 0.00     |
| 83   | Chrysene                   | 40.000 | 44.357 | -10.9  | 138   | -0.01    |
| 84   | Bis(2-ethylhexyl)phthalate | 40.000 | 43.508 | -8.8   | 130   | -0.02    |
| 85 c | Di-n-octyl phthalate       | 40.000 | 45.976 | -14.9  | 139   | -0.03    |
| 86   | Indeno(1,2,3-cd)pyrene     | 40.000 | 50.246 | -25.6# | 160   | 0.00     |
| 87 I | Perylene-d12               | 20.000 | 20.000 | 0.0    | 136   | -0.01    |

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|      | Compound               | Amount | Calc.  | %Dev  | Area% | Dev(min) |
|------|------------------------|--------|--------|-------|-------|----------|
| 88   | Benzo(b)fluoranthene   | 40.000 | 41.400 | -3.5  | 144   | -0.01    |
| 89   | Benzo(k)fluoranthene   | 40.000 | 42.770 | -6.9  | 149   | -0.01    |
| 90 C | Benzo(a)pyrene         | 40.000 | 43.496 | -8.7  | 151   | -0.01    |
| 91   | Dibenzo(a,h)anthracene | 40.000 | 45.063 | -12.7 | 160   | 0.00     |
| 92   | Benzo(q,h,i)perylene   | 40.000 | 45.573 | -13.9 | 161   | 0.00     |

(#) = Out of Range

SPCC's out = 0 CCC's out = 2