

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF071520\  
 Data File : BF120869.D  
 Acq On : 15 Jul 2020 13:05  
 Operator : JU/CG  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Jul 15 14:49:25 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070220.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jul 13 14:16:14 2020  
 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   | 134   | 0.00     |
| 2    | 1,4-Dioxane                  | 40.000 | 38.687 | 3.3   | 131   | -0.02    |
| 3    | Pyridine                     | 40.000 | 38.207 | 4.5   | 127   | -0.01    |
| 4    | n-Nitrosodimethylamine       | 40.000 | 32.064 | 19.8  | 109   | -0.01    |
| 5 S  | 2-Fluorophenol               | 80.000 | 74.264 | 7.2   | 124   | 0.00     |
| 6    | Aniline                      | 40.000 | 37.041 | 7.4   | 124   | 0.00     |
| 7 S  | Phenol-d6                    | 80.000 | 73.435 | 8.2   | 124   | 0.00     |
| 8    | 2-Chlorophenol               | 40.000 | 38.608 | 3.5   | 127   | 0.00     |
| 9    | Benzaldehyde                 | 40.000 | 35.950 | 10.1  | 133   | 0.00     |
| 10 C | Phenol                       | 40.000 | 39.772 | 0.6   | 134   | 0.00     |
| 11   | bis(2-Chloroethyl)ether      | 40.000 | 37.498 | 6.3   | 127   | 0.00     |
| 12   | 1,3-Dichlorobenzene          | 40.000 | 37.833 | 5.4   | 126   | 0.00     |
| 13 C | 1,4-Dichlorobenzene          | 40.000 | 37.460 | 6.3   | 126   | 0.00     |
| 14   | 1,2-Dichlorobenzene          | 40.000 | 36.846 | 7.9   | 123   | 0.00     |
| 15   | Benzyl Alcohol               | 40.000 | 34.473 | 13.8  | 115   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane) | 40.000 | 34.612 | 13.5  | 116   | 0.00     |
| 17   | 2-Methylphenol               | 40.000 | 37.106 | 7.2   | 125   | 0.00     |
| 18   | Hexachloroethane             | 40.000 | 37.244 | 6.9   | 124   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine   | 40.000 | 32.274 | 19.3  | 110   | 0.00     |
| 20   | 3+4-Methylphenols            | 40.000 | 34.575 | 13.6  | 114   | 0.00     |
| 21 I | Naphthalene-d8               | 20.000 | 20.000 | 0.0   | 127   | 0.00     |
| 22   | Acetophenone                 | 40.000 | 35.217 | 12.0  | 112   | 0.00     |
| 23 S | Nitrobenzene-d5              | 80.000 | 82.363 | -3.0  | 126   | 0.00     |
| 24   | Nitrobenzene                 | 40.000 | 39.441 | 1.4   | 121   | 0.00     |
| 25   | Isophorone                   | 40.000 | 36.234 | 9.4   | 116   | 0.00     |
| 26 C | 2-Nitrophenol                | 40.000 | 47.155 | -17.9 | 160   | 0.00     |
| 27   | 2,4-Dimethylphenol           | 40.000 | 39.142 | 2.1   | 124   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane   | 40.000 | 38.177 | 4.6   | 122   | 0.00     |
| 29 C | 2,4-Dichlorophenol           | 40.000 | 40.592 | -1.5  | 128   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene       | 40.000 | 40.109 | -0.3  | 126   | 0.00     |
| 31   | Naphthalene                  | 40.000 | 37.991 | 5.0   | 121   | 0.00     |
| 32   | Benzoic acid                 | 40.000 | 46.156 | -15.4 | 154   | 0.00     |
| 33   | 4-Chloroaniline              | 40.000 | 39.633 | 0.9   | 126   | 0.00     |
| 34 C | Hexachlorobutadiene          | 40.000 | 39.590 | 1.0   | 123   | 0.00     |
| 35   | Caprolactam                  | 40.000 | 42.034 | -5.1  | 133   | 0.00     |
| 36 C | 4-Chloro-3-methylphenol      | 40.000 | 38.162 | 4.6   | 119   | 0.00     |
| 37   | 2-Methylnaphthalene          | 40.000 | 38.055 | 4.9   | 119   | 0.00     |
| 38   | 1-Methylnaphthalene          | 40.000 | 38.320 | 4.2   | 120   | 0.00     |
| 39 I | Acenaphthene-d10             | 20.000 | 20.000 | 0.0   | 126   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene   | 40.000 | 41.252 | -3.1  | 125   | 0.00     |
| 41 P | Hexachlorocyclopentadiene    | 40.000 | 39.061 | 2.3   | 113   | 0.00     |
| 42 S | 2,4,6-Tribromophenol         | 80.000 | 88.348 | -10.4 | 129   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol        | 40.000 | 41.892 | -4.7  | 126   | 0.00     |
| 44   | 2,4,5-Trichlorophenol        | 40.000 | 42.057 | -5.1  | 127   | 0.00     |

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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) |
|------|----------------------------|--------|--------|--------|-------|----------|
| 45 S | 2-Fluorobiphenyl           | 80.000 | 74.544 | 6.8    | 115   | 0.00     |
| 46   | 1,1'-Biphenyl              | 40.000 | 39.177 | 2.1    | 119   | 0.00     |
| 47   | 2-Chloronaphthalene        | 40.000 | 40.066 | -0.2   | 123   | 0.00     |
| 48   | 2-Nitroaniline             | 40.000 | 41.589 | -4.0   | 133   | 0.00     |
| 49   | Acenaphthylene             | 40.000 | 38.941 | 2.6    | 122   | 0.00     |
| 50   | Dimethylphthalate          | 40.000 | 39.668 | 0.8    | 123   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 40.000 | 44.710 | -11.8  | 143   | 0.00     |
| 52 C | Acenaphthene               | 40.000 | 39.871 | 0.3    | 123   | 0.00     |
| 53   | 3-Nitroaniline             | 40.000 | 44.780 | -12.0  | 145   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 40.000 | 56.230 | -40.6# | 223   | 0.00     |
| 55   | Dibenzofuran               | 40.000 | 39.134 | 2.2    | 121   | 0.00     |
| 56 P | 4-Nitrophenol              | 40.000 | 45.326 | -13.3  | 144   | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 40.000 | 46.532 | -16.3  | 150   | 0.00     |
| 58   | Fluorene                   | 40.000 | 37.775 | 5.6    | 116   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 40.000 | 42.803 | -7.0   | 130   | 0.00     |
| 60   | Diethylphthalate           | 40.000 | 37.953 | 5.1    | 118   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 40.000 | 38.687 | 3.3    | 119   | 0.00     |
| 62   | 4-Nitroaniline             | 40.000 | 48.486 | -21.2  | 144   | 0.00     |
| 63   | Azobenzene                 | 40.000 | 35.772 | 10.6   | 113   | 0.00     |
| 64 I | Phenanthrene-d10           | 20.000 | 20.000 | 0.0    | 127   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 40.000 | 52.422 | -31.1# | 197   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 40.000 | 39.548 | 1.1    | 122   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 40.000 | 40.922 | -2.3   | 125   | 0.00     |
| 68   | Hexachlorobenzene          | 40.000 | 41.788 | -4.5   | 127   | 0.00     |
| 69   | Atrazine                   | 40.000 | 34.707 | 13.2   | 105   | 0.00     |
| 70 C | Pentachlorophenol          | 40.000 | 42.959 | -7.4   | 127   | 0.00     |
| 71   | Phenanthrene               | 40.000 | 38.313 | 4.2    | 120   | 0.00     |
| 72   | Anthracene                 | 40.000 | 38.650 | 3.4    | 120   | 0.00     |
| 73   | Carbazole                  | 40.000 | 38.869 | 2.8    | 122   | 0.00     |
| 74   | Di-n-butylphthalate        | 40.000 | 37.141 | 7.1    | 115   | 0.00     |
| 75 C | Fluoranthene               | 40.000 | 39.142 | 2.1    | 123   | 0.00     |
| 76 I | Chrysene-d12               | 20.000 | 20.000 | 0.0    | 127   | 0.00     |
| 77   | Benzidine                  | 40.000 | 36.533 | 8.7    | 107   | 0.00     |
| 78   | Pyrene                     | 40.000 | 38.822 | 2.9    | 121   | 0.00     |
| 79 S | Terphenyl-d14              | 80.000 | 71.685 | 10.4   | 114   | 0.00     |
| 80   | Butylbenzylphthalate       | 40.000 | 39.572 | 1.1    | 120   | 0.00     |
| 81   | Benzo(a)anthracene         | 40.000 | 38.677 | 3.3    | 119   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 40.000 | 43.533 | -8.8   | 130   | 0.00     |
| 83   | Chrysene                   | 40.000 | 39.504 | 1.2    | 122   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 40.000 | 36.594 | 8.5    | 113   | 0.00     |
| 85 c | Di-n-octyl phthalate       | 40.000 | 38.291 | 4.3    | 116   | 0.00     |
| 86 I | Perylene-d12               | 20.000 | 20.000 | 0.0    | 143   | 0.00     |
| 87   | Indeno(1,2,3-cd)pyrene     | 40.000 | 43.312 | -8.3   | 146   | 0.00     |

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|      | Compound               | Amount | Calc.  | %Dev  | Area% | Dev(min) |
|------|------------------------|--------|--------|-------|-------|----------|
| 88   | Benzo(b)fluoranthene   | 40.000 | 40.477 | -1.2  | 135   | 0.00     |
| 89   | Benzo(k)fluoranthene   | 40.000 | 37.066 | 7.3   | 124   | 0.00     |
| 90 C | Benzo(a)pyrene         | 40.000 | 40.046 | -0.1  | 133   | 0.00     |
| 91   | Dibenzo(a,h)anthracene | 40.000 | 43.094 | -7.7  | 144   | -0.01    |
| 92   | Benzo(q,h,i)perylene   | 40.000 | 44.541 | -11.4 | 151   | -0.01    |

(#) = Out of Range

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