

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF020923\  
 Data File : BF132372.D  
 Acq On : 09 Feb 2023 10:17  
 Operator : CG\JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Feb 10 00:56:09 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF013123.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Feb 09 01:56:26 2023  
 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    | 113   | 0.00     |
| 2    | 1,4-Dioxane                 | 40.000 | 40.748 | -1.9   | 113   | 0.02     |
| 3    | Pyridine                    | 40.000 | 36.762 | 8.1    | 101   | 0.02     |
| 4    | n-Nitrosodimethylamine      | 40.000 | 34.462 | 13.8   | 97    | 0.02     |
| 5 S  | 2-Fluorophenol              | 80.000 | 76.271 | 4.7    | 108   | 0.00     |
| 6    | Aniline                     | 40.000 | 36.318 | 9.2    | 103   | 0.00     |
| 7 S  | Phenol-d6                   | 80.000 | 73.632 | 8.0    | 104   | 0.00     |
| 8    | 2-Chlorophenol              | 40.000 | 40.092 | -0.2   | 112   | 0.00     |
| 9    | Benzaldehyde                | 40.000 | 45.093 | -12.7  | 121   | 0.00     |
| 10 C | Phenol                      | 40.000 | 36.625 | 8.4    | 103   | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 40.000 | 37.991 | 5.0    | 107   | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 40.000 | 40.756 | -1.9   | 115   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 40.000 | 40.615 | -1.5   | 114   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 40.000 | 40.643 | -1.6   | 114   | 0.00     |
| 15   | Benzyl Alcohol              | 40.000 | 36.829 | 7.9    | 102   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 40.000 | 32.038 | 19.9   | 89    | 0.00     |
| 17   | 2-Methylphenol              | 40.000 | 38.742 | 3.1    | 109   | 0.00     |
| 18   | Hexachloroethane            | 40.000 | 40.880 | -2.2   | 114   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 40.000 | 32.235 | 19.4   | 92    | 0.00     |
| 20   | 3+4-Methylphenols           | 40.000 | 37.565 | 6.1    | 103   | 0.00     |
| 21 I | Naphthalene-d8              | 20.000 | 20.000 | 0.0    | 111   | 0.00     |
| 22   | Acetophenone                | 40.000 | 37.325 | 6.7    | 105   | 0.00     |
| 23 S | Nitrobenzene-d5             | 80.000 | 73.533 | 8.1    | 102   | 0.00     |
| 24   | Nitrobenzene                | 40.000 | 35.836 | 10.4   | 99    | 0.00     |
| 25   | Isophorone                  | 40.000 | 36.454 | 8.9    | 103   | 0.00     |
| 26 C | 2-Nitrophenol               | 40.000 | 44.092 | -10.2  | 117   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 40.000 | 40.332 | -0.8   | 112   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 40.000 | 37.750 | 5.6    | 105   | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 40.000 | 42.349 | -5.9   | 117   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 40.000 | 42.914 | -7.3   | 119   | 0.00     |
| 31   | Naphthalene                 | 40.000 | 39.335 | 1.7    | 110   | 0.00     |
| 32   | Benzoic acid                | 40.000 | 60.645 | -51.6# | 163   | 0.02     |
| 33   | 4-Chloroaniline             | 40.000 | 39.084 | 2.3    | 108   | 0.00     |
| 34 C | Hexachlorobutadiene         | 40.000 | 46.699 | -16.7  | 130   | 0.00     |
| 35   | Caprolactam                 | 40.000 | 37.458 | 6.4    | 102   | 0.00     |
| 36 C | 4-Chloro-3-methylphenol     | 40.000 | 39.377 | 1.6    | 108   | 0.00     |
| 37   | 2-Methylnaphthalene         | 40.000 | 39.958 | 0.1    | 111   | 0.00     |
| 38   | 1-Methylnaphthalene         | 40.000 | 39.983 | 0.0    | 111   | 0.00     |
| 39 I | Acenaphthene-d10            | 20.000 | 20.000 | 0.0    | 109   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 40.000 | 44.131 | -10.3  | 121   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 40.000 | 40.051 | -0.1   | 106   | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 80.000 | 93.584 | -17.0  | 130   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 40.000 | 43.160 | -7.9   | 116   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 40.000 | 43.761 | -9.4   | 118   | 0.00     |
| 45 S | 2-Fluorobiphenyl            | 80.000 | 78.254 | 2.2    | 115   | 0.00     |
| 46   | 1,1'-Biphenyl               | 40.000 | 40.848 | -2.1   | 111   | 0.00     |
| 47   | 2-Chloronaphthalene         | 40.000 | 41.236 | -3.1   | 113   | 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) |
|------|----------------------------|--------|--------|-------|-------|----------|
| 48   | 2-Nitroaniline             | 40.000 | 33.995 | 15.0  | 90    | 0.00     |
| 49   | Acenaphthylene             | 40.000 | 39.706 | 0.7   | 108   | 0.00     |
| 50   | Dimethylphthalate          | 40.000 | 41.086 | -2.7  | 112   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 40.000 | 42.204 | -5.5  | 112   | 0.00     |
| 52 C | Acenaphthene               | 40.000 | 40.667 | -1.7  | 112   | 0.00     |
| 53   | 3-Nitroaniline             | 40.000 | 38.564 | 3.6   | 103   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 40.000 | 40.983 | -2.5  | 114   | 0.00     |
| 55   | Dibenzofuran               | 40.000 | 39.973 | 0.1   | 110   | 0.00     |
| 56 P | 4-Nitrophenol              | 40.000 | 36.616 | 8.5   | 96    | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 40.000 | 41.993 | -5.0  | 111   | 0.00     |
| 58   | Fluorene                   | 40.000 | 39.608 | 1.0   | 110   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 40.000 | 43.491 | -8.7  | 115   | 0.00     |
| 60   | Diethylphthalate           | 40.000 | 41.468 | -3.7  | 112   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 40.000 | 42.286 | -5.7  | 117   | 0.00     |
| 62   | 4-Nitroaniline             | 40.000 | 36.653 | 8.4   | 98    | 0.00     |
| 63   | Azobenzene                 | 40.000 | 34.238 | 14.4  | 92    | 0.00     |
| 64 I | Phenanthrene-d10           | 20.000 | 20.000 | 0.0   | 104   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 40.000 | 47.089 | -17.7 | 114   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 40.000 | 41.171 | -2.9  | 108   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 40.000 | 45.538 | -13.8 | 120   | 0.00     |
| 68   | Hexachlorobenzene          | 40.000 | 45.425 | -13.6 | 119   | 0.00     |
| 69   | Atrazine                   | 40.000 | 44.766 | -11.9 | 112   | 0.00     |
| 70 C | Pentachlorophenol          | 40.000 | 43.935 | -9.8  | 111   | 0.00     |
| 71   | Phenanthrene               | 40.000 | 39.622 | 0.9   | 105   | 0.00     |
| 72   | Anthracene                 | 40.000 | 40.210 | -0.5  | 105   | 0.00     |
| 73   | Carbazole                  | 40.000 | 37.976 | 5.1   | 100   | 0.00     |
| 74   | Di-n-butylphthalate        | 40.000 | 42.710 | -6.8  | 111   | 0.00     |
| 75 C | Fluoranthene               | 40.000 | 38.377 | 4.1   | 101   | 0.00     |
| 76 I | Chrysene-d12               | 20.000 | 20.000 | 0.0   | 86    | 0.00     |
| 77   | Benzydine                  | 40.000 | 29.225 | 26.9# | 66    | 0.00     |
| 78   | Pyrene                     | 40.000 | 45.870 | -14.7 | 100   | 0.00     |
| 79 S | Terphenyl-d14              | 80.000 | 96.531 | -20.7 | 108   | 0.00     |
| 80   | Butylbenzylphthalate       | 40.000 | 46.332 | -15.8 | 99    | 0.00     |
| 81   | Benzo(a)anthracene         | 40.000 | 41.125 | -2.8  | 88    | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 40.000 | 37.266 | 6.8   | 79    | 0.00     |
| 83   | Chrysene                   | 40.000 | 40.377 | -0.9  | 88    | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 40.000 | 44.757 | -11.9 | 96    | 0.00     |
| 85 c | Di-n-octyl phthalate       | 40.000 | 39.221 | 1.9   | 83    | 0.00     |
| 86 I | Perylene-d12               | 20.000 | 20.000 | 0.0   | 81    | 0.00     |
| 87   | Indeno(1,2,3-cd)pyrene     | 40.000 | 44.616 | -11.5 | 94    | 0.00     |
| 88   | Benzo(b)fluoranthene       | 40.000 | 41.630 | -4.1  | 81    | 0.00     |
| 89   | Benzo(k)fluoranthene       | 40.000 | 40.798 | -2.0  | 83    | 0.00     |
| 90 C | Benzo(a)pyrene             | 40.000 | 41.122 | -2.8  | 82    | 0.00     |
| 91   | Dibenzo(a,h)anthracene     | 40.000 | 44.781 | -12.0 | 93    | 0.00     |
| 92   | Benzo(g,h,i)perylene       | 40.000 | 44.586 | -11.5 | 95    | 0.00     |

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| Compound | Amount | Calc. | %Dev | Area% | Dev(min) |
|----------|--------|-------|------|-------|----------|
|----------|--------|-------|------|-------|----------|

(#) = Out of Range

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