

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM021224\  
 Data File : BM044066.D  
 Acq On : 12 Feb 2024 21:21  
 Operator : MA/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTDCCC040

Quant Time: Feb 13 00:50:32 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM020824.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Feb 09 02:42:01 2024  
 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                    | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|-----------------------------|-------|-------|------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0  | 125   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.546 | 0.549 | -0.5 | 131   | 0.00     |
| 3    | Pyridine                    | 1.441 | 1.458 | -1.2 | 129   | 0.00     |
| 4    | n-Nitrosodimethylamine      | 0.547 | 0.564 | -3.1 | 131   | 0.00     |
| 5 S  | 2-Fluorophenol              | 1.375 | 1.365 | 0.7  | 129   | 0.00     |
| 6    | Aniline                     | 1.720 | 1.709 | 0.6  | 125   | 0.00     |
| 7 S  | Phenol-d6                   | 1.750 | 1.763 | -0.7 | 129   | 0.00     |
| 8    | 2-Chlorophenol              | 1.436 | 1.398 | 2.6  | 126   | 0.00     |
| 9    | Benzaldehyde                | 0.899 | 0.921 | -2.4 | 132   | 0.00     |
| 10 C | Phenol                      | 1.766 | 1.775 | -0.5 | 129   | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 1.462 | 1.452 | 0.7  | 128   | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 1.630 | 1.586 | 2.7  | 126   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.643 | 1.596 | 2.9  | 126   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 1.566 | 1.523 | 2.7  | 126   | 0.00     |
| 15   | Benzyl Alcohol              | 1.139 | 1.135 | 0.4  | 126   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 1.854 | 1.807 | 2.5  | 126   | 0.00     |
| 17   | 2-Methylphenol              | 1.159 | 1.153 | 0.5  | 127   | 0.00     |
| 18   | Hexachloroethane            | 0.615 | 0.598 | 2.8  | 126   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 1.074 | 1.065 | 0.8  | 127   | 0.00     |
| 20   | 3+4-Methylphenols           | 1.595 | 1.559 | 2.3  | 124   | 0.00     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0  | 124   | 0.00     |
| 22   | Acetophenone                | 0.508 | 0.510 | -0.4 | 127   | 0.00     |
| 23 S | Nitrobenzene-d5             | 0.377 | 0.384 | -1.9 | 128   | 0.00     |
| 24   | Nitrobenzene                | 0.386 | 0.392 | -1.6 | 127   | 0.00     |
| 25   | Isophorone                  | 0.726 | 0.723 | 0.4  | 124   | 0.00     |
| 26 C | 2-Nitrophenol               | 0.179 | 0.178 | 0.6  | 123   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 0.229 | 0.227 | 0.9  | 123   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 0.469 | 0.459 | 2.1  | 123   | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.301 | 0.298 | 1.0  | 123   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 0.339 | 0.330 | 2.7  | 123   | 0.00     |
| 31   | Naphthalene                 | 1.104 | 1.083 | 1.9  | 125   | 0.00     |
| 32   | Benzoic acid                | 0.230 | 0.217 | 5.7  | 115   | 0.00     |
| 33   | 4-Chloroaniline             | 0.418 | 0.409 | 2.2  | 123   | 0.00     |
| 34 C | Hexachlorobutadiene         | 0.194 | 0.188 | 3.1  | 123   | 0.00     |
| 35   | Caprolactam                 | 0.097 | 0.095 | 2.1  | 122   | 0.00     |
| 36 C | 4-Chloro-3-methylphenol     | 0.326 | 0.324 | 0.6  | 123   | 0.00     |
| 37   | 2-Methylnaphthalene         | 0.735 | 0.711 | 3.3  | 122   | 0.00     |
| 38   | 1-Methylnaphthalene         | 0.722 | 0.690 | 4.4  | 122   | 0.00     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0  | 120   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.581 | 0.579 | 0.3  | 122   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 0.254 | 0.253 | 0.4  | 118   | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 0.217 | 0.212 | 2.3  | 116   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 0.396 | 0.399 | -0.8 | 120   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 0.434 | 0.441 | -1.6 | 121   | 0.00     |
| 45 S | 2-Fluorobiphenyl            | 1.417 | 1.421 | -0.3 | 123   | 0.00     |
| 46   | 1,1'-Biphenyl               | 1.578 | 1.581 | -0.2 | 122   | 0.00     |
| 47   | 2-Chloronaphthalene         | 1.250 | 1.251 | -0.1 | 122   | 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                   | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|----------------------------|-------|-------|------|-------|----------|
| 48   | 2-Nitroaniline             | 0.352 | 0.371 | -5.4 | 126   | 0.00     |
| 49   | Acenaphthylene             | 1.865 | 1.861 | 0.2  | 121   | 0.00     |
| 50   | Dimethylphthalate          | 1.483 | 1.450 | 2.2  | 119   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 0.327 | 0.324 | 0.9  | 118   | 0.00     |
| 52 C | Acenaphthene               | 1.154 | 1.139 | 1.3  | 121   | 0.00     |
| 53   | 3-Nitroaniline             | 0.352 | 0.358 | -1.7 | 120   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 0.170 | 0.165 | 2.9  | 118   | 0.00     |
| 55   | Dibenzofuran               | 1.787 | 1.761 | 1.5  | 121   | 0.00     |
| 56 P | 4-Nitrophenol              | 0.286 | 0.290 | -1.4 | 119   | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 0.428 | 0.432 | -0.9 | 119   | 0.00     |
| 58   | Fluorene                   | 1.381 | 1.365 | 1.2  | 122   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.341 | 0.339 | 0.6  | 119   | 0.00     |
| 60   | Diethylphthalate           | 1.533 | 1.470 | 4.1  | 117   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 0.669 | 0.658 | 1.6  | 121   | 0.00     |
| 62   | 4-Nitroaniline             | 0.354 | 0.360 | -1.7 | 121   | 0.00     |
| 63   | Azobenzene                 | 1.438 | 1.468 | -2.1 | 124   | 0.00     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0  | 118   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 0.119 | 0.117 | 1.7  | 119   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 0.608 | 0.599 | 1.5  | 119   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 0.198 | 0.183 | 7.6  | 112   | 0.00     |
| 68   | Hexachlorobenzene          | 0.242 | 0.234 | 3.3  | 117   | 0.00     |
| 69   | Atrazine                   | 0.165 | 0.162 | 1.8  | 112   | 0.00     |
| 70 C | Pentachlorophenol          | 0.158 | 0.156 | 1.3  | 114   | 0.00     |
| 71   | Phenanthrene               | 1.056 | 1.036 | 1.9  | 119   | 0.00     |
| 72   | Anthracene                 | 1.046 | 1.033 | 1.2  | 119   | 0.00     |
| 73   | Carbazole                  | 1.017 | 1.005 | 1.2  | 119   | 0.00     |
| 74   | Di-n-butylphthalate        | 1.322 | 1.275 | 3.6  | 114   | 0.00     |
| 75 C | Fluoranthene               | 1.186 | 1.162 | 2.0  | 119   | 0.00     |
| 76 I | Chrysene-d12               | 1.000 | 1.000 | 0.0  | 117   | 0.00     |
| 77   | Benzydine                  | 0.311 | 0.284 | 8.7  | 80    | 0.00     |
| 78   | Pyrene                     | 1.540 | 1.534 | 0.4  | 118   | 0.00     |
| 79 S | Terphenyl-d14              | 1.141 | 1.134 | 0.6  | 118   | 0.00     |
| 80   | Butylbenzylphthalate       | 0.685 | 0.663 | 3.2  | 111   | 0.00     |
| 81   | Benzo(a)anthracene         | 1.391 | 1.369 | 1.6  | 117   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 0.456 | 0.444 | 2.6  | 111   | 0.00     |
| 83   | Chrysene                   | 1.314 | 1.303 | 0.8  | 118   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 1.010 | 0.981 | 2.9  | 111   | 0.00     |
| 85 c | Di-n-octyl phthalate       | 1.764 | 1.645 | 6.7  | 107   | 0.00     |
| 86 I | Perylene-d12               | 1.000 | 1.000 | 0.0  | 116   | 0.00     |
| 87   | Indeno(1,2,3-cd)pyrene     | 1.416 | 1.394 | 1.6  | 115   | 0.00     |
| 88   | Benzo(b)fluoranthene       | 1.250 | 1.246 | 0.3  | 118   | 0.00     |
| 89   | Benzo(k)fluoranthene       | 1.188 | 1.164 | 2.0  | 115   | 0.00     |
| 90 C | Benzo(a)pyrene             | 1.092 | 1.074 | 1.6  | 115   | 0.00     |
| 91   | Dibenzo(a,h)anthracene     | 1.184 | 1.165 | 1.6  | 115   | 0.00     |
| 92   | Benzo(g,h,i)perylene       | 1.201 | 1.174 | 2.2  | 114   | 0.00     |

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| Compound | AvgRF | CCRF | %Dev Area | % Dev(min) |
|----------|-------|------|-----------|------------|
|----------|-------|------|-----------|------------|

(#) = Out of Range

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