

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM102523\  
 Data File : BM042433.D  
 Acq On : 25 Oct 2023 12:06  
 Operator : MA/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTDCCC040

Quant Time: Oct 25 12:46:24 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM102123.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sat Oct 21 02:32:53 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                    | AvgRF | CCRF  | %Dev   | Area% | Dev(min) |
|------|-----------------------------|-------|-------|--------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0    | 103   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.666 | 0.557 | 16.4   | 88    | 0.00     |
| 3    | Pyridine                    | 1.955 | 1.616 | 17.3   | 86    | 0.00     |
| 4    | n-Nitrosodimethylamine      | 0.987 | 0.790 | 20.0   | 81    | 0.00     |
| 5 S  | 2-Fluorophenol              | 1.415 | 1.282 | 9.4    | 92    | 0.00     |
| 6    | Aniline                     | 2.415 | 2.202 | 8.8    | 92    | 0.00     |
| 7 S  | Phenol-d6                   | 1.896 | 1.711 | 9.8    | 92    | 0.00     |
| 8    | 2-Chlorophenol              | 1.404 | 1.347 | 4.1    | 97    | 0.00     |
| 9    | Benzaldehyde                | 1.105 | 0.942 | 14.8   | 90    | 0.00     |
| 10 C | Phenol                      | 1.960 | 1.742 | 11.1   | 90    | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 1.609 | 1.418 | 11.9   | 91    | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 1.452 | 1.433 | 1.3    | 102   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.462 | 1.452 | 0.7    | 102   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 1.404 | 1.403 | 0.1    | 103   | 0.00     |
| 15   | Benzyl Alcohol              | 1.406 | 1.302 | 7.4    | 93    | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 2.546 | 2.249 | 11.7   | 91    | 0.00     |
| 17   | 2-Methylphenol              | 1.293 | 1.257 | 2.8    | 97    | 0.00     |
| 18   | Hexachloroethane            | 0.563 | 0.564 | -0.2   | 102   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 1.285 | 1.252 | 2.6    | 95    | 0.00     |
| 20   | 3+4-Methylphenols           | 1.734 | 1.701 | 1.9    | 99    | 0.00     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 105   | 0.00     |
| 22   | Acetophenone                | 0.546 | 0.526 | 3.7    | 97    | 0.00     |
| 23 S | Nitrobenzene-d5             | 0.452 | 0.431 | 4.6    | 94    | 0.00     |
| 24   | Nitrobenzene                | 0.452 | 0.428 | 5.3    | 94    | 0.00     |
| 25   | Isophorone                  | 0.769 | 0.805 | -4.7   | 103   | 0.00     |
| 26 C | 2-Nitrophenol               | 0.154 | 0.185 | -20.1# | 112   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 0.312 | 0.327 | -4.8   | 104   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 0.487 | 0.472 | 3.1    | 97    | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.271 | 0.309 | -14.0  | 113   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 0.291 | 0.323 | -11.0  | 113   | -0.01    |
| 31   | Naphthalene                 | 1.027 | 1.060 | -3.2   | 104   | 0.00     |
| 32   | Benzoic acid                | 0.201 | 0.267 | -32.8# | 126   | 0.00     |
| 33   | 4-Chloroaniline             | 0.446 | 0.462 | -3.6   | 103   | -0.01    |
| 34 C | Hexachlorobutadiene         | 0.157 | 0.196 | -24.8# | 126   | 0.00     |
| 35   | Caprolactam                 | 0.098 | 0.111 | -13.3  | 113   | 0.00     |
| 36 C | 4-Chloro-3-methylphenol     | 0.327 | 0.358 | -9.5   | 108   | 0.00     |
| 37   | 2-Methylnaphthalene         | 0.692 | 0.752 | -8.7   | 109   | 0.00     |
| 38   | 1-Methylnaphthalene         | 0.648 | 0.708 | -9.3   | 109   | 0.00     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 122   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.578 | 0.613 | -6.1   | 123   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 0.271 | 0.313 | -15.5  | 129   | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 0.193 | 0.258 | -33.7# | 152#  | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 0.374 | 0.415 | -11.0  | 125   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 0.442 | 0.485 | -9.7   | 125   | 0.00     |
| 45 S | 2-Fluorobiphenyl            | 1.485 | 1.468 | 1.1    | 114   | 0.00     |
| 46   | 1,1'-Biphenyl               | 1.587 | 1.549 | 2.4    | 113   | 0.00     |
| 47   | 2-Chloronaphthalene         | 1.164 | 1.133 | 2.7    | 113   | 0.00     |

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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sat Oct 21 02:32:53 2023  
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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             | 0.452 | 0.436 | 3.5    | 104   | 0.00     |
| 49   | Acenaphthylene             | 1.879 | 1.912 | -1.8   | 115   | 0.00     |
| 50   | Dimethylphthalate          | 1.467 | 1.536 | -4.7   | 121   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 0.299 | 0.330 | -10.4  | 121   | 0.00     |
| 52 C | Acenaphthene               | 1.136 | 1.139 | -0.3   | 116   | 0.00     |
| 53   | 3-Nitroaniline             | 0.348 | 0.373 | -7.2   | 116   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 0.165 | 0.204 | -23.6  | 141   | 0.00     |
| 55   | Dibenzofuran               | 1.797 | 1.845 | -2.7   | 119   | 0.00     |
| 56 P | 4-Nitrophenol              | 0.299 | 0.296 | 1.0    | 113   | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 0.383 | 0.449 | -17.2  | 128   | 0.00     |
| 58   | Fluorene                   | 1.418 | 1.470 | -3.7   | 119   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.332 | 0.406 | -22.3  | 138   | 0.00     |
| 60   | Diethylphthalate           | 1.448 | 1.550 | -7.0   | 122   | -0.01    |
| 61   | 4-Chlorophenyl-phenylether | 0.668 | 0.758 | -13.5  | 131   | 0.00     |
| 62   | 4-Nitroaniline             | 0.326 | 0.362 | -11.0  | 118   | 0.00     |
| 63   | Azobenzene                 | 1.737 | 1.669 | 3.9    | 105   | 0.00     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0    | 135   | -0.01    |
| 65   | 4,6-Dinitro-2-methylphenol | 0.115 | 0.130 | -13.0  | 140   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 0.602 | 0.604 | -0.3   | 125   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 0.185 | 0.213 | -15.1  | 143   | 0.00     |
| 68   | Hexachlorobenzene          | 0.198 | 0.231 | -16.7  | 147   | 0.00     |
| 69   | Atrazine                   | 0.153 | 0.183 | -19.6  | 147   | 0.00     |
| 70 C | Pentachlorophenol          | 0.142 | 0.177 | -24.6# | 156#  | 0.00     |
| 71   | Phenanthrene               | 1.088 | 1.088 | 0.0    | 126   | 0.00     |
| 72   | Anthracene                 | 1.078 | 1.097 | -1.8   | 125   | -0.01    |
| 73   | Carbazole                  | 1.011 | 1.010 | 0.1    | 123   | 0.00     |
| 74   | Di-n-butylphthalate        | 1.157 | 1.330 | -15.0  | 137   | -0.01    |
| 75 C | Fluoranthene               | 1.198 | 1.309 | -9.3   | 136   | 0.00     |
| 76 I | Chrysene-d12               | 1.000 | 1.000 | 0.0    | 126   | -0.01    |
| 77   | Benzidine                  | 0.289 | 0.356 | -23.2  | 178#  | 0.00     |
| 78   | Pyrene                     | 1.398 | 1.555 | -11.2  | 133   | 0.00     |
| 79 S | Terphenyl-d14              | 1.128 | 1.287 | -14.1  | 132   | -0.01    |
| 80   | Butylbenzylphthalate       | 0.564 | 0.708 | -25.5# | 142   | -0.01    |
| 81   | Benzo(a)anthracene         | 1.341 | 1.376 | -2.6   | 122   | -0.01    |
| 82   | 3,3'-Dichlorobenzidine     | 0.400 | 0.507 | -26.7# | 148   | -0.01    |
| 83   | Chrysene                   | 1.288 | 1.291 | -0.2   | 120   | -0.01    |
| 84   | Bis(2-ethylhexyl)phthalate | 0.795 | 1.033 | -29.9# | 144   | -0.01    |
| 85 c | Di-n-octyl phthalate       | 1.441 | 1.720 | -19.4  | 138   | -0.02    |
| 86 I | Perylene-d12               | 1.000 | 1.000 | 0.0    | 133   | -0.01    |
| 87   | Indeno(1,2,3-cd)pyrene     | 1.462 | 1.347 | 7.9    | 116   | -0.02    |
| 88   | Benzo(b)fluoranthene       | 1.204 | 1.191 | 1.1    | 122   | -0.01    |
| 89   | Benzo(k)fluoranthene       | 1.250 | 1.174 | 6.1    | 118   | -0.01    |
| 90 C | Benzo(a)pyrene             | 1.159 | 1.139 | 1.7    | 121   | -0.02    |
| 91   | Dibenzo(a,h)anthracene     | 1.215 | 1.127 | 7.2    | 116   | -0.02    |
| 92   | Benzo(g,h,i)perylene       | 1.180 | 1.106 | 6.3    | 119   | -0.03    |

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|----------|-------|------|------|-------|----------|
|----------|-------|------|------|-------|----------|

(#) = Out of Range

SPCC's out = 0 CCC's out = 3