

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM041425\  
 Data File : BM049910.D  
 Acq On : 14 Apr 2025 12:29  
 Operator : RC/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTDCCC040

Quant Time: Apr 14 13:34:25 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM040825.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Apr 09 04:00:55 2025  
 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    | 124   | -0.01    |
| 2    | 1,4-Dioxane                 | 0.485 | 0.471 | 2.9    | 120   | 0.00     |
| 3    | Pyridine                    | 1.274 | 1.232 | 3.3    | 119   | 0.00     |
| 4    | n-Nitrosodimethylamine      | 0.485 | 0.484 | 0.2    | 123   | 0.00     |
| 5 S  | 2-Fluorophenol              | 1.178 | 1.160 | 1.5    | 121   | 0.00     |
| 6    | Aniline                     | 1.256 | 0.612 | 51.3#  | 57    | 0.00     |
| 7 S  | Phenol-d6                   | 1.465 | 1.464 | 0.1    | 122   | 0.00     |
| 8    | 2-Chlorophenol              | 1.207 | 1.202 | 0.4    | 122   | 0.00     |
| 9    | Benzaldehyde                | 0.752 | 0.624 | 17.0   | 103   | 0.00     |
| 10 C | Phenol                      | 1.430 | 1.422 | 0.6    | 122   | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 1.121 | 1.630 | -45.4# | 178#  | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 1.427 | 1.435 | -0.6   | 124   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.436 | 1.433 | 0.2    | 122   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 1.352 | 1.353 | -0.1   | 122   | 0.00     |
| 15   | Benzyl Alcohol              | 0.923 | 0.857 | 7.2    | 111   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 1.250 | 1.278 | -2.2   | 125   | -0.01    |
| 17   | 2-Methylphenol              | 0.881 | 0.884 | -0.3   | 123   | 0.00     |
| 18   | Hexachloroethane            | 0.500 | 0.502 | -0.4   | 124   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 0.812 | 0.846 | -4.2   | 125   | 0.00     |
| 20   | 3+4-Methylphenols           | 1.201 | 1.193 | 0.7    | 121   | 0.00     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 123   | 0.00     |
| 22   | Acetophenone                | 0.486 | 0.497 | -2.3   | 123   | 0.00     |
| 23 S | Nitrobenzene-d5             | 0.359 | 0.368 | -2.5   | 124   | 0.00     |
| 24   | Nitrobenzene                | 0.346 | 0.348 | -0.6   | 121   | 0.00     |
| 25   | Isophorone                  | 0.602 | 0.621 | -3.2   | 124   | 0.00     |
| 26 C | 2-Nitrophenol               | 0.183 | 0.191 | -4.4   | 124   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 0.208 | 0.214 | -2.9   | 123   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 0.403 | 0.418 | -3.7   | 126   | -0.01    |
| 29 C | 2,4-Dichlorophenol          | 0.345 | 0.353 | -2.3   | 123   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 0.399 | 0.410 | -2.8   | 125   | 0.00     |
| 31   | Naphthalene                 | 1.028 | 1.042 | -1.4   | 123   | 0.00     |
| 32   | Benzoic acid                | 0.220 | 0.209 | 5.0    | 112   | 0.02     |
| 33   | 4-Chloroaniline             | 0.363 | 0.343 | 5.5    | 111   | 0.03     |
| 34 C | Hexachlorobutadiene         | 0.247 | 0.257 | -4.0   | 128   | -0.01    |
| 35   | Caprolactam                 | 0.099 | 0.096 | 3.0    | 116   | 0.00     |
| 36 C | 4-Chloro-3-methylphenol     | 0.302 | 0.308 | -2.0   | 123   | 0.00     |
| 37   | 2-Methylnaphthalene         | 0.724 | 0.749 | -3.5   | 126   | 0.00     |
| 38   | 1-Methylnaphthalene         | 0.705 | 0.722 | -2.4   | 125   | 0.00     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 126   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.700 | 0.717 | -2.4   | 128   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 0.245 | 0.268 | -9.4   | 128   | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 0.291 | 0.307 | -5.5   | 132   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 0.445 | 0.462 | -3.8   | 127   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 0.484 | 0.488 | -0.8   | 123   | 0.00     |
| 45 S | 2-Fluorobiphenyl            | 1.475 | 1.522 | -3.2   | 127   | -0.01    |
| 46   | 1,1'-Biphenyl               | 1.487 | 1.528 | -2.8   | 126   | 0.00     |
| 47   | 2-Chloronaphthalene         | 1.169 | 1.180 | -0.9   | 125   | 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             | 0.270 | 0.279 | -3.3  | 124   | 0.00     |
| 49   | Acenaphthylene             | 1.702 | 1.738 | -2.1  | 126   | 0.00     |
| 50   | Dimethylphthalate          | 1.411 | 1.459 | -3.4  | 128   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 0.296 | 0.312 | -5.4  | 127   | 0.00     |
| 52 C | Acenaphthene               | 1.083 | 1.100 | -1.6  | 125   | 0.00     |
| 53   | 3-Nitroaniline             | 0.286 | 0.288 | -0.7  | 118   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 0.175 | 0.186 | -6.3  | 128   | 0.00     |
| 55   | Dibenzofuran               | 1.810 | 1.851 | -2.3  | 127   | 0.00     |
| 56 P | 4-Nitrophenol              | 0.255 | 0.249 | 2.4   | 116   | 0.01     |
| 57   | 2,4-Dinitrotoluene         | 0.394 | 0.425 | -7.9  | 128   | 0.00     |
| 58   | Fluorene                   | 1.439 | 1.481 | -2.9  | 126   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.424 | 0.426 | -0.5  | 126   | 0.00     |
| 60   | Diethylphthalate           | 1.353 | 1.415 | -4.6  | 129   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 0.771 | 0.810 | -5.1  | 130   | -0.01    |
| 62   | 4-Nitroaniline             | 0.295 | 0.296 | -0.3  | 117   | 0.00     |
| 63   | Azobenzene                 | 1.158 | 1.198 | -3.5  | 126   | 0.00     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 129   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 0.126 | 0.130 | -3.2  | 127   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 0.599 | 0.609 | -1.7  | 126   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 0.232 | 0.243 | -4.7  | 131   | 0.00     |
| 68   | Hexachlorobenzene          | 0.266 | 0.275 | -3.4  | 130   | -0.01    |
| 69   | Atrazine                   | 0.155 | 0.133 | 14.2  | 132   | 0.00     |
| 70 C | Pentachlorophenol          | 0.196 | 0.190 | 3.1   | 119   | 0.00     |
| 71   | Phenanthrene               | 1.096 | 1.125 | -2.6  | 128   | 0.00     |
| 72   | Anthracene                 | 1.081 | 1.103 | -2.0  | 126   | 0.00     |
| 73   | Carbazole                  | 1.018 | 1.028 | -1.0  | 125   | 0.00     |
| 74   | Di-n-butylphthalate        | 1.172 | 1.251 | -6.7  | 132   | 0.00     |
| 75 C | Fluoranthene               | 1.348 | 1.409 | -4.5  | 131   | 0.00     |
| 76 I | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 128   | 0.00     |
| 77   | Benzidine                  | 0.265 | 0.182 | 31.3# | 58    | 0.00     |
| 78   | Pyrene                     | 1.378 | 1.441 | -4.6  | 131   | 0.00     |
| 79 S | Terphenyl-d14              | 1.067 | 1.221 | -14.4 | 132   | 0.00     |
| 80   | Butylbenzylphthalate       | 0.518 | 0.551 | -6.4  | 131   | 0.00     |
| 81   | Benzo(a)anthracene         | 1.329 | 1.370 | -3.1  | 129   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 0.502 | 0.500 | 0.4   | 122   | 0.00     |
| 83   | Chrysene                   | 1.264 | 1.289 | -2.0  | 128   | -0.01    |
| 84   | Bis(2-ethylhexyl)phthalate | 0.755 | 0.817 | -8.2  | 133   | 0.00     |
| 85 c | Di-n-octyl phthalate       | 1.320 | 1.389 | -5.2  | 130   | -0.01    |
| 86 I | Perylene-d12               | 1.000 | 1.000 | 0.0   | 125   | -0.02    |
| 87   | Indeno(1,2,3-cd)pyrene     | 1.491 | 1.476 | 1.0   | 122   | -0.02    |
| 88   | Benzo(b)fluoranthene       | 1.277 | 1.307 | -2.3  | 125   | 0.00     |
| 89   | Benzo(k)fluoranthene       | 1.238 | 1.257 | -1.5  | 126   | -0.01    |
| 90 C | Benzo(a)pyrene             | 1.122 | 1.136 | -1.2  | 126   | -0.01    |
| 91   | Dibenzo(a,h)anthracene     | 1.228 | 1.223 | 0.4   | 123   | -0.02    |
| 92   | Benzo(g,h,i)perylene       | 1.255 | 1.234 | 1.7   | 122   | -0.02    |

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

(#) = Out of Range

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