

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM043025\  
 Data File : BM050052.D  
 Acq On : 30 Apr 2025 19:20  
 Operator : RC/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTDCCC040

Quant Time: May 01 02:18:38 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM042825.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Apr 28 18:09:16 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   | 102   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.490 | 0.520 | -6.1  | 104   | 0.00     |
| 3    | Pyridine                    | 1.258 | 1.382 | -9.9  | 106   | 0.00     |
| 4    | n-Nitrosodimethylamine      | 0.493 | 0.542 | -9.9  | 106   | 0.00     |
| 5 S  | 2-Fluorophenol              | 1.142 | 1.228 | -7.5  | 104   | 0.00     |
| 6    | Aniline                     | 1.813 | 1.955 | -7.8  | 103   | 0.00     |
| 7 S  | Phenol-d6                   | 1.436 | 1.583 | -10.2 | 105   | 0.00     |
| 8    | 2-Chlorophenol              | 1.235 | 1.317 | -6.6  | 103   | 0.00     |
| 9    | Benzaldehyde                | 0.961 | 1.046 | -8.8  | 104   | 0.00     |
| 10 C | Phenol                      | 1.453 | 1.586 | -9.2  | 105   | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 1.213 | 1.311 | -8.1  | 105   | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 1.492 | 1.557 | -4.4  | 102   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.499 | 1.575 | -5.1  | 103   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 1.448 | 1.520 | -5.0  | 102   | 0.00     |
| 15   | Benzyl Alcohol              | 0.996 | 1.107 | -11.1 | 105   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 1.475 | 1.623 | -10.0 | 107   | 0.00     |
| 17   | 2-Methylphenol              | 0.957 | 1.041 | -8.8  | 103   | 0.00     |
| 18   | Hexachloroethane            | 0.533 | 0.563 | -5.6  | 104   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 0.921 | 1.034 | -12.3 | 105   | 0.00     |
| 20   | 3+4-Methylphenols           | 1.293 | 1.418 | -9.7  | 104   | 0.00     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 103   | 0.00     |
| 22   | Acetophenone                | 0.497 | 0.539 | -8.5  | 105   | 0.00     |
| 23 S | Nitrobenzene-d5             | 0.406 | 0.445 | -9.6  | 106   | 0.00     |
| 24   | Nitrobenzene                | 0.351 | 0.381 | -8.5  | 106   | 0.00     |
| 25   | Isophorone                  | 0.692 | 0.744 | -7.5  | 106   | 0.00     |
| 26 C | 2-Nitrophenol               | 0.179 | 0.195 | -8.9  | 103   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 0.297 | 0.325 | -9.4  | 104   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 0.428 | 0.461 | -7.7  | 105   | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.333 | 0.361 | -8.4  | 103   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 0.406 | 0.423 | -4.2  | 102   | 0.00     |
| 31   | Naphthalene                 | 1.013 | 1.068 | -5.4  | 104   | 0.00     |
| 32   | Benzoic acid                | 0.190 | 0.198 | -4.2  | 104   | 0.00     |
| 33   | 4-Chloroaniline             | 0.415 | 0.447 | -7.7  | 105   | 0.00     |
| 34 C | Hexachlorobutadiene         | 0.258 | 0.271 | -5.0  | 104   | 0.00     |
| 35   | Caprolactam                 | 0.097 | 0.103 | -6.2  | 102   | 0.00     |
| 36 C | 4-Chloro-3-methylphenol     | 0.308 | 0.329 | -6.8  | 105   | 0.00     |
| 37   | 2-Methylnaphthalene         | 0.679 | 0.717 | -5.6  | 104   | 0.00     |
| 38   | 1-Methylnaphthalene         | 0.719 | 0.755 | -5.0  | 104   | 0.00     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 103   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.692 | 0.754 | -9.0  | 104   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 0.423 | 0.448 | -5.9  | 101   | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 0.296 | 0.324 | -9.5  | 105   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 0.439 | 0.483 | -10.0 | 104   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 0.474 | 0.523 | -10.3 | 104   | 0.00     |
| 45 S | 2-Fluorobiphenyl            | 1.691 | 1.873 | -10.8 | 107   | 0.00     |
| 46   | 1,1'-Biphenyl               | 1.487 | 1.591 | -7.0  | 104   | 0.00     |
| 47   | 2-Chloronaphthalene         | 1.177 | 1.259 | -7.0  | 104   | 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             | 0.276 | 0.313 | -13.4 | 106   | 0.00     |
| 49   | Acenaphthylene             | 1.857 | 1.980 | -6.6  | 103   | 0.00     |
| 50   | Dimethylphthalate          | 1.462 | 1.544 | -5.6  | 104   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 0.303 | 0.328 | -8.3  | 103   | 0.00     |
| 52 C | Acenaphthene               | 1.075 | 1.153 | -7.3  | 105   | 0.00     |
| 53   | 3-Nitroaniline             | 0.289 | 0.313 | -8.3  | 100   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 0.164 | 0.166 | -1.2  | 96    | 0.00     |
| 55   | Dibenzofuran               | 1.788 | 1.908 | -6.7  | 105   | 0.00     |
| 56 P | 4-Nitrophenol              | 0.205 | 0.226 | -10.2 | 104   | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 0.418 | 0.462 | -10.5 | 104   | 0.00     |
| 58   | Fluorene                   | 1.463 | 1.608 | -9.9  | 106   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.410 | 0.447 | -9.0  | 106   | 0.00     |
| 60   | Diethylphthalate           | 1.377 | 1.463 | -6.2  | 105   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 0.804 | 0.892 | -10.9 | 107   | 0.00     |
| 62   | 4-Nitroaniline             | 0.279 | 0.313 | -12.2 | 103   | 0.00     |
| 63   | Azobenzene                 | 1.154 | 1.286 | -11.4 | 109   | 0.00     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 104   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 0.125 | 0.129 | -3.2  | 99    | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 0.614 | 0.663 | -8.0  | 105   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 0.242 | 0.259 | -7.0  | 104   | 0.00     |
| 68   | Hexachlorobenzene          | 0.279 | 0.295 | -5.7  | 104   | 0.00     |
| 69   | Atrazine                   | 0.222 | 0.240 | -8.1  | 105   | 0.00     |
| 70 C | Pentachlorophenol          | 0.157 | 0.171 | -8.9  | 106   | 0.00     |
| 71   | Phenanthrene               | 1.128 | 1.217 | -7.9  | 106   | 0.00     |
| 72   | Anthracene                 | 1.142 | 1.237 | -8.3  | 106   | 0.00     |
| 73   | Carbazole                  | 0.991 | 1.064 | -7.4  | 105   | 0.00     |
| 74   | Di-n-butylphthalate        | 1.180 | 1.270 | -7.6  | 106   | 0.00     |
| 75 C | Fluoranthene               | 1.324 | 1.445 | -9.1  | 106   | 0.00     |
| 76 I | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 109   | 0.00     |
| 77   | Benzidine                  | 0.710 | 0.670 | 5.6   | 94    | 0.00     |
| 78   | Pyrene                     | 1.404 | 1.489 | -6.1  | 107   | 0.00     |
| 79 S | Terphenyl-d14              | 1.352 | 1.553 | -14.9 | 111   | 0.00     |
| 80   | Butylbenzylphthalate       | 0.526 | 0.546 | -3.8  | 106   | 0.00     |
| 81   | Benzo(a)anthracene         | 1.377 | 1.478 | -7.3  | 109   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 0.527 | 0.557 | -5.7  | 104   | 0.00     |
| 83   | Chrysene                   | 1.275 | 1.356 | -6.4  | 110   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 0.786 | 0.841 | -7.0  | 108   | 0.00     |
| 85 c | Di-n-octyl phthalate       | 1.296 | 1.329 | -2.5  | 106   | 0.00     |
| 86 I | Perylene-d12               | 1.000 | 1.000 | 0.0   | 106   | -0.01    |
| 87   | Indeno(1,2,3-cd)pyrene     | 1.490 | 1.600 | -7.4  | 106   | -0.01    |
| 88   | Benzo(b)fluoranthene       | 1.301 | 1.421 | -9.2  | 107   | 0.00     |
| 89   | Benzo(k)fluoranthene       | 1.316 | 1.424 | -8.2  | 109   | 0.00     |
| 90 C | Benzo(a)pyrene             | 1.218 | 1.306 | -7.2  | 107   | 0.00     |
| 91   | Dibenzo(a,h)anthracene     | 1.215 | 1.313 | -8.1  | 107   | -0.02    |
| 92   | Benzo(g,h,i)perylene       | 1.163 | 1.237 | -6.4  | 106   | -0.01    |

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

(#) = Out of Range

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