

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF072523\  
 Data File : BF134459.D  
 Acq On : 25 Jul 2023 12:48  
 Operator : CG\JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Jul 25 17:24:25 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF072423.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jul 25 01:46:04 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    | 125   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.513 | 0.550 | -7.2   | 136   | 0.01     |
| 3    | Pyridine                    | 1.153 | 1.338 | -16.0  | 136   | 0.02     |
| 4    | n-Nitrosodimethylamine      | 0.681 | 0.830 | -21.9  | 150#  | 0.02     |
| 5 S  | 2-Fluorophenol              | 1.211 | 1.201 | 0.8    | 119   | 0.00     |
| 6    | Aniline                     | 1.748 | 1.742 | 0.3    | 122   | 0.00     |
| 7 S  | Phenol-d6                   | 1.483 | 1.484 | -0.1   | 125   | 0.00     |
| 8    | 2-Chlorophenol              | 1.226 | 1.233 | -0.6   | 123   | 0.00     |
| 9    | Benzaldehyde                | 0.794 | 0.737 | 7.2    | 123   | 0.00     |
| 10 C | Phenol                      | 1.543 | 1.531 | 0.8    | 122   | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 1.077 | 1.097 | -1.9   | 129   | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 1.319 | 1.334 | -1.1   | 122   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.357 | 1.336 | 1.5    | 121   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 1.275 | 1.262 | 1.0    | 124   | 0.00     |
| 15   | Benzyl Alcohol              | 1.145 | 1.194 | -4.3   | 123   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 1.586 | 1.655 | -4.4   | 130   | 0.00     |
| 17   | 2-Methylphenol              | 1.094 | 1.091 | 0.3    | 122   | 0.00     |
| 18   | Hexachloroethane            | 0.521 | 0.535 | -2.7   | 129   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 0.923 | 0.960 | -4.0   | 129   | 0.00     |
| 20   | 3+4-Methylphenols           | 1.400 | 1.388 | 0.9    | 122   | 0.00     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 130   | 0.00     |
| 22   | Acetophenone                | 0.497 | 0.492 | 1.0    | 127   | 0.00     |
| 23 S | Nitrobenzene-d5             | 0.395 | 0.390 | 1.3    | 125   | 0.00     |
| 24   | Nitrobenzene                | 0.393 | 0.388 | 1.3    | 128   | 0.00     |
| 25   | Isophorone                  | 0.659 | 0.670 | -1.7   | 130   | 0.00     |
| 26 C | 2-Nitrophenol               | 0.181 | 0.188 | -3.9   | 131   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 0.285 | 0.290 | -1.8   | 131   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 0.367 | 0.383 | -4.4   | 135   | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.285 | 0.288 | -1.1   | 128   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 0.308 | 0.309 | -0.3   | 127   | 0.00     |
| 31   | Naphthalene                 | 0.995 | 0.992 | 0.3    | 127   | 0.00     |
| 32   | Benzoic acid                | 0.191 | 0.206 | -7.9   | 132   | 0.02     |
| 33   | 4-Chloroaniline             | 0.411 | 0.381 | 7.3    | 117   | 0.00     |
| 34 C | Hexachlorobutadiene         | 0.206 | 0.210 | -1.9   | 129   | 0.00     |
| 35   | Caprolactam                 | 0.087 | 0.091 | -4.6   | 130   | 0.01     |
| 36 C | 4-Chloro-3-methylphenol     | 0.329 | 0.335 | -1.8   | 130   | 0.00     |
| 37   | 2-Methylnaphthalene         | 0.692 | 0.705 | -1.9   | 130   | 0.00     |
| 38   | 1-Methylnaphthalene         | 0.642 | 0.658 | -2.5   | 132   | 0.00     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 119   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.634 | 0.646 | -1.9   | 132   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 0.152 | 0.235 | -54.6# | 184#  | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 0.284 | 0.308 | -8.5   | 125   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 0.407 | 0.422 | -3.7   | 120   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 0.474 | 0.481 | -1.5   | 116   | 0.00     |
| 45 S | 2-Fluorobiphenyl            | 1.523 | 1.555 | -2.1   | 119   | 0.00     |
| 46   | 1,1'-Biphenyl               | 1.648 | 1.702 | -3.3   | 120   | 0.00     |
| 47   | 2-Chloronaphthalene         | 1.215 | 1.226 | -0.9   | 118   | 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             | 0.438 | 0.448 | -2.3   | 115   | 0.00     |
| 49   | Acenaphthylene             | 2.053 | 2.076 | -1.1   | 117   | 0.00     |
| 50   | Dimethylphthalate          | 1.483 | 1.529 | -3.1   | 119   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 0.325 | 0.338 | -4.0   | 118   | 0.00     |
| 52 C | Acenaphthene               | 1.213 | 1.226 | -1.1   | 116   | 0.00     |
| 53   | 3-Nitroaniline             | 0.373 | 0.366 | 1.9    | 112   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 0.143 | 0.150 | -4.9   | 120   | 0.00     |
| 55   | Dibenzofuran               | 1.867 | 1.892 | -1.3   | 117   | 0.00     |
| 56 P | 4-Nitrophenol              | 0.199 | 0.211 | -6.0   | 122   | 0.01     |
| 57   | 2,4-Dinitrotoluene         | 0.437 | 0.443 | -1.4   | 114   | 0.00     |
| 58   | Fluorene                   | 1.506 | 1.553 | -3.1   | 119   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.364 | 0.395 | -8.5   | 125   | 0.00     |
| 60   | Diethylphthalate           | 1.535 | 1.636 | -6.6   | 124   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 0.719 | 0.761 | -5.8   | 122   | 0.00     |
| 62   | 4-Nitroaniline             | 0.383 | 0.342 | 10.7   | 101   | 0.00     |
| 63   | Azobenzene                 | 1.413 | 1.490 | -5.4   | 123   | 0.00     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0    | 119   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 0.115 | 0.119 | -3.5   | 118   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 0.636 | 0.655 | -3.0   | 120   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 0.219 | 0.227 | -3.7   | 122   | 0.00     |
| 68   | Hexachlorobenzene          | 0.247 | 0.256 | -3.6   | 122   | 0.00     |
| 69   | Atrazine                   | 0.177 | 0.183 | -3.4   | 117   | 0.00     |
| 70 C | Pentachlorophenol          | 0.104 | 0.124 | -19.2  | 137   | 0.00     |
| 71   | Phenanthrene               | 1.045 | 1.050 | -0.5   | 117   | 0.00     |
| 72   | Anthracene                 | 1.079 | 1.087 | -0.7   | 117   | 0.00     |
| 73   | Carbazole                  | 1.016 | 1.004 | 1.2    | 116   | 0.00     |
| 74   | Di-n-butylphthalate        | 1.166 | 1.263 | -8.3   | 126   | 0.00     |
| 75 C | Fluoranthene               | 1.218 | 1.190 | 2.3    | 115   | 0.00     |
| 76 I | Chrysene-d12               | 1.000 | 1.000 | 0.0    | 109   | 0.00     |
| 77   | Benzidine                  | 0.427 | 0.433 | -1.4   | 108   | 0.00     |
| 78   | Pyrene                     | 1.590 | 1.643 | -3.3   | 112   | 0.00     |
| 79 S | Terphenyl-d14              | 1.146 | 1.241 | -8.3   | 118   | 0.00     |
| 80   | Butylbenzylphthalate       | 0.637 | 0.726 | -14.0  | 121   | 0.00     |
| 81   | Benzo(a)anthracene         | 1.395 | 1.414 | -1.4   | 112   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 0.442 | 0.434 | 1.8    | 107   | 0.00     |
| 83   | Chrysene                   | 1.335 | 1.353 | -1.3   | 110   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 0.774 | 0.968 | -25.1# | 134   | 0.00     |
| 85 c | Di-n-octyl phthalate       | 1.075 | 1.322 | -23.0# | 132   | 0.00     |
| 86 I | Perylene-d12               | 1.000 | 1.000 | 0.0    | 100   | 0.00     |
| 87   | Indeno(1,2,3-cd)pyrene     | 1.307 | 1.447 | -10.7  | 105   | 0.00     |
| 88   | Benzo(b)fluoranthene       | 1.251 | 1.283 | -2.6   | 105   | 0.00     |
| 89   | Benzo(k)fluoranthene       | 1.230 | 1.280 | -4.1   | 100   | 0.00     |
| 90 C | Benzo(a)pyrene             | 1.157 | 1.162 | -0.4   | 99    | 0.00     |
| 91   | Dibenzo(a,h)anthracene     | 1.055 | 1.208 | -14.5  | 110   | 0.00     |
| 92   | Benzo(g,h,i)perylene       | 1.083 | 1.957 | -80.7# | 174#  | 0.00     |

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

(#) = Out of Range

SPCC's out = 0 CCC's out = 1