

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG051122\  
 Data File : BG053522.D  
 Acq On : 11 May 2022 12:02  
 Operator : CG/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTDCCC040

Quant Time: May 11 13:39:13 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG050522.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed May 11 06:00:46 2022  
 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  | 108   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.618 | 0.613 | 0.8  | 108   | 0.00     |
| 3    | Pyridine                    | 1.510 | 1.562 | -3.4 | 107   | 0.00     |
| 4    | n-Nitrosodimethylamine      | 0.736 | 0.748 | -1.6 | 107   | 0.00     |
| 5 S  | 2-Fluorophenol              | 1.178 | 1.172 | 0.5  | 107   | 0.00     |
| 6    | Aniline                     | 1.981 | 2.041 | -3.0 | 108   | 0.00     |
| 7 S  | Phenol-d6                   | 1.788 | 1.833 | -2.5 | 109   | 0.00     |
| 8    | 2-Chlorophenol              | 1.232 | 1.213 | 1.5  | 106   | 0.00     |
| 9    | Benzaldehyde                | 1.136 | 1.128 | 0.7  | 113   | 0.00     |
| 10 C | Phenol                      | 1.749 | 1.779 | -1.7 | 106   | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 1.305 | 1.275 | 2.3  | 102   | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 1.453 | 1.438 | 1.0  | 107   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.467 | 1.447 | 1.4  | 109   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 1.377 | 1.362 | 1.1  | 107   | 0.00     |
| 15   | Benzyl Alcohol              | 1.495 | 1.557 | -4.1 | 109   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 2.123 | 2.249 | -5.9 | 113   | 0.00     |
| 17   | 2-Methylphenol              | 1.182 | 1.200 | -1.5 | 109   | 0.00     |
| 18   | Hexachloroethane            | 0.557 | 0.537 | 3.6  | 106   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 1.283 | 1.320 | -2.9 | 110   | 0.00     |
| 20   | 3+4-Methylphenols           | 1.670 | 1.701 | -1.9 | 106   | 0.00     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0  | 109   | 0.00     |
| 22   | Acetophenone                | 0.547 | 0.547 | 0.0  | 109   | 0.00     |
| 23 S | Nitrobenzene-d5             | 0.470 | 0.482 | -2.6 | 109   | 0.00     |
| 24   | Nitrobenzene                | 0.455 | 0.467 | -2.6 | 108   | 0.00     |
| 25   | Isophorone                  | 0.791 | 0.825 | -4.3 | 109   | 0.00     |
| 26 C | 2-Nitrophenol               | 0.184 | 0.188 | -2.2 | 108   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 0.277 | 0.284 | -2.5 | 108   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 0.455 | 0.465 | -2.2 | 109   | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.335 | 0.350 | -4.5 | 107   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 0.380 | 0.384 | -1.1 | 109   | 0.00     |
| 31   | Naphthalene                 | 1.017 | 1.039 | -2.2 | 110   | 0.00     |
| 32   | Benzoic acid                | 0.147 | 0.135 | 8.2  | 92    | 0.00     |
| 33   | 4-Chloroaniline             | 0.426 | 0.445 | -4.5 | 108   | 0.00     |
| 34 C | Hexachlorobutadiene         | 0.282 | 0.282 | 0.0  | 107   | 0.00     |
| 35   | Caprolactam                 | 0.122 | 0.116 | 4.9  | 101   | -0.01    |
| 36 C | 4-Chloro-3-methylphenol     | 0.400 | 0.403 | -0.8 | 104   | 0.00     |
| 37   | 2-Methylnaphthalene         | 0.759 | 0.771 | -1.6 | 108   | 0.00     |
| 38   | 1-Methylnaphthalene         | 0.741 | 0.745 | -0.5 | 110   | 0.00     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0  | 109   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.630 | 0.650 | -3.2 | 110   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 0.200 | 0.200 | 0.0  | 101   | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 0.295 | 0.296 | -0.3 | 103   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 0.412 | 0.421 | -2.2 | 104   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 0.438 | 0.457 | -4.3 | 106   | 0.00     |
| 45 S | 2-Fluorobiphenyl            | 1.369 | 1.384 | -1.1 | 109   | 0.00     |
| 46   | 1,1'-Biphenyl               | 1.363 | 1.389 | -1.9 | 110   | 0.00     |
| 47   | 2-Chloronaphthalene         | 1.068 | 1.065 | 0.3  | 108   | 0.00     |

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 QLast Update : Wed May 11 06:00:46 2022  
 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) |
|------|----------------------------|-------|-------|------|-------|----------|
| 48   | 2-Nitroaniline             | 0.398 | 0.420 | -5.5 | 109   | 0.00     |
| 49   | Acenaphthylene             | 1.704 | 1.711 | -0.4 | 106   | 0.00     |
| 50   | Dimethylphthalate          | 1.520 | 1.491 | 1.9  | 104   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 0.310 | 0.303 | 2.3  | 102   | 0.00     |
| 52 C | Acenaphthene               | 1.016 | 1.011 | 0.5  | 105   | 0.00     |
| 53   | 3-Nitroaniline             | 0.335 | 0.328 | 2.1  | 101   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 0.166 | 0.164 | 1.2  | 100   | 0.00     |
| 55   | Dibenzofuran               | 1.824 | 1.816 | 0.4  | 108   | 0.00     |
| 56 P | 4-Nitrophenol              | 0.239 | 0.224 | 6.3  | 96    | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 0.449 | 0.426 | 5.1  | 98    | 0.00     |
| 58   | Fluorene                   | 1.464 | 1.437 | 1.8  | 104   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.444 | 0.444 | 0.0  | 104   | 0.00     |
| 60   | Diethylphthalate           | 1.570 | 1.525 | 2.9  | 102   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 0.817 | 0.835 | -2.2 | 110   | 0.00     |
| 62   | 4-Nitroaniline             | 0.347 | 0.336 | 3.2  | 102   | 0.00     |
| 63   | Azobenzene                 | 1.608 | 1.607 | 0.1  | 107   | 0.00     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0  | 102   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 0.119 | 0.127 | -6.7 | 102   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 0.525 | 0.541 | -3.0 | 103   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 0.234 | 0.248 | -6.0 | 105   | 0.00     |
| 68   | Hexachlorobenzene          | 0.256 | 0.259 | -1.2 | 103   | 0.00     |
| 69   | Atrazine                   | 0.221 | 0.216 | 2.3  | 99    | 0.00     |
| 70 C | Pentachlorophenol          | 0.121 | 0.122 | -0.8 | 98    | 0.00     |
| 71   | Phenanthrene               | 1.017 | 1.032 | -1.5 | 103   | 0.00     |
| 72   | Anthracene                 | 1.000 | 1.011 | -1.1 | 104   | 0.00     |
| 73   | Carbazole                  | 0.980 | 0.968 | 1.2  | 100   | 0.00     |
| 74   | Di-n-butylphthalate        | 1.102 | 1.079 | 2.1  | 99    | 0.00     |
| 75 C | Fluoranthene               | 1.317 | 1.301 | 1.2  | 101   | 0.00     |
| 76 I | Chrysene-d12               | 1.000 | 1.000 | 0.0  | 101   | 0.00     |
| 77   | Benzydine                  | 0.431 | 0.458 | -6.3 | 109   | 0.00     |
| 78   | Pyrene                     | 1.295 | 1.292 | 0.2  | 99    | 0.00     |
| 79 S | Terphenyl-d14              | 1.074 | 1.070 | 0.4  | 101   | 0.00     |
| 80   | Butylbenzylphthalate       | 0.475 | 0.490 | -3.2 | 101   | 0.00     |
| 81   | Benzo(a)anthracene         | 1.276 | 1.288 | -0.9 | 102   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 0.323 | 0.329 | -1.9 | 99    | 0.00     |
| 83   | Chrysene                   | 1.191 | 1.210 | -1.6 | 102   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 0.655 | 0.682 | -4.1 | 103   | 0.00     |
| 85 c | Di-n-octyl phthalate       | 1.108 | 1.168 | -5.4 | 104   | 0.00     |
| 86 I | Perylene-d12               | 1.000 | 1.000 | 0.0  | 102   | 0.00     |
| 87   | Indeno(1,2,3-cd)pyrene     | 1.393 | 1.441 | -3.4 | 104   | 0.00     |
| 88   | Benzo(b)fluoranthene       | 1.184 | 1.217 | -2.8 | 102   | 0.00     |
| 89   | Benzo(k)fluoranthene       | 1.163 | 1.168 | -0.4 | 104   | 0.00     |
| 90 C | Benzo(a)pyrene             | 1.008 | 1.027 | -1.9 | 103   | 0.00     |
| 91   | Dibenzo(a,h)anthracene     | 1.148 | 1.176 | -2.4 | 104   | 0.00     |
| 92   | Benzo(g,h,i)perylene       | 1.138 | 1.174 | -3.2 | 104   | 0.00     |

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

(#) = Out of Range

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