

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062723\  
 Data File : BG057873.D  
 Acq On : 27 Jun 2023 19:15  
 Operator : CG/JU  
 Sample : SSTDICV040  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 ICVBG062723

Quant Time: Jun 28 01:13:52 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG062723.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Jun 28 01:12:33 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   | 104   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.599 | 0.587 | 2.0   | 104   | 0.00     |
| 3    | Pyridine                    | 1.729 | 1.772 | -2.5  | 108   | 0.00     |
| 4    | n-Nitrosodimethylamine      | 0.713 | 0.706 | 1.0   | 104   | 0.00     |
| 5 S  | 2-Fluorophenol              | 1.307 | 1.283 | 1.8   | 103   | 0.00     |
| 6    | Aniline                     | 2.442 | 2.479 | -1.5  | 110   | 0.00     |
| 7 S  | Phenol-d6                   | 1.955 | 2.000 | -2.3  | 109   | 0.00     |
| 8    | 2-Chlorophenol              | 1.312 | 1.317 | -0.4  | 106   | 0.00     |
| 9    | Benzaldehyde                | 1.095 | 1.075 | 1.8   | 109   | 0.00     |
| 10 C | Phenol                      | 1.899 | 1.933 | -1.8  | 108   | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 1.432 | 1.449 | -1.2  | 108   | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 1.350 | 1.323 | 2.0   | 104   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.356 | 1.336 | 1.5   | 104   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 1.312 | 1.311 | 0.1   | 106   | 0.00     |
| 15   | Benzyl Alcohol              | 1.440 | 1.513 | -5.1  | 115   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 2.385 | 2.354 | 1.3   | 108   | 0.00     |
| 17   | 2-Methylphenol              | 1.408 | 1.433 | -1.8  | 109   | 0.00     |
| 18   | Hexachloroethane            | 0.472 | 0.471 | 0.2   | 104   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 1.308 | 1.352 | -3.4  | 116   | 0.00     |
| 20   | 3+4-Methylphenols           | 1.962 | 2.068 | -5.4  | 114   | 0.00     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 111   | 0.00     |
| 22   | Acetophenone                | 0.557 | 0.564 | -1.3  | 112   | 0.00     |
| 23 S | Nitrobenzene-d5             | 0.408 | 0.416 | -2.0  | 112   | 0.00     |
| 24   | Nitrobenzene                | 0.410 | 0.423 | -3.2  | 113   | 0.00     |
| 25   | Isophorone                  | 0.881 | 0.923 | -4.8  | 120   | 0.00     |
| 26 C | 2-Nitrophenol               | 0.167 | 0.181 | -8.4  | 116   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 0.322 | 0.331 | -2.8  | 113   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 0.474 | 0.483 | -1.9  | 115   | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.316 | 0.328 | -3.8  | 114   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 0.340 | 0.339 | 0.3   | 109   | 0.00     |
| 31   | Naphthalene                 | 1.068 | 1.081 | -1.2  | 113   | 0.00     |
| 32   | Benzoic acid                | 0.252 | 0.280 | -11.1 | 126   | 0.00     |
| 33   | 4-Chloroaniline             | 0.502 | 0.533 | -6.2  | 119   | 0.00     |
| 34 C | Hexachlorobutadiene         | 0.210 | 0.207 | 1.4   | 109   | 0.00     |
| 35   | Caprolactam                 | 0.137 | 0.149 | -8.8  | 125   | 0.00     |
| 36 C | 4-Chloro-3-methylphenol     | 0.419 | 0.442 | -5.5  | 121   | 0.00     |
| 37   | 2-Methylnaphthalene         | 0.787 | 0.820 | -4.2  | 119   | 0.00     |
| 38   | 1-Methylnaphthalene         | 0.744 | 0.770 | -3.5  | 119   | 0.00     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 124   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.560 | 0.548 | 2.1   | 119   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 0.301 | 0.302 | -0.3  | 117   | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 0.336 | 0.339 | -0.9  | 121   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 0.412 | 0.417 | -1.2  | 123   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 0.491 | 0.497 | -1.2  | 122   | 0.00     |
| 45 S | 2-Fluorobiphenyl            | 1.311 | 1.263 | 3.7   | 120   | 0.00     |
| 46   | 1,1'-Biphenyl               | 1.362 | 1.320 | 3.1   | 120   | 0.00     |
| 47   | 2-Chloronaphthalene         | 1.046 | 1.025 | 2.0   | 121   | 0.00     |

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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
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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             | 0.417 | 0.433 | -3.8 | 123   | 0.00     |
| 49   | Acenaphthylene             | 1.835 | 1.794 | 2.2  | 122   | 0.00     |
| 50   | Dimethylphthalate          | 1.585 | 1.566 | 1.2  | 123   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 0.329 | 0.345 | -4.9 | 125   | 0.00     |
| 52 C | Acenaphthene               | 1.072 | 1.067 | 0.5  | 123   | 0.00     |
| 53   | 3-Nitroaniline             | 0.374 | 0.382 | -2.1 | 122   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 0.182 | 0.195 | -7.1 | 127   | 0.00     |
| 55   | Dibenzofuran               | 1.838 | 1.800 | 2.1  | 122   | 0.00     |
| 56 P | 4-Nitrophenol              | 0.307 | 0.317 | -3.3 | 117   | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 0.478 | 0.499 | -4.4 | 121   | 0.00     |
| 58   | Fluorene                   | 1.484 | 1.451 | 2.2  | 122   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.445 | 0.451 | -1.3 | 124   | 0.00     |
| 60   | Diethylphthalate           | 1.671 | 1.639 | 1.9  | 120   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 0.796 | 0.786 | 1.3  | 124   | 0.00     |
| 62   | 4-Nitroaniline             | 0.392 | 0.388 | 1.0  | 114   | 0.00     |
| 63   | Azobenzene                 | 1.598 | 1.593 | 0.3  | 123   | 0.00     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0  | 118   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 0.108 | 0.116 | -7.4 | 122   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 0.510 | 0.511 | -0.2 | 121   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 0.211 | 0.214 | -1.4 | 124   | 0.00     |
| 68   | Hexachlorobenzene          | 0.223 | 0.223 | 0.0  | 119   | 0.00     |
| 69   | Atrazine                   | 0.190 | 0.179 | 5.8  | 114   | 0.00     |
| 70 C | Pentachlorophenol          | 0.170 | 0.176 | -3.5 | 118   | 0.00     |
| 71   | Phenanthrene               | 0.960 | 0.940 | 2.1  | 115   | 0.00     |
| 72   | Anthracene                 | 0.990 | 0.966 | 2.4  | 114   | 0.00     |
| 73   | Carbazole                  | 0.941 | 0.904 | 3.9  | 109   | 0.00     |
| 74   | Di-n-butylphthalate        | 1.164 | 1.124 | 3.4  | 109   | 0.00     |
| 75 C | Fluoranthene               | 1.209 | 1.151 | 4.8  | 108   | 0.00     |
| 76 I | Chrysene-d12               | 1.000 | 1.000 | 0.0  | 107   | 0.00     |
| 77   | Benzydine                  | 0.469 | 0.457 | 2.6  | 95    | 0.00     |
| 78   | Pyrene                     | 1.428 | 1.404 | 1.7  | 107   | 0.00     |
| 79 S | Terphenyl-d14              | 1.097 | 1.060 | 3.4  | 106   | 0.00     |
| 80   | Butylbenzylphthalate       | 0.579 | 0.578 | 0.2  | 105   | 0.00     |
| 81   | Benzo(a)anthracene         | 1.384 | 1.365 | 1.4  | 106   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 0.492 | 0.482 | 2.0  | 105   | 0.00     |
| 83   | Chrysene                   | 1.310 | 1.301 | 0.7  | 108   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 0.822 | 0.805 | 2.1  | 105   | 0.00     |
| 85 c | Di-n-octyl phthalate       | 1.451 | 1.455 | -0.3 | 106   | 0.00     |
| 86 I | Perylene-d12               | 1.000 | 1.000 | 0.0  | 107   | 0.00     |
| 87   | Indeno(1,2,3-cd)pyrene     | 1.325 | 1.335 | -0.8 | 107   | -0.01    |
| 88   | Benzo(b)fluoranthene       | 1.091 | 1.081 | 0.9  | 106   | 0.00     |
| 89   | Benzo(k)fluoranthene       | 1.102 | 1.124 | -2.0 | 108   | 0.00     |
| 90 C | Benzo(a)pyrene             | 1.081 | 1.078 | 0.3  | 106   | 0.00     |
| 91   | Dibenzo(a,h)anthracene     | 1.099 | 1.110 | -1.0 | 108   | -0.01    |
| 92   | Benzo(g,h,i)perylene       | 1.092 | 1.101 | -0.8 | 107   | 0.00     |

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

(#) = Out of Range

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