

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP110620\  
 Data File : BP003851.D  
 Acq On : 07 Nov 2020 00:13  
 Operator : CG/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 LabSampleId :  
 SSTDCCC040

Quant Time: Nov 07 03:16:13 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\8270-BP110320.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Nov 06 11:33:14 2020  
 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   | 118   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.517 | 0.445 | 13.9  | 108   | 0.00     |
| 3    | Pyridine                    | 1.516 | 1.324 | 12.7  | 107   | 0.00     |
| 4    | n-Nitrosodimethylamine      | 0.698 | 0.612 | 12.3  | 109   | 0.00     |
| 5 S  | 2-Fluorophenol              | 1.198 | 1.134 | 5.3   | 117   | 0.00     |
| 6    | Aniline                     | 1.953 | 1.762 | 9.8   | 110   | 0.00     |
| 7 S  | Phenol-d6                   | 1.720 | 1.573 | 8.5   | 112   | 0.00     |
| 8    | 2-Chlorophenol              | 1.272 | 1.198 | 5.8   | 115   | 0.00     |
| 9    | Benzaldehyde                | 0.873 | 0.772 | 11.6  | 125   | 0.00     |
| 10 C | Phenol                      | 1.660 | 1.511 | 9.0   | 112   | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 1.334 | 1.180 | 11.5  | 109   | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 1.559 | 1.460 | 6.4   | 117   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.572 | 1.466 | 6.7   | 116   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 1.501 | 1.388 | 7.5   | 115   | 0.00     |
| 15   | Benzyl Alcohol              | 1.191 | 1.104 | 7.3   | 110   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 2.247 | 1.888 | 16.0  | 104   | 0.00     |
| 17   | 2-Methylphenol              | 1.090 | 0.996 | 8.6   | 111   | 0.00     |
| 18   | Hexachloroethane            | 0.559 | 0.471 | 15.7  | 102   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 1.028 | 0.926 | 9.9   | 108   | 0.00     |
| 20   | 3+4-Methylphenols           | 1.455 | 1.343 | 7.7   | 111   | 0.00     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 111   | 0.00     |
| 22   | Acetophenone                | 0.537 | 0.482 | 10.2  | 107   | 0.00     |
| 23 S | Nitrobenzene-d5             | 0.408 | 0.368 | 9.8   | 106   | 0.00     |
| 24   | Nitrobenzene                | 0.406 | 0.369 | 9.1   | 108   | 0.00     |
| 25   | Isophorone                  | 0.694 | 0.643 | 7.3   | 108   | 0.00     |
| 26 C | 2-Nitrophenol               | 0.158 | 0.160 | -1.3  | 113   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 0.264 | 0.248 | 6.1   | 112   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 0.476 | 0.418 | 12.2  | 105   | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.319 | 0.311 | 2.5   | 114   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 0.388 | 0.369 | 4.9   | 114   | 0.00     |
| 31   | Naphthalene                 | 1.073 | 0.984 | 8.3   | 111   | 0.00     |
| 32   | Benzoic acid                | 0.161 | 0.164 | -1.9  | 114   | 0.00     |
| 33   | 4-Chloroaniline             | 0.456 | 0.413 | 9.4   | 107   | 0.00     |
| 34 C | Hexachlorobutadiene         | 0.253 | 0.242 | 4.3   | 116   | 0.00     |
| 35   | Caprolactam                 | 0.098 | 0.097 | 1.0   | 111   | 0.00     |
| 36 C | 4-Chloro-3-methylphenol     | 0.344 | 0.323 | 6.1   | 110   | 0.00     |
| 37   | 2-Methylnaphthalene         | 0.768 | 0.694 | 9.6   | 109   | 0.00     |
| 38   | 1-Methylnaphthalene         | 0.733 | 0.655 | 10.6  | 109   | 0.00     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 107   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.662 | 0.632 | 4.5   | 111   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 0.334 | 0.080 | 76.0# | 27#   | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 0.250 | 0.256 | -2.4  | 114   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 0.409 | 0.418 | -2.2  | 114   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 0.431 | 0.432 | -0.2  | 113   | 0.00     |
| 45 S | 2-Fluorobiphenyl            | 1.431 | 1.320 | 7.8   | 108   | 0.00     |
| 46   | 1,1'-Biphenyl               | 1.553 | 1.423 | 8.4   | 107   | 0.00     |
| 47   | 2-Chloronaphthalene         | 1.274 | 1.183 | 7.1   | 108   | 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             | 0.380 | 0.396  | -4.2   | 113   | 0.00     |
| 49   | Acenaphthylene             | 1.871 | 1.738  | 7.1    | 107   | 0.00     |
| 50   | Dimethylphthalate          | 1.563 | 1.414  | 9.5    | 105   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 0.321 | 0.296  | 7.8    | 104   | 0.00     |
| 52 C | Acenaphthene               | 1.246 | 1.085  | 12.9   | 101   | 0.00     |
| 53   | 3-Nitroaniline             | 0.355 | 0.353  | 0.6    | 112   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 0.138 | 0.022# | 84.1#  | 18#   | 0.00     |
| 55   | Dibenzofuran               | 1.851 | 1.691  | 8.6    | 107   | 0.00     |
| 56 P | 4-Nitrophenol              | 0.256 | 0.251  | 2.0    | 107   | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 0.431 | 0.398  | 7.7    | 102   | 0.00     |
| 58   | Fluorene                   | 1.482 | 1.350  | 8.9    | 107   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.388 | 0.384  | 1.0    | 111   | 0.00     |
| 60   | Diethylphthalate           | 1.531 | 1.389  | 9.3    | 105   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 0.810 | 0.739  | 8.8    | 108   | 0.00     |
| 62   | 4-Nitroaniline             | 0.370 | 0.381  | -3.0   | 116   | 0.00     |
| 63   | Azobenzene                 | 1.445 | 1.277  | 11.6   | 102   | 0.00     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000  | 0.0    | 108   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 0.098 | 0.022  | 77.6#  | 25#   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 0.612 | 0.565  | 7.7    | 107   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 0.234 | 0.223  | 4.7    | 110   | 0.00     |
| 68   | Hexachlorobenzene          | 0.267 | 0.250  | 6.4    | 110   | 0.00     |
| 69   | Atrazine                   | 0.201 | 0.186  | 7.5    | 104   | 0.00     |
| 70 C | Pentachlorophenol          | 0.128 | 0.122  | 4.7    | 106   | 0.00     |
| 71   | Phenanthrene               | 1.122 | 1.019  | 9.2    | 107   | 0.00     |
| 72   | Anthracene                 | 1.084 | 1.004  | 7.4    | 108   | 0.00     |
| 73   | Carbazole                  | 1.006 | 0.934  | 7.2    | 108   | 0.00     |
| 74   | Di-n-butylphthalate        | 1.164 | 1.147  | 1.5    | 109   | 0.00     |
| 75 C | Fluoranthene               | 1.295 | 1.190  | 8.1    | 106   | 0.00     |
| 76 I | Chrysene-d12               | 1.000 | 1.000  | 0.0    | 101   | 0.00     |
| 77   | Benzydine                  | 0.472 | 0.623  | -32.0# | 132   | 0.00     |
| 78   | Pyrene                     | 1.357 | 1.327  | 2.2    | 105   | 0.00     |
| 79 S | Terphenyl-d14              | 1.035 | 1.028  | 0.7    | 109   | 0.00     |
| 80   | Butylbenzylphthalate       | 0.506 | 0.559  | -10.5  | 112   | 0.00     |
| 81   | Benzo(a)anthracene         | 1.319 | 1.243  | 5.8    | 102   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 0.455 | 0.473  | -4.0   | 111   | 0.00     |
| 83   | Chrysene                   | 1.267 | 1.168  | 7.8    | 101   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 0.745 | 0.778  | -4.4   | 108   | 0.00     |
| 85 c | Di-n-octyl phthalate       | 1.242 | 1.323  | -6.5   | 109   | 0.00     |
| 86 I | Perylene-d12               | 1.000 | 1.000  | 0.0    | 104   | 0.00     |
| 87   | Indeno(1,2,3-cd)pyrene     | 1.443 | 1.304  | 9.6    | 104   | 0.00     |
| 88   | Benzo(b)fluoranthene       | 1.309 | 1.159  | 11.5   | 104   | 0.00     |
| 89   | Benzo(k)fluoranthene       | 1.292 | 1.119  | 13.4   | 99    | 0.00     |
| 90 C | Benzo(a)pyrene             | 1.212 | 1.081  | 10.8   | 103   | 0.00     |
| 91   | Dibenzo(a,h)anthracene     | 1.203 | 1.089  | 9.5    | 104   | 0.00     |
| 92   | Benzo(g,h,i)perylene       | 1.164 | 1.065  | 8.5    | 106   | 0.00     |

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

(#) = Out of Range                      SPCC's out = 1    CCC's out = 0