

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF042220\  
 Data File : BF120122.D  
 Acq On : 22 Apr 2020 11:11  
 Operator : MA/SJ  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Apr 22 12:17:01 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF040820.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Apr 22 11:49:00 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   | 81    | 0.02     |
| 2    | 1,4-Dioxane                 | 0.515 | 0.537 | -4.3  | 83    | 0.01     |
| 3    | Pyridine                    | 1.414 | 1.504 | -6.4  | 89    | 0.02     |
| 4    | n-Nitrosodimethylamine      | 0.723 | 0.839 | -16.0 | 96    | 0.02     |
| 5 S  | 2-Fluorophenol              | 1.157 | 1.382 | -19.4 | 99    | 0.02     |
| 6    | Aniline                     | 1.957 | 1.851 | 5.4   | 77    | 0.02     |
| 7 S  | Phenol-d6                   | 1.491 | 1.791 | -20.1 | 96    | 0.02     |
| 8    | 2-Chlorophenol              | 1.293 | 1.517 | -17.3 | 96    | 0.02     |
| 9    | Benzaldehyde                | 0.824 | 0.915 | -11.0 | 92    | 0.02     |
| 10 C | Phenol                      | 1.692 | 1.942 | -14.8 | 94    | 0.02     |
| 11   | bis(2-Chloroethyl)ether     | 1.306 | 1.477 | -13.1 | 93    | 0.01     |
| 12   | 1,3-Dichlorobenzene         | 1.416 | 1.638 | -15.7 | 95    | 0.02     |
| 13 C | 1,4-Dichlorobenzene         | 1.442 | 1.631 | -13.1 | 94    | 0.02     |
| 14   | 1,2-Dichlorobenzene         | 1.329 | 1.540 | -15.9 | 94    | 0.02     |
| 15   | Benzyl Alcohol              | 1.158 | 1.318 | -13.8 | 93    | 0.01     |
| 16   | 2,2'-oxybis(1-Chloropropane | 2.542 | 2.685 | -5.6  | 85    | 0.01     |
| 17   | 2-Methylphenol              | 1.154 | 1.307 | -13.3 | 92    | 0.01     |
| 18   | Hexachloroethane            | 0.539 | 0.614 | -13.9 | 93    | 0.02     |
| 19 P | n-Nitroso-di-n-propylamine  | 0.959 | 1.032 | -7.6  | 85    | 0.02     |
| 20   | 3+4-Methylphenols           | 1.411 | 1.634 | -15.8 | 90    | 0.02     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 79    | 0.02     |
| 22   | Acetophenone                | 0.468 | 0.547 | -16.9 | 92    | 0.02     |
| 23 S | Nitrobenzene-d5             | 0.387 | 0.443 | -14.5 | 93    | 0.01     |
| 24   | Nitrobenzene                | 0.389 | 0.450 | -15.7 | 94    | 0.01     |
| 25   | Isophorone                  | 0.745 | 0.797 | -7.0  | 82    | 0.01     |
| 26 C | 2-Nitrophenol               | 0.194 | 0.225 | -16.0 | 86    | 0.02     |
| 27   | 2,4-Dimethylphenol          | 0.293 | 0.312 | -6.5  | 82    | 0.02     |
| 28   | bis(2-Chloroethoxy)methane  | 0.439 | 0.485 | -10.5 | 89    | 0.01     |
| 29 C | 2,4-Dichlorophenol          | 0.300 | 0.335 | -11.7 | 90    | 0.02     |
| 30   | 1,2,4-Trichlorobenzene      | 0.340 | 0.378 | -11.2 | 91    | 0.02     |
| 31   | Naphthalene                 | 1.052 | 1.176 | -11.8 | 92    | 0.02     |
| 32   | Benzoic acid                | 0.257 | 0.256 | 0.4   | 81    | 0.02     |
| 33   | 4-Chloroaniline             | 0.458 | 0.457 | 0.2   | 83    | 0.02     |
| 34 C | Hexachlorobutadiene         | 0.204 | 0.218 | -6.9  | 86    | 0.02     |
| 35   | Caprolactam                 | 0.092 | 0.099 | -7.6  | 90    | 0.02     |
| 36 C | 4-Chloro-3-methylphenol     | 0.325 | 0.354 | -8.9  | 89    | 0.02     |
| 37   | 2-Methylnaphthalene         | 0.713 | 0.766 | -7.4  | 89    | 0.02     |
| 38   | 1-Methylnaphthalene         | 0.672 | 0.723 | -7.6  | 83    | 0.02     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 74    | 0.02     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.632 | 0.732 | -15.8 | 82    | 0.02     |
| 41 P | Hexachlorocyclopentadiene   | 0.359 | 0.348 | 3.1   | 69    | 0.02     |
| 42 S | 2,4,6-Tribromophenol        | 0.245 | 0.246 | -0.4  | 72    | 0.02     |
| 43 C | 2,4,6-Trichlorophenol       | 0.436 | 0.486 | -11.5 | 83    | 0.02     |
| 44   | 2,4,5-Trichlorophenol       | 0.491 | 0.546 | -11.2 | 78    | 0.02     |

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 Max. RRF Dev : 25% Max. Rel. Area : 150%

|      | Compound                   | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|----------------------------|-------|-------|-------|-------|----------|
| 45 S | 2-Fluorobiphenyl           | 1.409 | 1.621 | -15.0 | 83    | 0.02     |
| 46   | 1,1'-Biphenyl              | 1.688 | 1.944 | -15.2 | 83    | 0.02     |
| 47   | 2-Chloronaphthalene        | 1.300 | 1.509 | -16.1 | 86    | 0.02     |
| 48   | 2-Nitroaniline             | 0.471 | 0.552 | -17.2 | 89    | 0.02     |
| 49   | Acenaphthylene             | 1.978 | 2.243 | -13.4 | 85    | 0.02     |
| 50   | Dimethylphthalate          | 1.547 | 1.696 | -9.6  | 83    | 0.02     |
| 51   | 2,6-Dinitrotoluene         | 0.339 | 0.379 | -11.8 | 84    | 0.02     |
| 52 C | Acenaphthene               | 1.319 | 1.348 | -2.2  | 78    | 0.02     |
| 53   | 3-Nitroaniline             | 0.387 | 0.444 | -14.7 | 90    | 0.02     |
| 54 P | 2,4-Dinitrophenol          | 0.203 | 0.207 | -2.0  | 77    | 0.02     |
| 55   | Dibenzofuran               | 1.810 | 2.024 | -11.8 | 84    | 0.02     |
| 56 P | 4-Nitrophenol              | 0.288 | 0.314 | -9.0  | 83    | 0.02     |
| 57   | 2,4-Dinitrotoluene         | 0.446 | 0.488 | -9.4  | 83    | 0.02     |
| 58   | Fluorene                   | 1.448 | 1.597 | -10.3 | 87    | 0.02     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.376 | 0.407 | -8.2  | 81    | 0.02     |
| 60   | Diethylphthalate           | 1.603 | 1.720 | -7.3  | 82    | 0.02     |
| 61   | 4-Chlorophenyl-phenylether | 0.742 | 0.778 | -4.9  | 81    | 0.02     |
| 62   | 4-Nitroaniline             | 0.369 | 0.460 | -24.7 | 94    | 0.02     |
| 63   | Azobenzene                 | 1.437 | 1.611 | -12.1 | 84    | 0.02     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 77    | 0.02     |
| 65   | 4,6-Dinitro-2-methylphenol | 0.132 | 0.148 | -12.1 | 86    | 0.02     |
| 66 c | n-Nitrosodiphenylamine     | 0.665 | 0.739 | -11.1 | 87    | 0.02     |
| 67   | 4-Bromophenyl-phenylether  | 0.249 | 0.249 | 0.0   | 76    | 0.02     |
| 68   | Hexachlorobenzene          | 0.252 | 0.264 | -4.8  | 80    | 0.02     |
| 69   | Atrazine                   | 0.185 | 0.175 | 5.4   | 70    | 0.02     |
| 70 C | Pentachlorophenol          | 0.145 | 0.132 | 9.0   | 68    | 0.02     |
| 71   | Phenanthrene               | 1.064 | 1.191 | -11.9 | 87    | 0.02     |
| 72   | Anthracene                 | 1.087 | 1.195 | -9.9  | 85    | 0.02     |
| 73   | Carbazole                  | 1.028 | 1.177 | -14.5 | 88    | 0.02     |
| 74   | Di-n-butylphthalate        | 1.352 | 1.382 | -2.2  | 77    | 0.02     |
| 75 C | Fluoranthene               | 1.148 | 1.304 | -13.6 | 90    | 0.02     |
| 76 I | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 91    | 0.02     |
| 77   | Benzidine                  | 0.676 | 0.161 | 76.2# | 23#   | 0.02     |
| 78   | Pyrene                     | 1.686 | 1.791 | -6.2  | 90    | 0.02     |
| 79 S | Terphenyl-d14              | 1.189 | 1.157 | 2.7   | 82    | 0.02     |
| 80   | Butylbenzylphthalate       | 0.818 | 0.787 | 3.8   | 84    | 0.02     |
| 81   | Benzo(a)anthracene         | 1.316 | 1.451 | -10.3 | 102   | 0.02     |
| 82   | 3,3'-Dichlorobenzidine     | 0.491 | 0.489 | 0.4   | 91    | 0.02     |
| 83   | Chrysene                   | 1.337 | 1.467 | -9.7  | 103   | 0.03     |
| 84   | Bis(2-ethylhexyl)phthalate | 1.108 | 0.993 | 10.4  | 83    | 0.02     |
| 85 c | Di-n-octyl phthalate       | 1.886 | 1.776 | 5.8   | 87    | 0.02     |
| 86   | Indeno(1,2,3-cd)pyrene     | 1.178 | 1.189 | -0.9  | 86    | 0.04     |
| 87 I | Perylene-d12               | 1.000 | 1.000 | 0.0   | 87    | 0.03     |

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| 88   | Benzo(b)fluoranthene   | 1.439 | 1.584 | -10.1 | 103   | 0.03     |
| 89   | Benzo(k)fluoranthene   | 1.241 | 1.468 | -18.3 | 90    | 0.03     |
| 90 C | Benzo(a)pyrene         | 1.210 | 1.391 | -15.0 | 101   | 0.03     |
| 91   | Dibenzo(a,h)anthracene | 1.072 | 1.130 | -5.4  | 83    | 0.04     |
| 92   | Benzo(g,h,i)perylene   | 1.058 | 1.117 | -5.6  | 87    | 0.05     |

(#) = Out of Range

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