

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF073119\  
 Data File : BF115868.D  
 Acq On : 31 Jul 2019 16:15  
 Operator : HP/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 LabSampleId :  
 SSTDCCC040

Quant Time: Aug 01 04:54:31 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Jul 31 15:31:51 2019  
 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                    | AvqRF | CCRF  | %Dev | Area% | Dev(min) |
|------|-----------------------------|-------|-------|------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0  | 129   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.604 | 0.592 | 2.0  | 129   | 0.00     |
| 3    | Pyridine                    | 1.686 | 1.649 | 2.2  | 128   | 0.00     |
| 4    | n-Nitrosodimethylamine      | 0.665 | 0.685 | -3.0 | 134   | 0.01     |
| 5 S  | 2-Fluorophenol              | 1.195 | 1.174 | 1.8  | 125   | 0.00     |
| 6    | Aniline                     | 2.159 | 2.196 | -1.7 | 129   | 0.00     |
| 7 S  | Phenol-d6                   | 1.507 | 1.509 | -0.1 | 129   | 0.00     |
| 8    | 2-Chlorophenol              | 1.321 | 1.334 | -1.0 | 129   | 0.00     |
| 9    | Benzaldehyde                | 1.040 | 1.020 | 1.9  | 125   | 0.00     |
| 10 C | Phenol                      | 1.775 | 1.787 | -0.7 | 129   | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 1.305 | 1.314 | -0.7 | 129   | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 1.527 | 1.523 | 0.3  | 128   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.529 | 1.516 | 0.9  | 127   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 1.408 | 1.411 | -0.2 | 125   | 0.00     |
| 15   | Benzyl Alcohol              | 1.144 | 1.142 | 0.2  | 125   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 1.780 | 1.769 | 0.6  | 126   | 0.00     |
| 17   | 2-Methylphenol              | 1.119 | 1.123 | -0.4 | 131   | 0.00     |
| 18   | Hexachloroethane            | 0.563 | 0.562 | 0.2  | 127   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 0.934 | 0.912 | 2.4  | 127   | 0.00     |
| 20   | 3+4-Methylphenols           | 1.324 | 1.307 | 1.3  | 124   | 0.00     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0  | 128   | 0.00     |
| 22   | Acetophenone                | 0.439 | 0.435 | 0.9  | 125   | 0.00     |
| 23 S | Nitrobenzene-d5             | 0.346 | 0.350 | -1.2 | 128   | 0.00     |
| 24   | Nitrobenzene                | 0.374 | 0.379 | -1.3 | 128   | 0.00     |
| 25   | Isophorone                  | 0.663 | 0.659 | 0.6  | 127   | 0.00     |
| 26 C | 2-Nitrophenol               | 0.176 | 0.181 | -2.8 | 129   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 0.265 | 0.269 | -1.5 | 128   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 0.411 | 0.410 | 0.2  | 127   | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.280 | 0.284 | -1.4 | 126   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 0.319 | 0.321 | -0.6 | 126   | 0.00     |
| 31   | Naphthalene                 | 0.941 | 0.951 | -1.1 | 125   | 0.00     |
| 32   | Benzoic acid                | 0.187 | 0.174 | 7.0  | 129   | 0.01     |
| 33   | 4-Chloroaniline             | 0.421 | 0.421 | 0.0  | 126   | 0.00     |
| 34 C | Hexachlorobutadiene         | 0.184 | 0.186 | -1.1 | 127   | 0.00     |
| 35   | Caprolactam                 | 0.091 | 0.093 | -2.2 | 129   | 0.00     |
| 36 C | 4-Chloro-3-methylphenol     | 0.291 | 0.292 | -0.3 | 127   | 0.00     |
| 37   | 2-Methylnaphthalene         | 0.620 | 0.619 | 0.2  | 125   | 0.00     |
| 38   | 1-Methylnaphthalene         | 0.586 | 0.581 | 0.9  | 124   | 0.00     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0  | 125   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.535 | 0.538 | -0.6 | 124   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 0.209 | 0.222 | -6.2 | 128   | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 0.194 | 0.189 | 2.6  | 120   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 0.368 | 0.370 | -0.5 | 124   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 0.380 | 0.378 | 0.5  | 122   | 0.00     |

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

|      | Compound                   | AvqRF | CCRF  | %Dev | Area% | Dev(min) |
|------|----------------------------|-------|-------|------|-------|----------|
| 45 S | 2-Fluorobiphenyl           | 1.068 | 1.051 | 1.6  | 120   | 0.00     |
| 46   | 1,1'-Biphenyl              | 1.412 | 1.399 | 0.9  | 122   | 0.00     |
| 47   | 2-Chloronaphthalene        | 1.175 | 1.177 | -0.2 | 124   | 0.00     |
| 48   | 2-Nitroaniline             | 0.388 | 0.392 | -1.0 | 125   | 0.00     |
| 49   | Acenaphthylene             | 1.715 | 1.714 | 0.1  | 122   | 0.00     |
| 50   | Dimethylphthalate          | 1.362 | 1.322 | 2.9  | 120   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 0.296 | 0.294 | 0.7  | 123   | 0.00     |
| 52 C | Acenaphthene               | 1.052 | 1.048 | 0.4  | 122   | 0.00     |
| 53   | 3-Nitroaniline             | 0.366 | 0.359 | 1.9  | 122   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 0.126 | 0.127 | -0.8 | 122   | 0.00     |
| 55   | Dibenzofuran               | 1.530 | 1.529 | 0.1  | 121   | 0.00     |
| 56 P | 4-Nitrophenol              | 0.234 | 0.234 | 0.0  | 122   | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 0.394 | 0.393 | 0.3  | 121   | 0.00     |
| 58   | Fluorene                   | 1.177 | 1.154 | 2.0  | 120   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.320 | 0.321 | -0.3 | 124   | 0.00     |
| 60   | Diethylphthalate           | 1.271 | 1.238 | 2.6  | 119   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 0.596 | 0.583 | 2.2  | 118   | 0.00     |
| 62   | 4-Nitroaniline             | 0.378 | 0.373 | 1.3  | 122   | 0.00     |
| 63   | Azobenzene                 | 1.308 | 1.270 | 2.9  | 119   | 0.00     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0  | 121   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 0.106 | 0.108 | -1.9 | 119   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 0.602 | 0.605 | -0.5 | 120   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 0.214 | 0.214 | 0.0  | 118   | 0.00     |
| 68   | Hexachlorobenzene          | 0.248 | 0.246 | 0.8  | 118   | 0.00     |
| 69   | Atrazine                   | 0.198 | 0.195 | 1.5  | 118   | 0.00     |
| 70 C | Pentachlorophenol          | 0.120 | 0.121 | -0.8 | 117   | 0.00     |
| 71   | Phenanthrene               | 1.022 | 1.010 | 1.2  | 117   | 0.00     |
| 72   | Anthracene                 | 1.035 | 1.032 | 0.3  | 119   | 0.00     |
| 73   | Carbazole                  | 0.939 | 0.943 | -0.4 | 119   | 0.00     |
| 74   | Di-n-butylphthalate        | 1.095 | 1.056 | 3.6  | 111   | 0.00     |
| 75 C | Fluoranthene               | 1.070 | 1.061 | 0.8  | 118   | 0.00     |
| 76 I | Chrysene-d12               | 1.000 | 1.000 | 0.0  | 112   | 0.00     |
| 77   | Benzidine                  | 0.676 | 0.704 | -4.1 | 109   | 0.00     |
| 78   | Pyrene                     | 1.508 | 1.544 | -2.4 | 117   | 0.00     |
| 79 S | Terphenyl-d14              | 0.949 | 0.939 | 1.1  | 113   | 0.00     |
| 80   | Butylbenzylphthalate       | 0.638 | 0.639 | -0.2 | 111   | 0.00     |
| 81   | Benzo(a)anthracene         | 1.278 | 1.284 | -0.5 | 112   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 0.444 | 0.448 | -0.9 | 110   | 0.00     |
| 83   | Chrysene                   | 1.241 | 1.222 | 1.5  | 110   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 0.829 | 0.803 | 3.1  | 106   | 0.00     |
| 85 c | Di-n-octyl phthalate       | 1.316 | 1.215 | 7.7  | 100   | 0.00     |
| 86   | Indeno(1,2,3-cd)pyrene     | 1.042 | 1.020 | 2.1  | 114   | 0.00     |
| 87 I | Perylene-d12               | 1.000 | 1.000 | 0.0  | 107   | 0.00     |

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|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88   | Benzo(b)fluoranthene   | 1.156 | 1.131 | 2.2  | 99    | 0.00     |
| 89   | Benzo(k)fluoranthene   | 1.126 | 1.128 | -0.2 | 111   | 0.00     |
| 90 C | Benzo(a)pyrene         | 1.034 | 1.027 | 0.7  | 105   | 0.00     |
| 91   | Dibenzo(a,h)anthracene | 0.895 | 0.926 | -3.5 | 114   | 0.00     |
| 92   | Benzo(g,h,i)perylene   | 0.879 | 0.921 | -4.8 | 119   | 0.00     |

(#) = Out of Range

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