

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031220\  
 Data File : BG044762.D  
 Acq On : 13 Mar 2020 9:23  
 Operator : CG/JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTDCCC040

Quant Time: Mar 13 10:01:19 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG031020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Mar 10 16:27:39 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                    | AvqRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|-------|-------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0   | 116   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.619 | 0.541 | 12.6  | 102   | 0.00     |
| 3    | Pyridine                    | 1.832 | 1.644 | 10.3  | 103   | 0.00     |
| 4    | n-Nitrosodimethylamine      | 0.695 | 0.689 | 0.9   | 113   | 0.00     |
| 5 S  | 2-Fluorophenol              | 1.285 | 1.194 | 7.1   | 109   | 0.00     |
| 6    | Aniline                     | 2.323 | 2.102 | 9.5   | 106   | 0.00     |
| 7 S  | Phenol-d6                   | 1.867 | 1.733 | 7.2   | 110   | 0.00     |
| 8    | 2-Chlorophenol              | 1.283 | 1.165 | 9.2   | 109   | 0.00     |
| 9    | Benzaldehyde                | 1.089 | 1.018 | 6.5   | 106   | 0.00     |
| 10 C | Phenol                      | 1.848 | 1.689 | 8.6   | 109   | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 1.475 | 1.287 | 12.7  | 104   | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 1.580 | 1.435 | 9.2   | 108   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.562 | 1.405 | 10.1  | 107   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 1.481 | 1.322 | 10.7  | 107   | 0.00     |
| 15   | Benzyl Alcohol              | 1.498 | 1.394 | 6.9   | 109   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 2.603 | 2.364 | 9.2   | 107   | 0.00     |
| 17   | 2-Methylphenol              | 1.252 | 1.158 | 7.5   | 109   | 0.00     |
| 18   | Hexachloroethane            | 0.615 | 0.571 | 7.2   | 110   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 1.318 | 1.171 | 11.2  | 103   | 0.00     |
| 20   | 3+4-Methylphenols           | 1.726 | 1.597 | 7.5   | 107   | 0.00     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 115   | 0.00     |
| 22   | Acetophenone                | 0.567 | 0.496 | 12.5  | 106   | 0.00     |
| 23 S | Nitrobenzene-d5             | 0.456 | 0.425 | 6.8   | 112   | 0.00     |
| 24   | Nitrobenzene                | 0.473 | 0.429 | 9.3   | 110   | 0.00     |
| 25   | Isophorone                  | 0.890 | 0.781 | 12.2  | 106   | 0.00     |
| 26 C | 2-Nitrophenol               | 0.154 | 0.166 | -7.8  | 130   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 0.290 | 0.261 | 10.0  | 110   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 0.531 | 0.469 | 11.7  | 105   | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.338 | 0.318 | 5.9   | 112   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 0.405 | 0.371 | 8.4   | 112   | 0.00     |
| 31   | Naphthalene                 | 1.108 | 1.002 | 9.6   | 109   | 0.00     |
| 32   | Benzoic acid                | 0.201 | 0.149 | 25.9# | 94    | 0.00     |
| 33   | 4-Chloroaniline             | 0.499 | 0.444 | 11.0  | 108   | 0.00     |
| 34 C | Hexachlorobutadiene         | 0.266 | 0.248 | 6.8   | 112   | 0.00     |
| 35   | Caprolactam                 | 0.128 | 0.122 | 4.7   | 110   | 0.00     |
| 36 C | 4-Chloro-3-methylphenol     | 0.404 | 0.375 | 7.2   | 110   | 0.00     |
| 37   | 2-Methylnaphthalene         | 0.829 | 0.745 | 10.1  | 107   | 0.00     |
| 38   | 1-Methylnaphthalene         | 0.779 | 0.702 | 9.9   | 107   | 0.00     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 114   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.659 | 0.599 | 9.1   | 110   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 0.360 | 0.333 | 7.5   | 113   | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 0.262 | 0.258 | 1.5   | 118   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 0.437 | 0.431 | 1.4   | 118   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 0.453 | 0.433 | 4.4   | 116   | 0.00     |

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|      | Compound                   | AvqRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|----------------------------|-------|-------|-------|-------|----------|
| 45 S | 2-Fluorobiphenyl           | 1.397 | 1.270 | 9.1   | 108   | 0.00     |
| 46   | 1,1'-Biphenyl              | 1.598 | 1.446 | 9.5   | 108   | 0.00     |
| 47   | 2-Chloronaphthalene        | 1.221 | 1.113 | 8.8   | 110   | 0.00     |
| 48   | 2-Nitroaniline             | 0.427 | 0.439 | -2.8  | 123   | 0.00     |
| 49   | Acenaphthylene             | 1.913 | 1.733 | 9.4   | 108   | 0.00     |
| 50   | Dimethylphthalate          | 1.593 | 1.448 | 9.1   | 109   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 0.321 | 0.317 | 1.2   | 116   | 0.00     |
| 52 C | Acenaphthene               | 1.253 | 1.122 | 10.5  | 109   | 0.00     |
| 53   | 3-Nitroaniline             | 0.369 | 0.354 | 4.1   | 112   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 0.143 | 0.082 | 42.7# | 72    | 0.00     |
| 55   | Dibenzofuran               | 1.817 | 1.667 | 8.3   | 110   | 0.00     |
| 56 P | 4-Nitrophenol              | 0.282 | 0.268 | 5.0   | 114   | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 0.440 | 0.447 | -1.6  | 120   | 0.00     |
| 58   | Fluorene                   | 1.494 | 1.365 | 8.6   | 110   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.417 | 0.410 | 1.7   | 116   | 0.00     |
| 60   | Diethylphthalate           | 1.661 | 1.525 | 8.2   | 109   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 0.805 | 0.749 | 7.0   | 113   | 0.00     |
| 62   | 4-Nitroaniline             | 0.394 | 0.373 | 5.3   | 113   | 0.00     |
| 63   | Azobenzene                 | 1.725 | 1.568 | 9.1   | 109   | 0.00     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 118   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 0.087 | 0.069 | 20.7  | 102   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 0.602 | 0.534 | 11.3  | 110   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 0.250 | 0.225 | 10.0  | 114   | 0.00     |
| 68   | Hexachlorobenzene          | 0.271 | 0.246 | 9.2   | 115   | 0.00     |
| 69   | Atrazine                   | 0.221 | 0.203 | 8.1   | 109   | 0.00     |
| 70 C | Pentachlorophenol          | 0.149 | 0.131 | 12.1  | 106   | 0.00     |
| 71   | Phenanthrene               | 1.069 | 0.959 | 10.3  | 111   | 0.00     |
| 72   | Anthracene                 | 1.054 | 0.946 | 10.2  | 111   | 0.00     |
| 73   | Carbazole                  | 1.012 | 0.897 | 11.4  | 109   | 0.00     |
| 74   | Di-n-butylphthalate        | 1.231 | 1.103 | 10.4  | 108   | 0.00     |
| 75 C | Fluoranthene               | 1.273 | 1.151 | 9.6   | 111   | 0.00     |
| 76 I | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 119   | 0.00     |
| 77   | Benzidine                  | 0.649 | 0.635 | 2.2   | 108   | 0.00     |
| 78   | Pyrene                     | 1.467 | 1.316 | 10.3  | 110   | 0.00     |
| 79 S | Terphenyl-d14              | 1.069 | 0.937 | 12.3  | 110   | 0.00     |
| 80   | Butylbenzylphthalate       | 0.653 | 0.606 | 7.2   | 113   | 0.00     |
| 81   | Benzo(a)anthracene         | 1.440 | 1.304 | 9.4   | 112   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine     | 0.557 | 0.499 | 10.4  | 108   | 0.00     |
| 83   | Chrysene                   | 1.376 | 1.233 | 10.4  | 111   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate | 0.901 | 0.817 | 9.3   | 111   | 0.00     |
| 85 c | Di-n-octyl phthalate       | 1.508 | 1.399 | 7.2   | 112   | -0.01    |
| 86   | Indeno(1,2,3-cd)pyrene     | 1.804 | 1.657 | 8.1   | 115   | -0.02    |
| 87 I | Perylene-d12               | 1.000 | 1.000 | 0.0   | 118   | 0.00     |

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|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88   | Benzo(b)fluoranthene   | 1.320 | 1.187 | 10.1 | 110   | 0.00     |
| 89   | Benzo(k)fluoranthene   | 1.250 | 1.148 | 8.2  | 116   | -0.01    |
| 90 C | Benzo(a)pyrene         | 1.235 | 1.123 | 9.1  | 114   | -0.01    |
| 91   | Dibenzo(a,h)anthracene | 1.219 | 1.119 | 8.2  | 115   | -0.02    |
| 92   | Benzo(q,h,i)perylene   | 1.231 | 1.129 | 8.3  | 115   | -0.03    |

(#) = Out of Range

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