

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG093019\  
 Data File : BG042966.D  
 Acq On : 1 Oct 2019 7:02  
 Operator : JU  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTDCCC040

Quant Time: Oct 01 08:24:56 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092519.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Sep 25 15:35:52 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   | 120   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.467 | 0.475 | -1.7  | 133   | 0.00     |
| 3    | Pyridine                    | 1.314 | 1.307 | 0.5   | 121   | 0.00     |
| 4    | n-Nitrosodimethylamine      | 0.632 | 0.654 | -3.5  | 125   | 0.00     |
| 5 S  | 2-Fluorophenol              | 1.121 | 1.151 | -2.7  | 127   | 0.00     |
| 6    | Aniline                     | 1.948 | 1.904 | 2.3   | 118   | 0.00     |
| 7 S  | Phenol-d6                   | 1.614 | 1.646 | -2.0  | 122   | 0.00     |
| 8    | 2-Chlorophenol              | 1.169 | 1.172 | -0.3  | 121   | 0.00     |
| 9    | Benzaldehyde                | 0.986 | 0.939 | 4.8   | 106   | 0.00     |
| 10 C | Phenol                      | 1.481 | 1.513 | -2.2  | 124   | 0.00     |
| 11   | bis(2-Chloroethyl)ether     | 1.146 | 1.110 | 3.1   | 118   | 0.00     |
| 12   | 1,3-Dichlorobenzene         | 1.491 | 1.491 | 0.0   | 123   | 0.00     |
| 13 C | 1,4-Dichlorobenzene         | 1.486 | 1.510 | -1.6  | 124   | 0.00     |
| 14   | 1,2-Dichlorobenzene         | 1.415 | 1.399 | 1.1   | 122   | 0.00     |
| 15   | Benzyl Alcohol              | 1.213 | 1.237 | -2.0  | 121   | 0.00     |
| 16   | 2,2'-oxybis(1-Chloropropane | 2.200 | 2.011 | 8.6   | 110   | 0.00     |
| 17   | 2-Methylphenol              | 1.036 | 1.022 | 1.4   | 116   | 0.00     |
| 18   | Hexachloroethane            | 0.560 | 0.543 | 3.0   | 120   | 0.00     |
| 19 P | n-Nitroso-di-n-propylamine  | 1.060 | 1.053 | 0.7   | 118   | 0.00     |
| 20   | 3+4-Methylphenols           | 1.420 | 1.456 | -2.5  | 122   | 0.00     |
| 21 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 118   | 0.00     |
| 22   | Acetophenone                | 0.516 | 0.508 | 1.6   | 107   | 0.00     |
| 23 S | Nitrobenzene-d5             | 0.411 | 0.406 | 1.2   | 117   | 0.00     |
| 24   | Nitrobenzene                | 0.392 | 0.394 | -0.5  | 120   | 0.00     |
| 25   | Isophorone                  | 0.723 | 0.704 | 2.6   | 115   | 0.00     |
| 26 C | 2-Nitrophenol               | 0.167 | 0.176 | -5.4  | 128   | 0.00     |
| 27   | 2,4-Dimethylphenol          | 0.257 | 0.253 | 1.6   | 119   | 0.00     |
| 28   | bis(2-Chloroethoxy)methane  | 0.410 | 0.397 | 3.2   | 114   | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.319 | 0.327 | -2.5  | 120   | 0.00     |
| 30   | 1,2,4-Trichlorobenzene      | 0.412 | 0.412 | 0.0   | 121   | 0.00     |
| 31   | Naphthalene                 | 0.985 | 0.968 | 1.7   | 119   | 0.00     |
| 32   | Benzoic acid                | 0.194 | 0.214 | -10.3 | 110   | 0.00     |
| 33   | 4-Chloroaniline             | 0.427 | 0.426 | 0.2   | 116   | 0.00     |
| 34 C | Hexachlorobutadiene         | 0.309 | 0.309 | 0.0   | 122   | 0.00     |
| 35   | Caprolactam                 | 0.113 | 0.107 | 5.3   | 112   | -0.01    |
| 36 C | 4-Chloro-3-methylphenol     | 0.347 | 0.346 | 0.3   | 116   | 0.00     |
| 37   | 2-Methylnaphthalene         | 0.751 | 0.742 | 1.2   | 117   | 0.00     |
| 38   | 1-Methylnaphthalene         | 0.711 | 0.692 | 2.7   | 118   | 0.00     |
| 39 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 113   | 0.00     |
| 40   | 1,2,4,5-Tetrachlorobenzene  | 0.752 | 0.732 | 2.7   | 100   | 0.00     |
| 41 P | Hexachlorocyclopentadiene   | 0.479 | 0.448 | 6.5   | 109   | 0.00     |
| 42 S | 2,4,6-Tribromophenol        | 0.344 | 0.344 | 0.0   | 119   | 0.00     |
| 43 C | 2,4,6-Trichlorophenol       | 0.440 | 0.442 | -0.5  | 117   | 0.00     |
| 44   | 2,4,5-Trichlorophenol       | 0.453 | 0.459 | -1.3  | 119   | 0.00     |

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|      | Compound                   | AvqRF | CCRF  | %Dev | Area% | Dev(min) |
|------|----------------------------|-------|-------|------|-------|----------|
| 45 S | 2-Fluorobiphenyl           | 1.380 | 1.366 | 1.0  | 115   | 0.00     |
| 46   | 1,1'-Biphenyl              | 1.517 | 1.453 | 4.2  | 102   | 0.00     |
| 47   | 2-Chloronaphthalene        | 1.137 | 1.107 | 2.6  | 114   | 0.00     |
| 48   | 2-Nitroaniline             | 0.378 | 0.376 | 0.5  | 116   | 0.00     |
| 49   | Acenaphthylene             | 1.740 | 1.713 | 1.6  | 116   | 0.00     |
| 50   | Dimethylphthalate          | 1.499 | 1.426 | 4.9  | 112   | 0.00     |
| 51   | 2,6-Dinitrotoluene         | 0.319 | 0.319 | 0.0  | 116   | 0.00     |
| 52 C | Acenaphthene               | 1.165 | 1.160 | 0.4  | 113   | 0.00     |
| 53   | 3-Nitroaniline             | 0.330 | 0.319 | 3.3  | 113   | 0.00     |
| 54 P | 2,4-Dinitrophenol          | 0.198 | 0.192 | 3.0  | 113   | 0.00     |
| 55   | Dibenzofuran               | 1.723 | 1.661 | 3.6  | 113   | 0.00     |
| 56 P | 4-Nitrophenol              | 0.251 | 0.237 | 5.6  | 108   | 0.00     |
| 57   | 2,4-Dinitrotoluene         | 0.467 | 0.455 | 2.6  | 113   | 0.00     |
| 58   | Fluorene                   | 1.471 | 1.416 | 3.7  | 112   | 0.00     |
| 59   | 2,3,4,6-Tetrachlorophenol  | 0.483 | 0.473 | 2.1  | 115   | 0.00     |
| 60   | Diethylphthalate           | 1.525 | 1.437 | 5.8  | 110   | 0.00     |
| 61   | 4-Chlorophenyl-phenylether | 0.882 | 0.834 | 5.4  | 112   | 0.00     |
| 62   | 4-Nitroaniline             | 0.360 | 0.347 | 3.6  | 115   | 0.00     |
| 63   | Azobenzene                 | 1.388 | 1.297 | 6.6  | 109   | 0.00     |
| 64 I | Phenanthrene-d10           | 1.000 | 1.000 | 0.0  | 111   | 0.00     |
| 65   | 4,6-Dinitro-2-methylphenol | 0.113 | 0.113 | 0.0  | 108   | 0.00     |
| 66 c | n-Nitrosodiphenylamine     | 0.528 | 0.540 | -2.3 | 113   | 0.00     |
| 67   | 4-Bromophenyl-phenylether  | 0.251 | 0.256 | -2.0 | 111   | 0.00     |
| 68   | Hexachlorobenzene          | 0.279 | 0.288 | -3.2 | 114   | 0.00     |
| 69   | Atrazine                   | 0.222 | 0.216 | 2.7  | 104   | 0.00     |
| 70 C | Pentachlorophenol          | 0.161 | 0.168 | -4.3 | 114   | 0.00     |
| 71   | Phenanthrene               | 0.976 | 0.961 | 1.5  | 108   | 0.00     |
| 72   | Anthracene                 | 0.975 | 0.964 | 1.1  | 108   | 0.00     |
| 73   | Carbazole                  | 0.904 | 0.887 | 1.9  | 109   | 0.00     |
| 74   | Di-n-butylphthalate        | 1.078 | 1.054 | 2.2  | 107   | 0.00     |
| 75 C | Fluoranthene               | 1.291 | 1.242 | 3.8  | 106   | 0.00     |
| 76 I | Chrysene-d12               | 1.000 | 1.000 | 0.0  | 106   | 0.00     |
| 77   | Benzidine                  | 0.501 | 0.479 | 4.4  | 96    | -0.01    |
| 78   | Pyrene                     | 1.194 | 1.212 | -1.5 | 106   | 0.00     |
| 79 S | Terphenyl-d14              | 0.939 | 0.991 | -5.5 | 106   | 0.00     |
| 80   | Butylbenzylphthalate       | 0.453 | 0.456 | -0.7 | 107   | -0.01    |
| 81   | Benzo(a)anthracene         | 1.246 | 1.251 | -0.4 | 107   | -0.01    |
| 82   | 3,3'-Dichlorobenzidine     | 0.489 | 0.497 | -1.6 | 107   | 0.00     |
| 83   | Chrysene                   | 1.209 | 1.198 | 0.9  | 106   | -0.01    |
| 84   | Bis(2-ethylhexyl)phthalate | 0.645 | 0.649 | -0.6 | 109   | -0.01    |
| 85 c | Di-n-octyl phthalate       | 1.054 | 1.076 | -2.1 | 110   | -0.02    |
| 86   | Indeno(1,2,3-cd)pyrene     | 1.506 | 1.517 | -0.7 | 109   | -0.05    |
| 87 I | Perylene-d12               | 1.000 | 1.000 | 0.0  | 108   | -0.03    |

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|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88   | Benzo(b)fluoranthene   | 1.132 | 1.126 | 0.5  | 109   | -0.02    |
| 89   | Benzo(k)fluoranthene   | 1.120 | 1.101 | 1.7  | 107   | -0.02    |
| 90 C | Benzo(a)pyrene         | 1.068 | 1.065 | 0.3  | 109   | -0.03    |
| 91   | Dibenzo(a,h)anthracene | 1.095 | 1.093 | 0.2  | 109   | -0.05    |
| 92   | Benzo(g,h,i)perylene   | 1.061 | 1.056 | 0.5  | 110   | -0.05    |

(#) = Out of Range

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