

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN101123\  
 Data File : BN028247.D  
 Acq On : 11 Oct 2023 23:51  
 Operator : MA/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Oct 12 04:07:45 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN100223.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Oct 12 04:03:12 2023  
 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   | 56    | -0.04    |
| 2       | 1,4-Dioxane                | 0.449 | 0.491 | -9.4  | 60    | -0.03    |
| 3       | n-Nitrosodimethylamine     | 0.514 | 0.533 | -3.7  | 58    | -0.02    |
| 4 S     | 2-Fluorophenol             | 1.193 | 1.165 | 2.3   | 58    | -0.03    |
| 5 S     | Phenol-d6                  | 1.506 | 1.455 | 3.4   | 59    | -0.02    |
| 6       | bis(2-Chloroethyl)ether    | 1.212 | 1.100 | 9.2   | 51    | -0.03    |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0   | 58    | -0.03    |
| 8 S     | Nitrobenzene-d5            | 0.340 | 0.318 | 6.5   | 57    | -0.02    |
| 9       | Naphthalene                | 1.034 | 1.048 | -1.4  | 59    | -0.03    |
| 10      | Hexachlorobutadiene        | 0.184 | 0.174 | 5.4   | 55    | -0.04    |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.593 | 0.616 | -3.9  | 61    | -0.03    |
| 12      | 2-Methylnaphthalene        | 0.723 | 0.760 | -5.1  | 60    | -0.03    |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0   | 64    | -0.03    |
| 14 S    | 2,4,6-Tribromophenol       | 0.183 | 0.181 | 1.1   | 68    | -0.01    |
| 15 S    | 2-Fluorobiphenyl           | 1.442 | 1.346 | 6.7   | 61    | -0.03    |
| 16      | Acenaphthylene             | 1.772 | 1.748 | 1.4   | 63    | -0.03    |
| 17      | Acenaphthene               | 1.130 | 1.168 | -3.4  | 65    | -0.02    |
| 18      | Fluorene                   | 1.467 | 1.556 | -6.1  | 66    | -0.02    |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 68    | -0.02    |
| 20      | 4,6-Dinitro-2-methylphenol | 0.053 | 0.005 | 90.6# | 9#    | 1.04#    |
| 21      | 4-Bromophenyl-phenylether  | 0.209 | 0.200 | 4.3   | 65    | -0.01    |
| 22      | Hexachlorobenzene          | 0.219 | 0.212 | 3.2   | 64    | -0.02    |
| 23      | Atrazine                   | 0.182 | 0.177 | 2.7   | 66    | -0.01    |
| 24      | Pentachlorophenol          | 0.103 | 0.077 | 25.2# | 60    | -0.01    |
| 25      | Phenanthrene               | 1.078 | 1.106 | -2.6  | 69    | -0.02    |
| 26      | Anthracene                 | 1.019 | 1.011 | 0.8   | 67    | -0.02    |
| 27 SURR | Fluoranthene-d10           | 1.005 | 1.100 | -9.5  | 71    | -0.02    |
| 28      | Fluoranthene               | 1.245 | 1.402 | -12.6 | 73    | -0.02    |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 77    | -0.02    |
| 30      | Pyrene                     | 1.467 | 1.367 | 6.8   | 73    | -0.02    |
| 31 S    | Terphenyl-d14              | 0.933 | 0.799 | 14.4  | 70    | -0.02    |
| 32      | Benzo(a)anthracene         | 1.361 | 1.424 | -4.6  | 79    | -0.02    |
| 33      | Chrysene                   | 1.312 | 1.421 | -8.3  | 82    | -0.02    |
| 34      | Bis(2-ethylhexyl)phthalate | 0.895 | 0.953 | -6.5  | 80    | -0.02    |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0   | 96    | -0.02    |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.345 | 1.508 | -12.1 | 117   | -0.03    |
| 37      | Benzo(b)fluoranthene       | 1.292 | 1.257 | 2.7   | 93    | -0.02    |
| 38      | Benzo(k)fluoranthene       | 1.310 | 1.259 | 3.9   | 92    | -0.02    |
| 39 C    | Benzo(a)pyrene             | 1.222 | 1.221 | 0.1   | 97    | -0.02    |
| 40      | Dibenzo(a,h)anthracene     | 1.101 | 1.210 | -9.9  | 115   | -0.03    |
| 41      | Benzo(g,h,i)perylene       | 1.130 | 1.286 | -13.8 | 121   | -0.03    |

(#) = Out of Range

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

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