

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091423\  
 Data File : BN027686.D  
 Acq On : 14 Sep 2023 18:50  
 Operator : MA/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 52 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Sep 14 19:32:31 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN082923.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Sep 12 02:45:58 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   | 133   | 0.00     |
| 2       | 1,4-Dioxane                | 0.488 | 0.496 | -1.6  | 127   | 0.00     |
| 3       | n-Nitrosodimethylamine     | 0.518 | 0.555 | -7.1  | 145   | 0.00     |
| 4 S     | 2-Fluorophenol             | 1.043 | 1.128 | -8.1  | 147   | -0.01    |
| 5 S     | Phenol-d6                  | 1.268 | 1.270 | -0.2  | 136   | 0.00     |
| 6       | bis(2-Chloroethyl)ether    | 1.226 | 1.230 | -0.3  | 132   | 0.00     |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0   | 136   | -0.01    |
| 8 S     | Nitrobenzene-d5            | 0.292 | 0.325 | -11.3 | 145   | -0.01    |
| 9       | Naphthalene                | 1.091 | 1.105 | -1.3  | 139   | -0.01    |
| 10      | Hexachlorobutadiene        | 0.203 | 0.208 | -2.5  | 129   | -0.01    |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.587 | 0.513 | 12.6  | 115   | -0.01    |
| 12      | 2-Methylnaphthalene        | 0.776 | 0.670 | 13.7  | 113   | 0.00     |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0   | 96    | -0.01    |
| 14 S    | 2,4,6-Tribromophenol       | 0.148 | 0.114 | 23.0  | 80    | 0.00     |
| 15 S    | 2-Fluorobiphenyl           | 1.521 | 1.619 | -6.4  | 101   | -0.01    |
| 16      | Acenaphthylene             | 1.821 | 1.664 | 8.6   | 89    | -0.01    |
| 17      | Acenaphthene               | 1.215 | 1.211 | 0.3   | 94    | -0.01    |
| 18      | Fluorene                   | 1.620 | 1.488 | 8.1   | 85    | -0.01    |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 81    | 0.00     |
| 20      | 4,6-Dinitro-2-methylphenol | 0.006 | 0.005 | 16.7  | 85    | 0.00     |
| 21      | 4-Bromophenyl-phenylether  | 0.211 | 0.209 | 0.9   | 83    | -0.01    |
| 22      | Hexachlorobenzene          | 0.250 | 0.251 | -0.4  | 82    | -0.01    |
| 23      | Atrazine                   | 0.158 | 0.140 | 11.4  | 76    | 0.00     |
| 24      | Pentachlorophenol          | 0.096 | 0.075 | 21.9  | 70    | -0.01    |
| 25      | Phenanthrene               | 1.176 | 1.126 | 4.3   | 78    | -0.01    |
| 26      | Anthracene                 | 0.999 | 0.917 | 8.2   | 78    | 0.00     |
| 27 SURR | Fluoranthene-d10           | 0.994 | 0.846 | 14.9  | 69    | 0.00     |
| 28      | Fluoranthene               | 1.384 | 1.160 | 16.2  | 68    | 0.00     |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 62    | 0.00     |
| 30      | Pyrene                     | 1.590 | 1.707 | -7.4  | 66    | 0.00     |
| 31 S    | Terphenyl-d14              | 0.804 | 0.836 | -4.0  | 68    | 0.00     |
| 32      | Benzo(a)anthracene         | 1.371 | 1.322 | 3.6   | 62    | 0.00     |
| 33      | Chrysene                   | 1.507 | 1.519 | -0.8  | 60    | 0.00     |
| 34      | Bis(2-ethylhexyl)phthalate | 0.662 | 0.558 | 15.7  | 52    | 0.00     |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0   | 56    | -0.01    |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.624 | 1.606 | 1.1   | 63    | 0.00     |
| 37      | Benzo(b)fluoranthene       | 1.548 | 1.517 | 2.0   | 59    | 0.00     |
| 38      | Benzo(k)fluoranthene       | 1.591 | 1.518 | 4.6   | 58    | 0.00     |
| 39 C    | Benzo(a)pyrene             | 1.447 | 1.289 | 10.9  | 53    | 0.00     |
| 40      | Dibenzo(a,h)anthracene     | 1.273 | 1.296 | -1.8  | 65    | -0.01    |
| 41      | Benzo(g,h,i)perylene       | 1.476 | 1.509 | -2.2  | 62    | -0.02    |

(#) = Out of Range

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

**Instrument :**  
BNA\_N  
**LabSampleId :**  
SSTDCCC0.4