

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN090924\  
 Data File : BN033811.D  
 Acq On : 09 Sep 2024 10:36  
 Operator : MA/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Sep 10 03:14:10 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN090424.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Sep 04 18:36:01 2024  
 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  | 86    | 0.00     |
| 2       | 1,4-Dioxane                | 0.455 | 0.424 | 6.8  | 84    | 0.00     |
| 3       | n-Nitrosodimethylamine     | 0.534 | 0.537 | -0.6 | 90    | 0.00     |
| 4 S     | 2-Fluorophenol             | 1.225 | 1.079 | 11.9 | 80    | 0.00     |
| 5 S     | Phenol-d6                  | 1.450 | 1.259 | 13.2 | 80    | 0.00     |
| 6       | bis(2-Chloroethyl)ether    | 1.094 | 1.036 | 5.3  | 84    | -0.01    |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0  | 85    | -0.01    |
| 8 S     | Nitrobenzene-d5            | 0.341 | 0.317 | 7.0  | 84    | -0.01    |
| 9       | Naphthalene                | 1.068 | 1.031 | 3.5  | 85    | -0.01    |
| 10      | Hexachlorobutadiene        | 0.222 | 0.223 | -0.5 | 87    | -0.01    |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.566 | 0.529 | 6.5  | 83    | -0.01    |
| 12      | 2-Methylnaphthalene        | 0.655 | 0.613 | 6.4  | 84    | -0.01    |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0  | 83    | -0.01    |
| 14 S    | 2,4,6-Tribromophenol       | 0.202 | 0.170 | 15.8 | 76    | 0.00     |
| 15 S    | 2-Fluorobiphenyl           | 1.668 | 1.652 | 1.0  | 84    | -0.01    |
| 16      | Acenaphthylene             | 1.729 | 1.587 | 8.2  | 82    | -0.01    |
| 17      | Acenaphthene               | 1.185 | 1.142 | 3.6  | 83    | 0.00     |
| 18      | Fluorene                   | 1.462 | 1.406 | 3.8  | 84    | 0.00     |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0  | 86    | -0.01    |
| 20      | 4,6-Dinitro-2-methylphenol | 0.063 | 0.055 | 12.7 | 93    | 0.00     |
| 21      | 4-Bromophenyl-phenylether  | 0.232 | 0.219 | 5.6  | 83    | -0.01    |
| 22      | Hexachlorobenzene          | 0.266 | 0.263 | 1.1  | 85    | -0.01    |
| 23      | Atrazine                   | 0.185 | 0.166 | 10.3 | 80    | -0.01    |
| 24      | Pentachlorophenol          | 0.109 | 0.106 | 2.8  | 93    | -0.01    |
| 25      | Phenanthrene               | 1.102 | 1.064 | 3.4  | 85    | 0.00     |
| 26      | Anthracene                 | 0.958 | 0.886 | 7.5  | 83    | -0.01    |
| 27 SURR | Fluoranthene-d10           | 0.987 | 0.927 | 6.1  | 84    | 0.00     |
| 28      | Fluoranthene               | 1.235 | 1.162 | 5.9  | 85    | 0.00     |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0  | 88    | 0.00     |
| 30      | Pyrene                     | 1.612 | 1.582 | 1.9  | 85    | 0.00     |
| 31 S    | Terphenyl-d14              | 0.853 | 0.840 | 1.5  | 85    | 0.00     |
| 32      | Benzo(a)anthracene         | 1.428 | 1.353 | 5.3  | 84    | 0.00     |
| 33      | Chrysene                   | 1.412 | 1.463 | -3.6 | 91    | 0.00     |
| 34      | Bis(2-ethylhexyl)phthalate | 0.779 | 0.696 | 10.7 | 79    | 0.00     |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0  | 90    | 0.00     |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.653 | 1.621 | 1.9  | 89    | 0.00     |
| 37      | Benzo(b)fluoranthene       | 1.470 | 1.460 | 0.7  | 89    | 0.00     |
| 38      | Benzo(k)fluoranthene       | 1.433 | 1.470 | -2.6 | 94    | 0.00     |
| 39 C    | Benzo(a)pyrene             | 1.229 | 1.184 | 3.7  | 89    | 0.00     |
| 40      | Dibenzo(a,h)anthracene     | 1.325 | 1.268 | 4.3  | 87    | 0.00     |
| 41      | Benzo(g,h,i)perylene       | 1.437 | 1.422 | 1.0  | 88    | 0.00     |

(#) = Out of Range

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

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