

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN062722\  
 Data File : BN020501.D  
 Acq On : 27 Jun 2022 15:27  
 Operator : CG/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Jun 28 01:28:27 2022  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\8270-SIM-BN062622.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jun 27 02:03:21 2022  
 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  | 92    | -0.09    |
| 2       | 1,4-Dioxane                | 0.524 | 0.495 | 5.5  | 91    | -0.03    |
| 3       | n-Nitrosodimethylamine     | 0.478 | 0.420 | 12.1 | 86    | -0.04    |
| 4 S     | 2-Fluorophenol             | 1.110 | 1.162 | -4.7 | 98    | -0.06    |
| 5 S     | Phenol-d6                  | 1.327 | 1.433 | -8.0 | 104   | -0.09    |
| 6       | bis(2-Chloroethyl)ether    | 1.142 | 1.154 | -1.1 | 98    | -0.09    |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0  | 98    | -0.11    |
| 8 S     | Nitrobenzene-d5            | 0.324 | 0.302 | 6.8  | 100   | -0.10    |
| 9       | Naphthalene                | 1.195 | 1.180 | 1.3  | 101   | -0.11    |
| 10      | Hexachlorobutadiene        | 0.218 | 0.213 | 2.3  | 99    | -0.10    |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.686 | 0.682 | 0.6  | 103   | -0.11    |
| 12      | 2-Methylnaphthalene        | 0.794 | 0.787 | 0.9  | 102   | -0.11    |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0  | 104   | -0.11    |
| 14 S    | 2,4,6-Tribromophenol       | 0.151 | 0.143 | 5.3  | 114   | -0.11    |
| 15 S    | 2-Fluorobiphenyl           | 1.651 | 1.591 | 3.6  | 102   | -0.11    |
| 16      | Acenaphthylene             | 1.923 | 1.827 | 5.0  | 107   | -0.11    |
| 17      | Acenaphthene               | 1.308 | 1.272 | 2.8  | 104   | -0.11    |
| 18      | Fluorene                   | 1.707 | 1.667 | 2.3  | 106   | -0.12    |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0  | 109   | -0.11    |
| 20      | 4,6-Dinitro-2-methylphenol | 0.051 | 0.039 | 23.5 | 113   | -0.12    |
| 21      | 4-Bromophenyl-phenylether  | 0.233 | 0.216 | 7.3  | 109   | -0.11    |
| 22      | Hexachlorobenzene          | 0.257 | 0.239 | 7.0  | 109   | -0.11    |
| 23      | Atrazine                   | 0.183 | 0.156 | 14.8 | 103   | -0.11    |
| 24      | Pentachlorophenol          | 0.061 | 0.066 | -8.2 | 160#  | -0.11    |
| 25      | Phenanthrene               | 1.349 | 1.246 | 7.6  | 108   | -0.11    |
| 26      | Anthracene                 | 1.163 | 1.065 | 8.4  | 110   | -0.11    |
| 27 SURR | Fluoranthene-d10           | 1.188 | 1.038 | 12.6 | 103   | -0.12    |
| 28      | Fluoranthene               | 1.471 | 1.294 | 12.0 | 104   | -0.12    |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0  | 92    | -0.13    |
| 30      | Pyrene                     | 1.694 | 1.818 | -7.3 | 101   | -0.12    |
| 31 S    | Terphenyl-d14              | 0.956 | 1.026 | -7.3 | 101   | -0.13    |
| 32      | Benzo(a)anthracene         | 1.429 | 1.385 | 3.1  | 95    | -0.13    |
| 33      | Chrysene                   | 1.582 | 1.535 | 3.0  | 93    | -0.13    |
| 34      | Bis(2-ethylhexyl)phthalate | 0.805 | 0.796 | 1.1  | 103   | -0.13    |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0  | 76    | -0.20    |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.782 | 1.504 | 15.6 | 68    | -0.30    |
| 37      | Benzo(b)fluoranthene       | 1.541 | 1.545 | -0.3 | 80    | -0.18    |
| 38      | Benzo(k)fluoranthene       | 1.641 | 1.572 | 4.2  | 78    | -0.18    |
| 39 C    | Benzo(a)pyrene             | 1.337 | 1.275 | 4.6  | 76    | -0.20    |
| 40      | Dibenzo(a,h)anthracene     | 1.429 | 1.214 | 15.0 | 69    | -0.32    |
| 41      | Benzo(g,h,i)perylene       | 1.561 | 1.336 | 14.4 | 67    | -0.34    |

(#) = Out of Range

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

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