

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061925\  
 Data File : BN037338.D  
 Acq On : 19 Jun 2025 07:06  
 Operator : RC/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Jun 19 08:40:04 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN061925.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Jun 18 18:40:16 2025  
 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   | 173#  | 0.00     |
| 2       | 1,4-Dioxane                | 0.357 | 0.318 | 10.9  | 137   | 0.00     |
| 3       | n-Nitrosodimethylamine     | 0.414 | 0.450 | -8.7  | 193#  | 0.00     |
| 4 S     | 2-Fluorophenol             | 0.727 | 0.740 | -1.8  | 178#  | 0.00     |
| 5 S     | Phenol-d6                  | 0.715 | 0.739 | -3.4  | 182#  | -0.01    |
| 6       | bis(2-Chloroethyl)ether    | 0.612 | 0.683 | -11.6 | 212#  | 0.00     |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0   | 166#  | -0.01    |
| 8 S     | Nitrobenzene-d5            | 0.333 | 0.352 | -5.7  | 178#  | -0.01    |
| 9       | Naphthalene                | 1.035 | 1.046 | -1.1  | 168#  | 0.00     |
| 10      | Hexachlorobutadiene        | 0.524 | 0.530 | -1.1  | 157#  | -0.01    |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.650 | 0.706 | -8.6  | 170#  | 0.00     |
| 12      | 2-Methylnaphthalene        | 0.729 | 0.748 | -2.6  | 175#  | -0.01    |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0   | 175#  | -0.01    |
| 14 S    | 2,4,6-Tribromophenol       | 0.295 | 0.257 | 12.9  | 155#  | 0.00     |
| 15 S    | 2-Fluorobiphenyl           | 1.753 | 1.672 | 4.6   | 166#  | 0.00     |
| 16      | Acenaphthylene             | 1.683 | 1.485 | 11.8  | 160#  | 0.00     |
| 17      | Acenaphthene               | 1.117 | 0.981 | 12.2  | 158#  | -0.01    |
| 18      | Fluorene                   | 1.584 | 1.410 | 11.0  | 160#  | -0.01    |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 164#  | 0.00     |
| 20      | 4,6-Dinitro-2-methylphenol | 0.123 | 0.112 | 8.9   | 195#  | -0.01    |
| 21      | 4-Bromophenyl-phenylether  | 0.332 | 0.314 | 5.4   | 163#  | -0.01    |
| 22      | Hexachlorobenzene          | 0.348 | 0.337 | 3.2   | 152#  | 0.00     |
| 23      | Atrazine                   | 0.259 | 0.250 | 3.5   | 167#  | 0.00     |
| 24      | Pentachlorophenol          | 0.193 | 0.180 | 6.7   | 173#  | -0.01    |
| 25      | Phenanthrene               | 1.126 | 1.094 | 2.8   | 167#  | 0.00     |
| 26      | Anthracene                 | 1.056 | 0.997 | 5.6   | 164#  | -0.01    |
| 27 SURR | Fluoranthene-d10           | 1.268 | 1.307 | -3.1  | 162#  | 0.00     |
| 28      | Fluoranthene               | 1.564 | 1.517 | 3.0   | 166#  | 0.00     |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 187#  | 0.00     |
| 30      | Pyrene                     | 1.523 | 1.468 | 3.6   | 171#  | 0.00     |
| 31 S    | Terphenyl-d14              | 0.918 | 0.879 | 4.2   | 172#  | 0.00     |
| 32      | Benzo(a)anthracene         | 1.288 | 1.299 | -0.9  | 203#  | 0.00     |
| 33      | Chrysene                   | 1.644 | 1.568 | 4.6   | 178#  | 0.00     |
| 34      | Bis(2-ethylhexyl)phthalate | 0.528 | 0.512 | 3.0   | 179#  | 0.00     |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0   | 195#  | 0.00     |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.682 | 1.673 | 0.5   | 204#  | -0.01    |
| 37      | Benzo(b)fluoranthene       | 1.397 | 1.374 | 1.6   | 208#  | 0.00     |
| 38      | Benzo(k)fluoranthene       | 1.535 | 1.374 | 10.5  | 176#  | -0.05    |
| 39 C    | Benzo(a)pyrene             | 1.294 | 1.246 | 3.7   | 196#  | 0.00     |
| 40      | Dibenzo(a,h)anthracene     | 1.351 | 1.304 | 3.5   | 205#  | -0.02    |
| 41      | Benzo(g,h,i)perylene       | 1.542 | 1.492 | 3.2   | 197#  | -0.01    |

(#) = Out of Range

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

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