

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100625\  
 Data File : BN038039.D  
 Acq On : 07 Oct 2025 09:47  
 Operator : RC/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 39 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Oct 07 10:39:40 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN092325.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Sep 24 11:00:01 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#  | -0.02    |
| 2       | 1,4-Dioxane                | 0.406 | 0.426 | -4.9  | 180#  | 0.00     |
| 3       | n-Nitrosodimethylamine     | 0.540 | 0.545 | -0.9  | 183#  | 0.00     |
| 4 S     | 2-Fluorophenol             | 0.913 | 0.850 | 6.9   | 172#  | -0.01    |
| 5 S     | Phenol-d6                  | 1.199 | 1.108 | 7.6   | 171#  | -0.01    |
| 6       | bis(2-Chloroethyl)ether    | 1.113 | 1.093 | 1.8   | 175#  | -0.02    |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0   | 178#  | -0.02    |
| 8 S     | Nitrobenzene-d5            | 0.305 | 0.291 | 4.6   | 179#  | -0.02    |
| 9       | Naphthalene                | 1.102 | 1.051 | 4.6   | 179#  | -0.02    |
| 10      | Hexachlorobutadiene        | 0.188 | 0.185 | 1.6   | 181#  | -0.02    |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.556 | 0.515 | 7.4   | 184#  | -0.02    |
| 12      | 2-Methylnaphthalene        | 0.720 | 0.679 | 5.7   | 181#  | -0.02    |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0   | 182#  | -0.02    |
| 14 S    | 2,4,6-Tribromophenol       | 0.152 | 0.117 | 23.0  | 157#  | -0.02    |
| 15 S    | 2-Fluorobiphenyl           | 1.638 | 1.521 | 7.1   | 178#  | -0.01    |
| 16      | Acenaphthylene             | 1.816 | 1.702 | 6.3   | 186#  | -0.02    |
| 17      | Acenaphthene               | 1.294 | 1.184 | 8.5   | 180#  | -0.02    |
| 18      | Fluorene                   | 1.565 | 1.333 | 14.8  | 167#  | -0.01    |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 175#  | -0.01    |
| 20      | 4,6-Dinitro-2-methylphenol | 0.057 | 0.032 | 43.9# | 128   | -0.01    |
| 21      | 4-Bromophenyl-phenylether  | 0.261 | 0.248 | 5.0   | 177#  | -0.01    |
| 22      | Hexachlorobenzene          | 0.316 | 0.307 | 2.8   | 176#  | -0.02    |
| 23      | Atrazine                   | 0.199 | 0.154 | 22.6  | 156#  | -0.02    |
| 24      | Pentachlorophenol          | 0.108 | 0.071 | 34.3# | 135   | -0.01    |
| 25      | Phenanthrene               | 1.316 | 1.222 | 7.1   | 173#  | -0.01    |
| 26      | Anthracene                 | 1.142 | 1.017 | 10.9  | 173#  | -0.02    |
| 27 SURR | Fluoranthene-d10           | 1.006 | 0.820 | 18.5  | 161#  | -0.01    |
| 28      | Fluoranthene               | 1.450 | 1.188 | 18.1  | 158#  | -0.02    |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 120   | 0.00     |
| 30      | Pyrene                     | 1.579 | 1.961 | -24.2 | 154#  | -0.01    |
| 31 S    | Terphenyl-d14              | 0.797 | 0.944 | -18.4 | 150   | -0.01    |
| 32      | Benzo(a)anthracene         | 1.398 | 1.322 | 5.4   | 122   | 0.00     |
| 33      | Chrysene                   | 1.512 | 1.478 | 2.2   | 121   | 0.00     |
| 34      | Bis(2-ethylhexyl)phthalate | 0.553 | 0.437 | 21.0  | 107   | 0.00     |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0   | 101   | -0.02    |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.828 | 1.860 | -1.8  | 113   | -0.04    |
| 37      | Benzo(b)fluoranthene       | 1.725 | 1.717 | 0.5   | 105   | -0.02    |
| 38      | Benzo(k)fluoranthene       | 1.736 | 1.646 | 5.2   | 104   | -0.02    |
| 39 C    | Benzo(a)pyrene             | 1.364 | 1.296 | 5.0   | 106   | -0.02    |
| 40      | Dibenzo(a,h)anthracene     | 1.427 | 1.456 | -2.0  | 112   | -0.03    |
| 41      | Benzo(g,h,i)perylene       | 1.499 | 1.563 | -4.3  | 112   | -0.03    |

(#) = Out of Range

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

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