

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN090524\  
 Data File : BN033760.D  
 Acq On : 05 Sep 2024 09:56  
 Operator : MA/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Sep 05 11:27:48 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                   | Amount | Calc. | %Dev | Area% | Dev(min) |
|---------|----------------------------|--------|-------|------|-------|----------|
| 1 I     | 1,4-Dichlorobenzene-d4     | 0.400  | 0.400 | 0.0  | 100   | 0.00     |
| 2       | 1,4-Dioxane                | 0.400  | 0.390 | 2.5  | 101   | 0.00     |
| 3       | n-Nitrosodimethylamine     | 0.400  | 0.371 | 7.3  | 97    | 0.00     |
| 4 S     | 2-Fluorophenol             | 0.400  | 0.353 | 11.8 | 92    | 0.00     |
| 5 S     | Phenol-d6                  | 0.400  | 0.347 | 13.3 | 93    | 0.00     |
| 6       | bis(2-Chloroethyl)ether    | 0.400  | 0.385 | 3.8  | 99    | 0.00     |
| 7 I     | Naphthalene-d8             | 0.400  | 0.400 | 0.0  | 100   | 0.00     |
| 8 S     | Nitrobenzene-d5            | 0.400  | 0.367 | 8.3  | 97    | 0.00     |
| 9       | Naphthalene                | 0.400  | 0.385 | 3.8  | 100   | 0.00     |
| 10      | Hexachlorobutadiene        | 0.400  | 0.390 | 2.5  | 100   | -0.01    |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.400  | 0.379 | 5.3  | 99    | 0.00     |
| 12      | 2-Methylnaphthalene        | 0.400  | 0.378 | 5.5  | 99    | 0.00     |
| 13 I    | Acenaphthene-d10           | 0.400  | 0.400 | 0.0  | 98    | -0.01    |
| 14 S    | 2,4,6-Tribromophenol       | 0.400  | 0.325 | 18.8 | 87    | 0.00     |
| 15 S    | 2-Fluorobiphenyl           | 0.400  | 0.397 | 0.8  | 100   | 0.00     |
| 16      | Acenaphthylene             | 0.400  | 0.371 | 7.3  | 98    | 0.00     |
| 17      | Acenaphthene               | 0.400  | 0.390 | 2.5  | 100   | 0.00     |
| 18      | Fluorene                   | 0.400  | 0.386 | 3.5  | 100   | 0.00     |
| 19 I    | Phenanthrene-d10           | 0.400  | 0.400 | 0.0  | 102   | 0.00     |
| 20      | 4,6-Dinitro-2-methylphenol | 0.400  | 0.328 | 18.0 | 104   | 0.00     |
| 21      | 4-Bromophenyl-phenylether  | 0.400  | 0.379 | 5.3  | 99    | 0.00     |
| 22      | Hexachlorobenzene          | 0.400  | 0.403 | -0.8 | 103   | -0.01    |
| 23      | Atrazine                   | 0.400  | 0.349 | 12.8 | 92    | -0.01    |
| 24      | Pentachlorophenol          | 0.400  | 0.342 | 14.5 | 96    | -0.01    |
| 25      | Phenanthrene               | 0.400  | 0.389 | 2.8  | 102   | 0.00     |
| 26      | Anthracene                 | 0.400  | 0.373 | 6.8  | 99    | -0.01    |
| 27 SURR | Fluoranthene-d10           | 0.400  | 0.383 | 4.3  | 102   | 0.00     |
| 28      | Fluoranthene               | 0.400  | 0.382 | 4.5  | 103   | 0.00     |
| 29 I    | Chrysene-d12               | 0.400  | 0.400 | 0.0  | 107   | 0.00     |
| 30      | Pyrene                     | 0.400  | 0.393 | 1.8  | 102   | 0.00     |
| 31 S    | Terphenyl-d14              | 0.400  | 0.394 | 1.5  | 103   | 0.00     |
| 32      | Benzo(a)anthracene         | 0.400  | 0.385 | 3.8  | 104   | 0.00     |
| 33      | Chrysene                   | 0.400  | 0.408 | -2.0 | 109   | 0.00     |
| 34      | Bis(2-ethylhexyl)phthalate | 0.400  | 0.338 | 15.5 | 91    | 0.00     |
| 35 I    | Perylene-d12               | 0.400  | 0.400 | 0.0  | 109   | 0.00     |
| 36      | Indeno(1,2,3-cd)pyrene     | 0.400  | 0.397 | 0.8  | 110   | 0.00     |
| 37      | Benzo(b)fluoranthene       | 0.400  | 0.402 | -0.5 | 110   | 0.00     |
| 38      | Benzo(k)fluoranthene       | 0.400  | 0.397 | 0.8  | 110   | 0.00     |
| 39 C    | Benzo(a)pyrene             | 0.400  | 0.382 | 4.5  | 107   | 0.00     |
| 40      | Dibenzo(a,h)anthracene     | 0.400  | 0.394 | 1.5  | 109   | 0.00     |
| 41      | Benzo(g,h,i)perylene       | 0.400  | 0.403 | -0.8 | 109   | 0.00     |

(#) = Out of Range

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

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