

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN022323\  
 Data File : BN024098.D  
 Acq On : 23 Feb 2023 13:39  
 Operator : CG/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Feb 24 02:18:05 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN020723.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Feb 23 03:52:37 2023  
 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   | 78    | 0.00     |
| 2       | 1,4-Dioxane                | 0.379 | 0.394 | -4.0  | 84    | 0.00     |
| 3       | n-Nitrosodimethylamine     | 0.425 | 0.384 | 9.6   | 79    | 0.00     |
| 4 S     | 2-Fluorophenol             | 0.831 | 0.809 | 2.6   | 81    | 0.00     |
| 5 S     | Phenol-d6                  | 0.980 | 0.974 | 0.6   | 85    | 0.00     |
| 6       | bis(2-Chloroethyl)ether    | 0.979 | 1.067 | -9.0  | 90    | 0.00     |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0   | 81    | 0.00     |
| 8 S     | Nitrobenzene-d5            | 0.321 | 0.316 | 1.6   | 92    | 0.00     |
| 9       | Naphthalene                | 0.998 | 1.027 | -2.9  | 93    | 0.00     |
| 10      | Hexachlorobutadiene        | 0.262 | 0.268 | -2.3  | 91    | 0.00     |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.556 | 0.533 | 4.1   | 87    | 0.00     |
| 12      | 2-Methylnaphthalene        | 0.664 | 0.662 | 0.3   | 90    | 0.00     |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0   | 73    | 0.00     |
| 14 S    | 2,4,6-Tribromophenol       | 0.206 | 0.150 | 27.2# | 61    | 0.00     |
| 15 S    | 2-Fluorobiphenyl           | 1.582 | 1.676 | -5.9  | 86    | -0.01    |
| 16      | Acenaphthylene             | 1.558 | 1.466 | 5.9   | 81    | 0.00     |
| 17      | Acenaphthene               | 1.060 | 1.099 | -3.7  | 85    | 0.00     |
| 18      | Fluorene                   | 1.438 | 1.386 | 3.6   | 76    | 0.00     |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 59    | 0.00     |
| 20      | 4,6-Dinitro-2-methylphenol | 0.065 | 0.045 | 30.8# | 53    | 0.00     |
| 21      | 4-Bromophenyl-phenylether  | 0.260 | 0.255 | 1.9   | 69    | 0.00     |
| 22      | Hexachlorobenzene          | 0.313 | 0.343 | -9.6  | 75    | 0.00     |
| 23      | Atrazine                   | 0.170 | 0.153 | 10.0  | 63    | 0.00     |
| 24      | Pentachlorophenol          | 0.111 | 0.087 | 21.6  | 58    | 0.00     |
| 25      | Phenanthrene               | 1.104 | 1.120 | -1.4  | 68    | 0.00     |
| 26      | Anthracene                 | 0.908 | 0.828 | 8.8   | 63    | -0.01    |
| 27 SURR | Fluoranthene-d10           | 0.971 | 0.867 | 10.7  | 57    | 0.00     |
| 28      | Fluoranthene               | 1.265 | 1.139 | 10.0  | 58    | 0.00     |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 42#   | 0.00     |
| 30      | Pyrene                     | 1.387 | 1.548 | -11.6 | 57    | 0.00     |
| 31 S    | Terphenyl-d14              | 0.746 | 0.781 | -4.7  | 52    | 0.00     |
| 32      | Benzo(a)anthracene         | 1.280 | 1.169 | 8.7   | 44#   | 0.00     |
| 33      | Chrysene                   | 1.381 | 1.439 | -4.2  | 50    | 0.00     |
| 34      | Bis(2-ethylhexyl)phthalate | 0.585 | 0.547 | 6.5   | 47#   | 0.00     |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0   | 39#   | 0.00     |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.818 | 1.746 | 4.0   | 46#   | 0.00     |
| 37      | Benzo(b)fluoranthene       | 1.619 | 1.640 | -1.3  | 46#   | 0.00     |
| 38      | Benzo(k)fluoranthene       | 1.600 | 1.614 | -0.9  | 45#   | -0.01    |
| 39 C    | Benzo(a)pyrene             | 1.228 | 1.169 | 4.8   | 44#   | 0.00     |
| 40      | Dibenzo(a,h)anthracene     | 1.455 | 1.412 | 3.0   | 46#   | 0.00     |
| 41      | Benzo(g,h,i)perylene       | 1.543 | 1.588 | -2.9  | 49#   | 0.01     |

(#) = Out of Range

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

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