

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN110624\  
 Data File : BN034869.D  
 Acq On : 06 Nov 2024 11:18  
 Operator : RC/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Nov 06 12:16:31 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN103024.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Nov 05 00:03:28 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                   | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|---------|----------------------------|-------|-------|-------|-------|----------|
| 1 I     | 1,4-Dichlorobenzene-d4     | 1.000 | 1.000 | 0.0   | 109   | -0.01    |
| 2       | 1,4-Dioxane                | 0.439 | 0.442 | -0.7  | 118   | 0.00     |
| 3       | n-Nitrosodimethylamine     | 0.494 | 0.513 | -3.8  | 128   | 0.00     |
| 4 S     | 2-Fluorophenol             | 1.212 | 1.189 | 1.9   | 118   | -0.01    |
| 5 S     | Phenol-d6                  | 1.574 | 1.544 | 1.9   | 120   | -0.01    |
| 6       | bis(2-Chloroethyl)ether    | 1.149 | 1.216 | -5.8  | 125   | -0.01    |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0   | 110   | -0.01    |
| 8 S     | Nitrobenzene-d5            | 0.314 | 0.321 | -2.2  | 127   | -0.02    |
| 9       | Naphthalene                | 1.106 | 1.028 | 7.1   | 114   | -0.02    |
| 10      | Hexachlorobutadiene        | 0.182 | 0.164 | 9.9   | 110   | -0.02    |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.579 | 0.516 | 10.9  | 110   | -0.02    |
| 12      | 2-Methylnaphthalene        | 0.723 | 0.637 | 11.9  | 108   | -0.02    |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0   | 99    | -0.01    |
| 14 S    | 2,4,6-Tribromophenol       | 0.198 | 0.164 | 17.2  | 103   | -0.01    |
| 15 S    | 2-Fluorobiphenyl           | 1.575 | 1.468 | 6.8   | 107   | -0.01    |
| 16      | Acenaphthylene             | 2.007 | 1.661 | 17.2  | 98    | -0.01    |
| 17      | Acenaphthene               | 1.337 | 1.187 | 11.2  | 104   | -0.02    |
| 18      | Fluorene                   | 1.667 | 1.484 | 11.0  | 103   | -0.01    |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 98    | -0.01    |
| 20      | 4,6-Dinitro-2-methylphenol | 0.059 | 0.048 | 18.6  | 106   | -0.01    |
| 21      | 4-Bromophenyl-phenylether  | 0.231 | 0.205 | 11.3  | 95    | -0.01    |
| 22      | Hexachlorobenzene          | 0.259 | 0.227 | 12.4  | 93    | -0.01    |
| 23      | Atrazine                   | 0.199 | 0.194 | 2.5   | 106   | -0.01    |
| 24      | Pentachlorophenol          | 0.112 | 0.120 | -7.1  | 129   | -0.01    |
| 25      | Phenanthrene               | 1.176 | 1.137 | 3.3   | 104   | -0.01    |
| 26      | Anthracene                 | 1.101 | 1.031 | 6.4   | 102   | -0.01    |
| 27 SURR | Fluoranthene-d10           | 0.938 | 0.941 | -0.3  | 109   | -0.01    |
| 28      | Fluoranthene               | 1.297 | 1.308 | -0.8  | 109   | -0.01    |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 113   | 0.00     |
| 30      | Pyrene                     | 2.091 | 1.914 | 8.5   | 112   | -0.01    |
| 31 S    | Terphenyl-d14              | 0.859 | 0.724 | 15.7  | 104   | -0.01    |
| 32      | Benzo(a)anthracene         | 1.592 | 1.624 | -2.0  | 130   | 0.00     |
| 33      | Chrysene                   | 1.559 | 1.577 | -1.2  | 126   | 0.00     |
| 34      | Bis(2-ethylhexyl)phthalate | 1.331 | 1.301 | 2.3   | 114   | 0.00     |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0   | 128   | -0.02    |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.471 | 1.660 | -12.8 | 159#  | -0.03    |
| 37      | Benzo(b)fluoranthene       | 1.610 | 1.664 | -3.4  | 146   | -0.02    |
| 38      | Benzo(k)fluoranthene       | 1.571 | 1.608 | -2.4  | 147   | -0.01    |
| 39 C    | Benzo(a)pyrene             | 1.329 | 1.339 | -0.8  | 145   | -0.02    |
| 40      | Dibenzo(a,h)anthracene     | 1.147 | 1.277 | -11.3 | 157#  | -0.04    |
| 41      | Benzo(g,h,i)perylene       | 1.240 | 1.395 | -12.5 | 156#  | -0.04    |

(#) = Out of Range

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

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