

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN082923\  
 Data File : BN027232.D  
 Acq On : 29 Aug 2023 09:21  
 Operator : MA/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Aug 30 00:02:39 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN082923.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Aug 29 05:37:30 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   | 109   | 0.00     |
| 2       | 1,4-Dioxane                | 0.488 | 0.470 | 3.7   | 99    | 0.00     |
| 3       | n-Nitrosodimethylamine     | 0.518 | 0.550 | -6.2  | 118   | 0.00     |
| 4 S     | 2-Fluorophenol             | 1.043 | 0.990 | 5.1   | 106   | 0.00     |
| 5 S     | Phenol-d6                  | 1.268 | 1.179 | 7.0   | 104   | 0.00     |
| 6       | bis(2-Chloroethyl)ether    | 1.226 | 1.276 | -4.1  | 112   | 0.00     |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0   | 110   | 0.00     |
| 8 S     | Nitrobenzene-d5            | 0.292 | 0.316 | -8.2  | 114   | 0.00     |
| 9       | Naphthalene                | 1.091 | 1.108 | -1.6  | 112   | -0.01    |
| 10      | Hexachlorobutadiene        | 0.203 | 0.223 | -9.9  | 111   | -0.01    |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.587 | 0.613 | -4.4  | 111   | 0.00     |
| 12      | 2-Methylnaphthalene        | 0.776 | 0.815 | -5.0  | 111   | 0.00     |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0   | 111   | 0.00     |
| 14 S    | 2,4,6-Tribromophenol       | 0.148 | 0.143 | 3.4   | 116   | 0.00     |
| 15 S    | 2-Fluorobiphenyl           | 1.521 | 1.555 | -2.2  | 113   | 0.00     |
| 16      | Acenaphthylene             | 1.821 | 1.795 | 1.4   | 112   | 0.00     |
| 17      | Acenaphthene               | 1.215 | 1.225 | -0.8  | 110   | 0.00     |
| 18      | Fluorene                   | 1.620 | 1.715 | -5.9  | 115   | 0.00     |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 112   | 0.00     |
| 20      | 4,6-Dinitro-2-methylphenol | 0.006 | 0.007 | -16.7 | 143   | 0.00     |
| 21      | 4-Bromophenyl-phenylether  | 0.211 | 0.215 | -1.9  | 118   | -0.01    |
| 22      | Hexachlorobenzene          | 0.250 | 0.256 | -2.4  | 116   | 0.00     |
| 23      | Atrazine                   | 0.158 | 0.158 | 0.0   | 119   | 0.00     |
| 24      | Pentachlorophenol          | 0.096 | 0.094 | 2.1   | 122   | -0.01    |
| 25      | Phenanthrene               | 1.176 | 1.198 | -1.9  | 116   | -0.01    |
| 26      | Anthracene                 | 0.999 | 0.991 | 0.8   | 118   | 0.00     |
| 27 SURR | Fluoranthene-d10           | 0.994 | 0.985 | 0.9   | 111   | 0.00     |
| 28      | Fluoranthene               | 1.384 | 1.364 | 1.4   | 112   | 0.00     |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 104   | 0.00     |
| 30      | Pyrene                     | 1.590 | 1.694 | -6.5  | 110   | 0.00     |
| 31 S    | Terphenyl-d14              | 0.804 | 0.804 | 0.0   | 110   | 0.00     |
| 32      | Benzo(a)anthracene         | 1.371 | 1.324 | 3.4   | 104   | 0.00     |
| 33      | Chrysene                   | 1.507 | 1.559 | -3.5  | 104   | 0.00     |
| 34      | Bis(2-ethylhexyl)phthalate | 0.662 | 0.562 | 15.1  | 87    | 0.00     |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0   | 97    | 0.00     |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.624 | 1.482 | 8.7   | 100   | 0.00     |
| 37      | Benzo(b)fluoranthene       | 1.548 | 1.545 | 0.2   | 103   | 0.00     |
| 38      | Benzo(k)fluoranthene       | 1.591 | 1.528 | 4.0   | 101   | 0.00     |
| 39 C    | Benzo(a)pyrene             | 1.447 | 1.350 | 6.7   | 96    | 0.00     |
| 40      | Dibenzo(a,h)anthracene     | 1.273 | 1.157 | 9.1   | 100   | 0.00     |
| 41      | Benzo(g,h,i)perylene       | 1.476 | 1.416 | 4.1   | 101   | 0.00     |

(#) = Out of Range

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

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