

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN120122\  
 Data File : BN023027.D  
 Acq On : 01 Dec 2022 20:25  
 Operator : CG/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Dec 02 01:46:58 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN112522.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Dec 02 01:41:37 2022  
 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   | 86    | 0.00     |
| 2       | 1,4-Dioxane                | 0.508 | 0.437 | 14.0  | 72    | 0.00     |
| 3       | n-Nitrosodimethylamine     | 0.547 | 0.450 | 17.7  | 69    | 0.00     |
| 4 S     | 2-Fluorophenol             | 1.039 | 0.802 | 22.8  | 64    | 0.00     |
| 5 S     | Phenol-d6                  | 1.324 | 1.052 | 20.5  | 68    | 0.00     |
| 6       | bis(2-Chloroethyl)ether    | 1.243 | 1.154 | 7.2   | 76    | 0.00     |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0   | 85    | 0.00     |
| 8 S     | Nitrobenzene-d5            | 0.311 | 0.288 | 7.4   | 79    | 0.00     |
| 9       | Naphthalene                | 1.232 | 1.119 | 9.2   | 75    | 0.00     |
| 10      | Hexachlorobutadiene        | 0.253 | 0.235 | 7.1   | 76    | 0.00     |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.868 | 0.827 | 4.7   | 82    | 0.00     |
| 12      | 2-Methylnaphthalene        | 0.251 | 0.178 | 29.1# | 64    | 0.00     |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0   | 75    | 0.01     |
| 14 S    | 2,4,6-Tribromophenol       | 0.193 | 0.137 | 29.0# | 54    | 0.00     |
| 15 S    | 2-Fluorobiphenyl           | 1.830 | 1.747 | 4.5   | 68    | 0.00     |
| 16      | Acenaphthylene             | 2.118 | 1.685 | 20.4  | 60    | 0.00     |
| 17      | Acenaphthene               | 1.393 | 1.259 | 9.6   | 65    | 0.00     |
| 18      | Fluorene                   | 1.785 | 1.510 | 15.4  | 60    | 0.00     |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 67    | 0.00     |
| 20      | 4,6-Dinitro-2-methylphenol | 0.045 | 0.043 | 4.4   | 69    | 0.00     |
| 21      | 4-Bromophenyl-phenylether  | 0.289 | 0.245 | 15.2  | 55    | 0.00     |
| 22      | Hexachlorobenzene          | 0.352 | 0.339 | 3.7   | 61    | 0.00     |
| 23      | Atrazine                   | 0.204 | 0.150 | 26.5# | 50    | 0.00     |
| 24      | Pentachlorophenol          | 0.096 | 0.071 | 26.0# | 55    | 0.00     |
| 25      | Phenanthrene               | 1.391 | 1.273 | 8.5   | 59    | 0.00     |
| 26      | Anthracene                 | 1.207 | 0.974 | 19.3  | 54    | 0.00     |
| 27 SURR | Fluoranthene-d10           | 1.087 | 0.975 | 10.3  | 59    | 0.00     |
| 28      | Fluoranthene               | 1.479 | 1.195 | 19.2  | 53    | 0.00     |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 55    | 0.00     |
| 30      | Pyrene                     | 2.082 | 2.042 | 1.9   | 52    | 0.00     |
| 31 S    | Terphenyl-d14              | 0.940 | 0.833 | 11.4  | 45#   | 0.00     |
| 32      | Benzo(a)anthracene         | 1.616 | 1.378 | 14.7  | 46#   | 0.00     |
| 33      | Chrysene                   | 1.676 | 1.623 | 3.2   | 51    | 0.00     |
| 34      | Bis(2-ethylhexyl)phthalate | 0.649 | 0.610 | 6.0   | 58    | 0.00     |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0   | 54    | 0.00     |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.660 | 1.936 | -16.6 | 63    | 0.00     |
| 37      | Benzo(b)fluoranthene       | 1.786 | 1.869 | -4.6  | 56    | 0.00     |
| 38      | Benzo(k)fluoranthene       | 1.934 | 1.868 | 3.4   | 50#   | 0.00     |
| 39 C    | Benzo(a)pyrene             | 1.363 | 1.318 | 3.3   | 52    | 0.00     |
| 40      | Dibenzo(a,h)anthracene     | 1.318 | 1.540 | -16.8 | 63    | 0.00     |
| 41      | Benzo(g,h,i)perylene       | 1.420 | 1.530 | -7.7  | 56    | 0.00     |

(#) = Out of Range

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

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BNA\_N  
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SSTDCCC0.4