

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN092822\  
 Data File : BN021939.D  
 Acq On : 27 Sep 2022 17:36  
 Operator : CG/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Sep 28 04:29:49 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN091422.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Sep 28 04:21:32 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   | 106   | 0.00     |
| 2       | 1,4-Dioxane                | 0.517 | 0.470 | 9.1   | 102   | 0.00     |
| 3       | n-Nitrosodimethylamine     | 0.525 | 0.561 | -6.9  | 120   | 0.00     |
| 4 S     | 2-Fluorophenol             | 1.046 | 1.157 | -10.6 | 122   | 0.00     |
| 5 S     | Phenol-d6                  | 1.327 | 1.454 | -9.6  | 123   | 0.00     |
| 6       | bis(2-Chloroethyl)ether    | 1.316 | 1.275 | 3.1   | 107   | 0.00     |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0   | 106   | 0.00     |
| 8 S     | Nitrobenzene-d5            | 0.331 | 0.340 | -2.7  | 118   | 0.01     |
| 9       | Naphthalene                | 1.181 | 1.175 | 0.5   | 111   | 0.00     |
| 10      | Hexachlorobutadiene        | 0.197 | 0.199 | -1.0  | 111   | 0.00     |
| 11 SURR | 2-Methylnaphthalene-d10    | 1.009 | 0.846 | 16.2  | 91    | 0.00     |
| 12      | 2-Methylnaphthalene        | 0.254 | 0.208 | 18.1  | 99    | 0.00     |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0   | 90    | 0.00     |
| 14 S    | 2,4,6-Tribromophenol       | 0.183 | 0.127 | 30.6# | 69    | 0.00     |
| 15 S    | 2-Fluorobiphenyl           | 1.633 | 1.725 | -5.6  | 99    | 0.00     |
| 16      | Acenaphthylene             | 1.910 | 1.741 | 8.8   | 91    | 0.00     |
| 17      | Acenaphthene               | 1.305 | 1.278 | 2.1   | 93    | 0.00     |
| 18      | Fluorene                   | 1.735 | 1.569 | 9.6   | 85    | 0.00     |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 69    | 0.01     |
| 20      | 4,6-Dinitro-2-methylphenol | 0.000 | 0.000 | 0.0   | 0#    | -15.71#  |
| 21      | 4-Bromophenyl-phenylether  | 0.246 | 0.253 | -2.8  | 77    | 0.00     |
| 22      | Hexachlorobenzene          | 0.277 | 0.324 | -17.0 | 87    | 0.00     |
| 23      | Atrazine                   | 0.204 | 0.183 | 10.3  | 69    | 0.00     |
| 24      | Pentachlorophenol          | 0.104 | 0.071 | 31.7# | 57    | 0.00     |
| 25      | Phenanthrene               | 1.341 | 1.313 | 2.1   | 72    | 0.00     |
| 26      | Anthracene                 | 1.136 | 1.023 | 9.9   | 69    | 0.00     |
| 27 SURR | Fluoranthene-d10           | 1.076 | 0.870 | 19.1  | 60    | 0.00     |
| 28      | Fluoranthene               | 1.447 | 1.237 | 14.5  | 63    | 0.00     |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 54    | 0.00     |
| 30      | Pyrene                     | 1.986 | 2.155 | -8.5  | 60    | 0.00     |
| 31 S    | Terphenyl-d14              | 0.914 | 1.004 | -9.8  | 61    | 0.00     |
| 32      | Benzo(a)anthracene         | 1.534 | 1.434 | 6.5   | 54    | 0.00     |
| 33      | Chrysene                   | 1.595 | 1.623 | -1.8  | 58    | 0.00     |
| 34      | Bis(2-ethylhexyl)phthalate | 1.025 | 0.936 | 8.7   | 55    | 0.00     |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0   | 54    | 0.00     |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.900 | 1.817 | 4.4   | 55    | 0.00     |
| 37      | Benzo(b)fluoranthene       | 1.817 | 1.766 | 2.8   | 56    | 0.00     |
| 38      | Benzo(k)fluoranthene       | 1.776 | 1.714 | 3.5   | 54    | 0.00     |
| 39 C    | Benzo(a)pyrene             | 1.448 | 1.334 | 7.9   | 54    | 0.00     |
| 40      | Dibenzo(a,h)anthracene     | 1.504 | 1.380 | 8.2   | 52    | 0.00     |
| 41      | Benzo(g,h,i)perylene       | 1.548 | 1.623 | -4.8  | 60    | 0.00     |

(#) = Out of Range

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

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