

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN032122\  
 Data File : BN019052.D  
 Acq On : 22 Mar 2022 10:04  
 Operator : CG/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Mar 22 12:18:59 2022  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\8270-SIM-BN031722.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Mar 17 13:18:23 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   | 151#  | 0.00     |
| 2       | 1,4-Dioxane                | 0.494 | 0.450 | 8.9   | 131   | 0.00     |
| 3       | n-Nitrosodimethylamine     | 0.577 | 0.465 | 19.4  | 128   | 0.00     |
| 4 S     | 2-Fluorophenol             | 0.896 | 0.835 | 6.8   | 142   | 0.00     |
| 5 S     | Phenol-d6                  | 0.940 | 0.867 | 7.8   | 150   | 0.00     |
| 6       | bis(2-Chloroethyl)ether    | 1.118 | 1.065 | 4.7   | 146   | 0.00     |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0   | 153#  | 0.00     |
| 8 S     | Nitrobenzene-d5            | 0.289 | 0.239 | 17.3  | 142   | -0.01    |
| 9       | Naphthalene                | 1.138 | 1.096 | 3.7   | 151#  | -0.01    |
| 10      | Hexachlorobutadiene        | 0.278 | 0.257 | 7.6   | 142   | -0.01    |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.664 | 0.586 | 11.7  | 138   | 0.00     |
| 12      | 2-Methylnaphthalene        | 0.774 | 0.694 | 10.3  | 141   | 0.00     |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0   | 123   | -0.01    |
| 14 S    | 2,4,6-Tribromophenol       | 0.193 | 0.112 | 42.0# | 80    | -0.01    |
| 15 S    | 2-Fluorobiphenyl           | 1.603 | 1.657 | -3.4  | 133   | -0.01    |
| 16      | Acenaphthylene             | 1.856 | 1.703 | 8.2   | 124   | 0.00     |
| 17      | Acenaphthene               | 1.302 | 1.198 | 8.0   | 118   | -0.01    |
| 18      | Fluorene                   | 1.641 | 1.390 | 15.3  | 106   | -0.01    |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 90    | -0.01    |
| 20      | 4,6-Dinitro-2-methylphenol | 0.049 | 0.014 | 71.4# | 29#   | 0.01     |
| 21      | 4-Bromophenyl-phenylether  | 0.252 | 0.258 | -2.4  | 96    | 0.00     |
| 22      | Hexachlorobenzene          | 0.325 | 0.341 | -4.9  | 95    | 0.00     |
| 23      | Atrazine                   | 0.188 | 0.150 | 20.2  | 80    | -0.01    |
| 24      | Pentachlorophenol          | 0.100 | 0.045 | 55.0# | 47#   | 0.00     |
| 25      | Phenanthrene               | 1.251 | 1.165 | 6.9   | 88    | -0.01    |
| 26      | Anthracene                 | 1.057 | 0.891 | 15.7  | 83    | -0.01    |
| 27 SURR | Fluoranthene-d10           | 1.187 | 0.921 | 22.4  | 71    | 0.00     |
| 28      | Fluoranthene               | 1.468 | 1.133 | 22.8  | 71    | 0.00     |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 71    | 0.00     |
| 30      | Pyrene                     | 1.993 | 1.973 | 1.0   | 71    | 0.00     |
| 31 S    | Terphenyl-d14              | 0.982 | 0.847 | 13.7  | 63    | 0.00     |
| 32      | Benzo(a)anthracene         | 1.233 | 1.103 | 10.5  | 71    | 0.00     |
| 33      | Chrysene                   | 1.662 | 1.578 | 5.1   | 67    | 0.00     |
| 34      | Bis(2-ethylhexyl)phthalate | 0.817 | 0.740 | 9.4   | 70    | 0.00     |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0   | 87    | 0.00     |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.443 | 1.348 | 6.6   | 85    | -0.02    |
| 37      | Benzo(b)fluoranthene       | 1.459 | 1.214 | 16.8  | 80    | 0.00     |
| 38      | Benzo(k)fluoranthene       | 2.000 | 1.822 | 8.9   | 79    | 0.00     |
| 39 C    | Benzo(a)pyrene             | 1.328 | 1.219 | 8.2   | 82    | 0.00     |
| 40      | Dibenzo(a,h)anthracene     | 1.124 | 1.056 | 6.0   | 86    | -0.02    |
| 41      | Benzo(g,h,i)perylene       | 1.417 | 1.556 | -9.8  | 106   | -0.02    |

(#) = Out of Range

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

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