

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN011422\  
 Data File : BN018402.D  
 Acq On : 14 Jan 2022 11:22  
 Operator : CG/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Jan 14 12:13:14 2022  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\8270-SIM-BN010722.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Jan 14 12:12:13 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   | 75    | -0.02    |
| 2       | 1,4-Dioxane                | 0.225 | 0.181 | 19.6  | 62    | -0.03    |
| 3       | n-Nitrosodimethylamine     | 0.361 | 0.390 | -8.0  | 87    | 0.00     |
| 4 S     | 2-Fluorophenol             | 0.984 | 0.968 | 1.6   | 78    | -0.02    |
| 5 S     | Phenol-d6                  | 0.986 | 0.967 | 1.9   | 78    | -0.02    |
| 6       | bis(2-Chloroethyl)ether    | 1.068 | 1.077 | -0.8  | 76    | -0.02    |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0   | 76    | -0.03    |
| 8 S     | Nitrobenzene-d5            | 0.255 | 0.230 | 9.8   | 84    | -0.01    |
| 9       | Naphthalene                | 1.016 | 1.017 | -0.1  | 76    | -0.01    |
| 10      | Hexachlorobutadiene        | 0.189 | 0.197 | -4.2  | 80    | -0.01    |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.626 | 0.628 | -0.3  | 77    | -0.02    |
| 12      | 2-Methylnaphthalene        | 0.622 | 0.642 | -3.2  | 78    | -0.01    |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0   | 80    | -0.01    |
| 14 S    | 2,4,6-Tribromophenol       | 0.158 | 0.140 | 11.4  | 101   | -0.01    |
| 15 S    | 2-Fluorobiphenyl           | 1.444 | 1.463 | -1.3  | 81    | -0.03    |
| 16      | Acenaphthylene             | 1.588 | 1.464 | 7.8   | 79    | -0.01    |
| 17      | Acenaphthene               | 1.109 | 1.088 | 1.9   | 79    | -0.01    |
| 18      | Fluorene                   | 1.246 | 1.272 | -2.1  | 83    | -0.01    |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 86    | -0.01    |
| 20      | 4,6-Dinitro-2-methylphenol | 0.028 | 0.012 | 57.1# | 98    | -0.02    |
| 21      | 4-Bromophenyl-phenylether  | 0.220 | 0.213 | 3.2   | 87    | -0.01    |
| 22      | Hexachlorobenzene          | 0.293 | 0.292 | 0.3   | 86    | -0.01    |
| 23      | Atrazine                   | 0.146 | 0.126 | 13.7  | 94    | -0.01    |
| 24      | Pentachlorophenol          | 0.071 | 0.046 | 35.2# | 114   | -0.01    |
| 25      | Phenanthrene               | 1.124 | 1.145 | -1.9  | 88    | 0.00     |
| 26      | Anthracene                 | 0.929 | 0.892 | 4.0   | 88    | -0.01    |
| 27 SURR | Fluoranthene-d10           | 1.142 | 1.130 | 1.1   | 87    | 0.00     |
| 28      | Fluoranthene               | 1.194 | 1.275 | -6.8  | 90    | 0.00     |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 90    | 0.00     |
| 30      | Pyrene                     | 1.416 | 1.459 | -3.0  | 92    | 0.00     |
| 31 S    | Terphenyl-d14              | 0.866 | 0.910 | -5.1  | 94    | 0.00     |
| 32      | Benzo(a)anthracene         | 1.178 | 1.107 | 6.0   | 94    | 0.00     |
| 33      | Chrysene                   | 1.338 | 1.377 | -2.9  | 90    | -0.01    |
| 34      | Bis(2-ethylhexyl)phthalate | 0.707 | 0.516 | 27.0# | 96    | 0.00     |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0   | 93    | 0.00     |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.271 | 1.108 | 12.8  | 90    | 0.00     |
| 37      | Benzo(b)fluoranthene       | 1.297 | 1.302 | -0.4  | 91    | 0.00     |
| 38      | Benzo(k)fluoranthene       | 1.268 | 1.315 | -3.7  | 96    | 0.00     |
| 39 C    | Benzo(a)pyrene             | 1.199 | 1.144 | 4.6   | 92    | 0.00     |
| 40      | Dibenzo(a,h)anthracene     | 0.951 | 0.780 | 18.0  | 89    | 0.00     |
| 41      | Benzo(g,h,i)perylene       | 1.163 | 1.038 | 10.7  | 87    | -0.01    |

(#) = Out of Range

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

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