

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN011322\  
 Data File : BN018395.D  
 Acq On : 13 Jan 2022 13:04  
 Operator : CG/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Jan 13 20:20:17 2022  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\8270-SIM-BN010722.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jan 10 13:58:28 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   | 83    | -0.02    |
| 2       | 1,4-Dioxane                | 0.225 | 0.181 | 19.6  | 68    | -0.03    |
| 3       | n-Nitrosodimethylamine     | 0.361 | 0.400 | -10.8 | 98    | 0.00     |
| 4 S     | 2-Fluorophenol             | 0.984 | 1.124 | -14.2 | 99    | 0.00     |
| 5 S     | Phenol-d6                  | 0.986 | 1.082 | -9.7  | 97    | -0.02    |
| 6       | bis(2-Chloroethyl)ether    | 1.068 | 1.182 | -10.7 | 92    | -0.02    |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0   | 84    | -0.03    |
| 8 S     | Nitrobenzene-d5            | 0.255 | 0.262 | -2.7  | 105   | -0.01    |
| 9       | Naphthalene                | 1.016 | 1.078 | -6.1  | 89    | -0.01    |
| 10      | Hexachlorobutadiene        | 0.189 | 0.206 | -9.0  | 92    | -0.01    |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.626 | 0.626 | 0.0   | 84    | -0.02    |
| 12      | 2-Methylnaphthalene        | 0.622 | 0.662 | -6.4  | 88    | -0.01    |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0   | 86    | -0.01    |
| 14 S    | 2,4,6-Tribromophenol       | 0.158 | 0.140 | 11.4  | 108   | -0.01    |
| 15 S    | 2-Fluorobiphenyl           | 1.444 | 1.556 | -7.8  | 92    | -0.01    |
| 16      | Acenaphthylene             | 1.588 | 1.526 | 3.9   | 88    | 0.00     |
| 17      | Acenaphthene               | 1.109 | 1.125 | -1.4  | 88    | -0.01    |
| 18      | Fluorene                   | 1.246 | 1.313 | -5.4  | 92    | -0.01    |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 91    | -0.01    |
| 20      | 4,6-Dinitro-2-methylphenol | 0.028 | 0.015 | 46.4# | 128   | -0.03    |
| 21      | 4-Bromophenyl-phenylether  | 0.220 | 0.220 | 0.0   | 95    | -0.01    |
| 22      | Hexachlorobenzene          | 0.293 | 0.298 | -1.7  | 93    | -0.01    |
| 23      | Atrazine                   | 0.146 | 0.126 | 13.7  | 99    | 0.00     |
| 24      | Pentachlorophenol          | 0.071 | 0.049 | 31.0# | 130   | -0.01    |
| 25      | Phenanthrene               | 1.124 | 1.181 | -5.1  | 96    | 0.00     |
| 26      | Anthracene                 | 0.929 | 0.913 | 1.7   | 96    | -0.01    |
| 27 SURR | Fluoranthene-d10           | 1.142 | 1.121 | 1.8   | 92    | -0.01    |
| 28      | Fluoranthene               | 1.194 | 1.306 | -9.4  | 98    | -0.01    |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 97    | 0.00     |
| 30      | Pyrene                     | 1.416 | 1.466 | -3.5  | 100   | 0.00     |
| 31 S    | Terphenyl-d14              | 0.866 | 0.943 | -8.9  | 106   | 0.00     |
| 32      | Benzo(a)anthracene         | 1.178 | 1.159 | 1.6   | 107   | -0.01    |
| 33      | Chrysene                   | 1.338 | 1.425 | -6.5  | 101   | -0.01    |
| 34      | Bis(2-ethylhexyl)phthalate | 0.707 | 0.541 | 23.5  | 109   | 0.00     |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0   | 107   | 0.00     |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.271 | 1.256 | 1.2   | 117   | -0.01    |
| 37      | Benzo(b)fluoranthene       | 1.297 | 1.340 | -3.3  | 107   | 0.00     |
| 38      | Benzo(k)fluoranthene       | 1.268 | 1.385 | -9.2  | 116   | 0.00     |
| 39 C    | Benzo(a)pyrene             | 1.199 | 1.225 | -2.2  | 112   | 0.00     |
| 40      | Dibenzo(a,h)anthracene     | 0.951 | 0.901 | 5.3   | 117   | -0.01    |
| 41      | Benzo(g,h,i)perylene       | 1.163 | 1.153 | 0.9   | 110   | -0.01    |

(#) = Out of Range

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

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