

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN081622\  
 Data File : BN021249.D  
 Acq On : 16 Aug 2022 09:05  
 Operator : CG/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Aug 16 11:55:20 2022  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\8270-SIM-BN081622.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Aug 16 09:19:59 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   | 125   | 0.00     |
| 2       | 1,4-Dioxane                | 0.558 | 0.464 | 16.8  | 115   | 0.00     |
| 3       | n-Nitrosodimethylamine     | 0.455 | 0.350 | 23.1  | 127   | 0.00     |
| 4 S     | 2-Fluorophenol             | 0.998 | 0.978 | 2.0   | 136   | 0.00     |
| 5 S     | Phenol-d6                  | 1.165 | 1.160 | 0.4   | 135   | 0.00     |
| 6       | bis(2-Chloroethyl)ether    | 0.983 | 0.897 | 8.7   | 115   | 0.00     |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0   | 119   | -0.01    |
| 8 S     | Nitrobenzene-d5            | 0.296 | 0.241 | 18.6  | 136   | -0.01    |
| 9       | Naphthalene                | 1.210 | 1.122 | 7.3   | 117   | 0.00     |
| 10      | Hexachlorobutadiene        | 0.217 | 0.195 | 10.1  | 117   | -0.01    |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.696 | 0.618 | 11.2  | 119   | 0.00     |
| 12      | 2-Methylnaphthalene        | 0.825 | 0.726 | 12.0  | 118   | 0.00     |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0   | 112   | 0.00     |
| 14 S    | 2,4,6-Tribromophenol       | 0.131 | 0.093 | 29.0# | 101   | 0.00     |
| 15 S    | 2-Fluorobiphenyl           | 1.591 | 1.574 | 1.1   | 117   | 0.00     |
| 16      | Acenaphthylene             | 1.922 | 1.683 | 12.4  | 114   | 0.00     |
| 17      | Acenaphthene               | 1.321 | 1.166 | 11.7  | 107   | 0.00     |
| 18      | Fluorene                   | 1.680 | 1.564 | 6.9   | 116   | 0.00     |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 120   | 0.00     |
| 20      | 4,6-Dinitro-2-methylphenol | 0.040 | 0.020 | 50.0# | 80    | 0.00     |
| 21      | 4-Bromophenyl-phenylether  | 0.246 | 0.203 | 17.5  | 118   | 0.00     |
| 22      | Hexachlorobenzene          | 0.277 | 0.256 | 7.6   | 130   | 0.00     |
| 23      | Atrazine                   | 0.158 | 0.132 | 16.5  | 128   | 0.00     |
| 24      | Pentachlorophenol          | 0.062 | 0.061 | 1.6   | 152#  | -0.01    |
| 25      | Phenanthrene               | 1.331 | 1.199 | 9.9   | 129   | 0.00     |
| 26      | Anthracene                 | 1.112 | 0.961 | 13.6  | 128   | -0.01    |
| 27 SURR | Fluoranthene-d10           | 1.086 | 1.098 | -1.1  | 141   | 0.00     |
| 28      | Fluoranthene               | 1.392 | 1.379 | 0.9   | 142   | 0.00     |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 138   | 0.00     |
| 30      | Pyrene                     | 2.060 | 1.980 | 3.9   | 142   | 0.00     |
| 31 S    | Terphenyl-d14              | 1.091 | 1.046 | 4.1   | 142   | 0.00     |
| 32      | Benzo(a)anthracene         | 1.301 | 1.085 | 16.6  | 131   | 0.00     |
| 33      | Chrysene                   | 1.617 | 1.542 | 4.6   | 145   | 0.00     |
| 34      | Bis(2-ethylhexyl)phthalate | 0.889 | 0.712 | 19.9  | 123   | 0.00     |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0   | 110   | 0.00     |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.769 | 1.479 | 16.4  | 104   | 0.00     |
| 37      | Benzo(b)fluoranthene       | 1.669 | 1.528 | 8.4   | 108   | 0.00     |
| 38      | Benzo(k)fluoranthene       | 1.684 | 1.522 | 9.6   | 110   | 0.00     |
| 39 C    | Benzo(a)pyrene             | 1.392 | 1.204 | 13.5  | 101   | 0.00     |
| 40      | Dibenzo(a,h)anthracene     | 1.368 | 1.142 | 16.5  | 103   | 0.00     |
| 41      | Benzo(g,h,i)perylene       | 1.590 | 1.499 | 5.7   | 115   | 0.00     |

(#) = Out of Range

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

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