

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN032123\  
 Data File : BN024470.D  
 Acq On : 22 Mar 2023 05:51  
 Operator : CG/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Mar 22 06:20:41 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN031323.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Mar 22 04:14:21 2023  
 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   | 132   | 0.00     |
| 2       | 1,4-Dioxane                | 0.435 | 0.424 | 2.5   | 123   | 0.00     |
| 3       | n-Nitrosodimethylamine     | 0.638 | 0.643 | -0.8  | 131   | 0.00     |
| 4 S     | 2-Fluorophenol             | 0.886 | 0.909 | -2.6  | 136   | 0.00     |
| 5 S     | Phenol-d6                  | 1.138 | 1.130 | 0.7   | 135   | 0.00     |
| 6       | bis(2-Chloroethyl)ether    | 1.129 | 1.140 | -1.0  | 130   | 0.00     |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0   | 133   | 0.00     |
| 8 S     | Nitrobenzene-d5            | 0.267 | 0.313 | -17.2 | 164#  | 0.00     |
| 9       | Naphthalene                | 1.017 | 1.024 | -0.7  | 135   | -0.01    |
| 10      | Hexachlorobutadiene        | 0.193 | 0.201 | -4.1  | 137   | -0.01    |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.492 | 0.480 | 2.4   | 133   | 0.00     |
| 12      | 2-Methylnaphthalene        | 0.687 | 0.663 | 3.5   | 131   | 0.00     |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0   | 125   | 0.00     |
| 14 S    | 2,4,6-Tribromophenol       | 0.204 | 0.160 | 21.6  | 107   | 0.00     |
| 15 S    | 2-Fluorobiphenyl           | 1.511 | 1.526 | -1.0  | 128   | -0.01    |
| 16      | Acenaphthylene             | 1.897 | 1.714 | 9.6   | 118   | 0.00     |
| 17      | Acenaphthene               | 1.208 | 1.138 | 5.8   | 120   | -0.01    |
| 18      | Fluorene                   | 1.443 | 1.293 | 10.4  | 115   | -0.01    |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 118   | -0.01    |
| 20      | 4,6-Dinitro-2-methylphenol | 0.053 | 0.025 | 52.8# | 70    | -0.01    |
| 21      | 4-Bromophenyl-phenylether  | 0.238 | 0.226 | 5.0   | 116   | 0.00     |
| 22      | Hexachlorobenzene          | 0.292 | 0.285 | 2.4   | 118   | 0.00     |
| 23      | Atrazine                   | 0.151 | 0.135 | 10.6  | 119   | -0.01    |
| 24      | Pentachlorophenol          | 0.136 | 0.083 | 39.0# | 84    | -0.01    |
| 25      | Phenanthrene               | 1.185 | 1.130 | 4.6   | 115   | 0.00     |
| 26      | Anthracene                 | 1.094 | 0.987 | 9.8   | 113   | 0.00     |
| 27 SURR | Fluoranthene-d10           | 0.848 | 0.805 | 5.1   | 116   | 0.00     |
| 28      | Fluoranthene               | 1.297 | 1.203 | 7.2   | 114   | 0.00     |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 105   | 0.00     |
| 30      | Pyrene                     | 1.621 | 1.793 | -10.6 | 115   | 0.00     |
| 31 S    | Terphenyl-d14              | 0.674 | 0.644 | 4.5   | 100   | 0.00     |
| 32      | Benzo(a)anthracene         | 1.336 | 1.283 | 4.0   | 105   | 0.00     |
| 33      | Chrysene                   | 1.425 | 1.407 | 1.3   | 103   | 0.00     |
| 34      | Bis(2-ethylhexyl)phthalate | 0.629 | 0.713 | -13.4 | 130   | 0.00     |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0   | 102   | 0.00     |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.498 | 1.407 | 6.1   | 101   | 0.00     |
| 37      | Benzo(b)fluoranthene       | 1.556 | 1.549 | 0.4   | 108   | 0.00     |
| 38      | Benzo(k)fluoranthene       | 1.605 | 1.486 | 7.4   | 100   | 0.00     |
| 39 C    | Benzo(a)pyrene             | 1.427 | 1.336 | 6.4   | 102   | 0.00     |
| 40      | Dibenzo(a,h)anthracene     | 1.161 | 1.055 | 9.1   | 99    | -0.01    |
| 41      | Benzo(g,h,i)perylene       | 1.354 | 1.336 | 1.3   | 106   | -0.01    |

(#) = Out of Range

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

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BNA\_N  
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SSTDCCC0.4