

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN050224\  
 Data File : BN031382.D  
 Acq On : 02 May 2024 21:41  
 Operator : MA/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: May 03 01:12:19 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN041524.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu May 02 15:02:22 2024  
 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   | 61    | 0.00     |
| 2       | 1,4-Dioxane                | 0.492 | 0.485 | 1.4   | 63    | 0.00     |
| 3       | n-Nitrosodimethylamine     | 0.461 | 0.497 | -7.8  | 71    | 0.00     |
| 4 S     | 2-Fluorophenol             | 0.891 | 0.998 | -12.0 | 72    | 0.00     |
| 5 S     | Phenol-d6                  | 1.079 | 1.073 | 0.6   | 67    | 0.00     |
| 6       | bis(2-Chloroethyl)ether    | 1.067 | 0.933 | 12.6  | 57    | 0.00     |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0   | 56    | 0.00     |
| 8 S     | Nitrobenzene-d5            | 0.277 | 0.298 | -7.6  | 65    | 0.00     |
| 9       | Naphthalene                | 1.103 | 1.008 | 8.6   | 53    | -0.01    |
| 10      | Hexachlorobutadiene        | 0.221 | 0.250 | -13.1 | 65    | 0.00     |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.616 | 0.571 | 7.3   | 55    | 0.00     |
| 12      | 2-Methylnaphthalene        | 0.733 | 0.639 | 12.8  | 51    | 0.00     |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0   | 51    | 0.00     |
| 14 S    | 2,4,6-Tribromophenol       | 0.155 | 0.162 | -4.5  | 60    | 0.00     |
| 15 S    | 2-Fluorobiphenyl           | 1.345 | 1.487 | -10.6 | 58    | 0.00     |
| 16      | Acenaphthylene             | 1.763 | 1.580 | 10.4  | 50    | 0.00     |
| 17      | Acenaphthene               | 1.244 | 1.109 | 10.9  | 48#   | 0.00     |
| 18      | Fluorene                   | 1.647 | 1.409 | 14.5  | 46#   | 0.00     |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 47#   | 0.00     |
| 20      | 4,6-Dinitro-2-methylphenol | 0.057 | 0.057 | 0.0   | 60    | 0.00     |
| 21      | 4-Bromophenyl-phenylether  | 0.238 | 0.218 | 8.4   | 46#   | 0.00     |
| 22      | Hexachlorobenzene          | 0.278 | 0.257 | 7.6   | 45#   | 0.00     |
| 23      | Atrazine                   | 0.172 | 0.177 | -2.9  | 55    | 0.00     |
| 24      | Pentachlorophenol          | 0.097 | 0.110 | -13.4 | 70    | 0.00     |
| 25      | Phenanthrene               | 1.181 | 1.020 | 13.6  | 43#   | 0.00     |
| 26      | Anthracene                 | 0.977 | 0.883 | 9.6   | 47#   | 0.00     |
| 27 SURR | Fluoranthene-d10           | 1.077 | 1.051 | 2.4   | 49#   | 0.00     |
| 28      | Fluoranthene               | 1.416 | 1.246 | 12.0  | 45#   | 0.00     |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 52    | 0.00     |
| 30      | Pyrene                     | 1.553 | 1.351 | 13.0  | 47#   | 0.00     |
| 31 S    | Terphenyl-d14              | 0.929 | 0.914 | 1.6   | 54    | 0.00     |
| 32      | Benzo(a)anthracene         | 1.393 | 1.248 | 10.4  | 51    | 0.00     |
| 33      | Chrysene                   | 1.515 | 1.314 | 13.3  | 47#   | 0.00     |
| 34      | Bis(2-ethylhexyl)phthalate | 0.638 | 0.770 | -20.7 | 71    | 0.00     |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0   | 63    | 0.00     |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.860 | 1.592 | 14.4  | 59    | -0.01    |
| 37      | Benzo(b)fluoranthene       | 1.650 | 1.347 | 18.4  | 56    | 0.00     |
| 38      | Benzo(k)fluoranthene       | 1.616 | 1.276 | 21.0  | 54    | 0.00     |
| 39 C    | Benzo(a)pyrene             | 1.374 | 1.129 | 17.8  | 58    | 0.00     |
| 40      | Dibenzo(a,h)anthracene     | 1.490 | 1.285 | 13.8  | 60    | 0.00     |
| 41      | Benzo(g,h,i)perylene       | 1.640 | 1.437 | 12.4  | 58    | 0.00     |

(#) = Out of Range

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

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