

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061523\  
 Data File : BN025867.D  
 Acq On : 15 Jun 2023 21:09  
 Operator : MA/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC0.4

Quant Time: Jun 16 01:47:38 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\8270-SIM-BN053123.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Jun 16 01:41:46 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   | 79    | 0.00     |
| 2       | 1,4-Dioxane                | 0.469 | 0.485 | -3.4  | 74    | 0.00     |
| 3       | n-Nitrosodimethylamine     | 0.599 | 0.404 | 32.6# | 50#   | 0.01     |
| 4 S     | 2-Fluorophenol             | 1.125 | 1.005 | 10.7  | 64    | 0.01     |
| 5 S     | Phenol-d6                  | 1.381 | 1.152 | 16.6  | 61    | 0.00     |
| 6       | bis(2-Chloroethyl)ether    | 1.254 | 1.097 | 12.5  | 62    | 0.00     |
| 7 I     | Naphthalene-d8             | 1.000 | 1.000 | 0.0   | 69    | 0.00     |
| 8 S     | Nitrobenzene-d5            | 0.372 | 0.342 | 8.1   | 62    | 0.01     |
| 9       | Naphthalene                | 1.196 | 1.128 | 5.7   | 62    | 0.01     |
| 10      | Hexachlorobutadiene        | 0.184 | 0.217 | -17.9 | 76    | 0.00     |
| 11 SURR | 2-Methylnaphthalene-d10    | 0.576 | 0.553 | 4.0   | 62    | 0.00     |
| 12      | 2-Methylnaphthalene        | 0.761 | 0.736 | 3.3   | 62    | 0.00     |
| 13 I    | Acenaphthene-d10           | 1.000 | 1.000 | 0.0   | 62    | 0.00     |
| 14 S    | 2,4,6-Tribromophenol       | 0.123 | 0.125 | -1.6  | 63    | 0.00     |
| 15 S    | 2-Fluorobiphenyl           | 1.691 | 1.720 | -1.7  | 61    | 0.00     |
| 16      | Acenaphthylene             | 1.926 | 1.918 | 0.4   | 60    | 0.00     |
| 17      | Acenaphthene               | 1.263 | 1.293 | -2.4  | 61    | 0.00     |
| 18      | Fluorene                   | 1.626 | 1.658 | -2.0  | 58    | 0.01     |
| 19 I    | Phenanthrene-d10           | 1.000 | 1.000 | 0.0   | 64    | 0.00     |
| 20      | 4,6-Dinitro-2-methylphenol | 0.072 | 0.055 | 23.6  | 48#   | 0.01     |
| 21      | 4-Bromophenyl-phenylether  | 0.232 | 0.232 | 0.0   | 61    | 0.00     |
| 22      | Hexachlorobenzene          | 0.198 | 0.230 | -16.2 | 70    | 0.00     |
| 23      | Atrazine                   | 0.189 | 0.187 | 1.1   | 58    | 0.01     |
| 24      | Pentachlorophenol          | 0.074 | 0.045 | 39.2# | 39#   | 0.01     |
| 25      | Phenanthrene               | 1.229 | 1.256 | -2.2  | 61    | 0.00     |
| 26      | Anthracene                 | 1.125 | 1.097 | 2.5   | 59    | 0.01     |
| 27 SURR | Fluoranthene-d10           | 0.823 | 0.941 | -14.3 | 65    | 0.00     |
| 28      | Fluoranthene               | 1.213 | 1.406 | -15.9 | 66    | 0.00     |
| 29 I    | Chrysene-d12               | 1.000 | 1.000 | 0.0   | 84    | 0.00     |
| 30      | Pyrene                     | 2.267 | 1.947 | 14.1  | 68    | 0.00     |
| 31 S    | Terphenyl-d14              | 0.970 | 0.819 | 15.6  | 68    | 0.00     |
| 32      | Benzo(a)anthracene         | 1.544 | 1.591 | -3.0  | 83    | 0.00     |
| 33      | Chrysene                   | 1.635 | 1.781 | -8.9  | 84    | 0.00     |
| 34      | Bis(2-ethylhexyl)phthalate | 0.920 | 0.966 | -5.0  | 75    | 0.00     |
| 35 I    | Perylene-d12               | 1.000 | 1.000 | 0.0   | 105   | 0.00     |
| 36      | Indeno(1,2,3-cd)pyrene     | 1.730 | 2.007 | -16.0 | 122   | 0.00     |
| 37      | Benzo(b)fluoranthene       | 1.578 | 1.922 | -21.8 | 117   | 0.00     |
| 38      | Benzo(k)fluoranthene       | 1.612 | 1.842 | -14.3 | 113   | 0.00     |
| 39 C    | Benzo(a)pyrene             | 1.495 | 1.678 | -12.2 | 112   | 0.00     |
| 40      | Dibenzo(a,h)anthracene     | 1.389 | 1.583 | -14.0 | 121   | 0.01     |
| 41      | Benzo(g,h,i)perylene       | 1.552 | 1.851 | -19.3 | 126   | 0.00     |

(#) = Out of Range

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

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BNA\_N  
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SSTDCCC0.4