

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091522\
 Data File : BN021782.D
 Acq On : 16 Sep 2022 00:48
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 16 06:28:26 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 14 01:55:04 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 | 106 | 0.00 |
| 2 | 1,4-Dioxane | 0.517 | 0.487 | 5.8 | 106 | 0.00 |
| 3 | n-Nitrosodimethylamine | 0.525 | 0.523 | 0.4 | 112 | 0.00 |
| 4 S | 2-Fluorophenol | 1.046 | 1.029 | 1.6 | 109 | 0.00 |
| 5 S | Phenol-d6 | 1.327 | 1.317 | 0.8 | 112 | 0.00 |
| 6 | bis(2-Chloroethyl)ether | 1.316 | 1.263 | 4.0 | 107 | 0.00 |
| 7 I | Naphthalene-d8 | 1.000 | 1.000 | 0.0 | 106 | -0.01 |
| 8 S | Nitrobenzene-d5 | 0.331 | 0.308 | 6.9 | 106 | -0.01 |
| 9 | Naphthalene | 1.181 | 1.146 | 3.0 | 108 | -0.01 |
| 10 | Hexachlorobutadiene | 0.197 | 0.195 | 1.0 | 108 | 0.00 |
| 11 SURR | 2-Methylnaphthalene-d10 | 1.009 | 0.953 | 5.6 | 103 | 0.00 |
| 12 | 2-Methylnaphthalene | 0.254 | 0.229 | 9.8 | 109 | 0.00 |
| 13 I | Acenaphthene-d10 | 1.000 | 1.000 | 0.0 | 101 | 0.00 |
| 14 S | 2,4,6-Tribromophenol | 0.183 | 0.138 | 24.6 | 85 | -0.01 |
| 15 S | 2-Fluorobiphenyl | 1.633 | 1.614 | 1.2 | 104 | 0.00 |
| 16 | Acenaphthylene | 1.910 | 1.735 | 9.2 | 102 | 0.00 |
| 17 | Acenaphthene | 1.305 | 1.240 | 5.0 | 102 | 0.00 |
| 18 | Fluorene | 1.735 | 1.653 | 4.7 | 101 | -0.01 |
| 19 I | Phenanthrene-d10 | 1.000 | 1.000 | 0.0 | 96 | 0.00 |
| 20 | 4,6-Dinitro-2-methylphenol | 0.000 | 0.000 | 0.0 | 0# | -15.71# |
| 21 | 4-Bromophenyl-phenylether | 0.246 | 0.233 | 5.3 | 98 | 0.00 |
| 22 | Hexachlorobenzene | 0.277 | 0.277 | 0.0 | 103 | 0.00 |
| 23 | Atrazine | 0.204 | 0.173 | 15.2 | 90 | 0.00 |
| 24 | Pentachlorophenol | 0.104 | 0.058 | 44.2# | 64 | 0.00 |
| 25 | Phenanthrene | 1.341 | 1.282 | 4.4 | 97 | 0.00 |
| 26 | Anthracene | 1.136 | 1.029 | 9.4 | 96 | 0.00 |
| 27 SURR | Fluoranthene-d10 | 1.076 | 0.979 | 9.0 | 93 | 0.00 |
| 28 | Fluoranthene | 1.447 | 1.331 | 8.0 | 95 | 0.00 |
| 29 I | Chrysene-d12 | 1.000 | 1.000 | 0.0 | 89 | -0.01 |
| 30 | Pyrene | 1.986 | 2.014 | -1.4 | 92 | 0.00 |
| 31 S | Terphenyl-d14 | 0.914 | 0.927 | -1.4 | 92 | 0.00 |
| 32 | Benzo(a)anthracene | 1.534 | 1.409 | 8.1 | 88 | -0.01 |
| 33 | Chrysene | 1.595 | 1.577 | 1.1 | 92 | -0.01 |
| 34 | Bis(2-ethylhexyl)phthalate | 1.025 | 0.813 | 20.7 | 78 | -0.01 |
| 35 I | Perylene-d12 | 1.000 | 1.000 | 0.0 | 81 | -0.02 |
| 36 | Indeno(1,2,3-cd)pyrene | 1.900 | 1.837 | 3.3 | 83 | -0.02 |
| 37 | Benzo(b)fluoranthene | 1.817 | 1.807 | 0.6 | 86 | -0.01 |
| 38 | Benzo(k)fluoranthene | 1.776 | 1.763 | 0.7 | 84 | -0.01 |
| 39 C | Benzo(a)pyrene | 1.448 | 1.415 | 2.3 | 85 | -0.01 |
| 40 | Dibenzo(a,h)anthracene | 1.504 | 1.450 | 3.6 | 83 | -0.02 |
| 41 | Benzo(g,h,i)perylene | 1.548 | 1.485 | 4.1 | 83 | -0.02 |

(#) = Out of Range

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

Instrument :

BNA_N

LabSampleId :

SSTDCCC0.4