Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN081219\
Data File : BN007128.D
Acq On : 13 Aug 2019 12:55
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Aug 13 13:57:42 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN081019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 13 12:12:02 2019
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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	155	0.00
2		n-Nitrosodimethylamine	0.400	0.373	6.8	184	0.00
3	S	2-Fluorophenol	0.400	0.426	-6.5	172	0.00
4	S	Phenol-d6	0.400	0.437	-9.2	182	0.00
5		bis(2-Chloroethyl)ether	0.400	0.392	2.0	159	0.00
6	I	Naphthalene-d8	0.400	0.400	0.0	167	0.00
7	S	Nitrobenzene-d5	0.400	0.410	-2.5	187	0.00
8		Naphthalene	0.400	0.376	6.0	168	0.00
9		Hexachlorobutadiene	0.400	0.379	5.3	166	-0.01
10	SURR	2-Methylnaphthalene-d10	0.400	0.370	7.5	167	0.00
11		2-Methylnaphthalene	0.400	0.372	7.0	168	0.00
12	I	Acenaphthene-d10	0.400	0.400	0.0	162	0.00
13	S	2,4,6-Tribromophenol	0.400	0.344	14.0	191	0.00
14	S	2-Fluorobiphenyl	0.400	0.230	42.5#	85	0.00
15		Acenaphthylene	0.400	0.487	-21.7	210	0.00
16		Acenaphthene	0.400	0.385	3.8	167	-0.01
17		Fluorene	0.400	0.352	12.0	172	0.00
18	I	Phenanthrene-d10	0.400	0.400	0.0	146	0.00
19		4-Bromophenyl-phenylether	0.400	0.394	1.5	146	-0.01
20		Hexachlorobenzene	0.400	0.402	-0.5	149	0.00
21		Pentachlorophenol	0.400	0.385	3.8	163	0.00
22		Phenanthrene	0.400	0.390	2.5	148	0.00
23		Anthracene	0.400	0.402	-0.5	158	-0.01
24	SURR	Fluoranthene-d10	0.400	0.357	10.8	137	0.00
25		Fluoranthene	0.400	0.364	9.0	138	0.00
26	I	Chrysene-d12	0.400	0.400	0.0	131	-0.01
27		Pyrene	0.400	0.423	-5.7	143	0.00
28	S	Terphenyl-d14	0.400	0.413	-3.2	140	-0.01
29		Benzo(a)anthracene	0.400	0.400	0.0	134	0.00
30		Chrysene	0.400	0.399	0.3	136	-0.01
31		Indeno(1,2,3-cd)pyrene	0.400	0.429	-7.2	149	-0.01
32	I	Perylene-d12	0.400	0.400	0.0	146	-0.01
33		Benzo(b)fluoranthene	0.400	0.390	2.5	145	-0.01
34		Benzo(k)fluoranthene	0.400	0.373	6.8	136	0.00
35	C	Benzo(a)pyrene	0.400	0.397	0.8	151	-0.01
36		Dibenzo(a,h)anthracene	0.400	0.394	1.5	150	-0.01
37		Benzo(g,h,i)perylene	0.400	0.389	2.8	144	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 0