Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN081319\
Data File : BN007166.D
Acq On : 14 Aug 2019 12:58
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Aug 14 13:37:28 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	185	0.00
2		n-Nitrosodimethylamine	0.400	0.368	8.0	217	0.00
3	S	2-Fluorophenol	0.400	0.436	-9.0	211	0.00
4	S	Phenol-d6	0.400	0.445	-11.2	222	0.00
5		bis(2-Chloroethyl)ether	0.400	0.389	2.8	188	0.00
6	I	Naphthalene-d8	0.400	0.400	0.0	195	0.00
7	S	Nitrobenzene-d5	0.400	0.423	-5.7	225	0.00
8		Naphthalene	0.400	0.380	5.0	198	0.00
9		Hexachlorobutadiene	0.400	0.394	1.5	202	-0.01
10	SURR	2-Methylnaphthalene-d10	0.400	0.372	7.0	196	0.00
11		2-Methylnaphthalene	0.400	0.386	3.5	203	0.00
12	I	Acenaphthene-d10	0.400	0.400	0.0	191	0.00
13	S	2,4,6-Tribromophenol	0.400	0.373	6.8	245	0.00
14	S	2-Fluorobiphenyl	0.400	0.222	44.5#	96	0.00
15		Acenaphthylene	0.400	0.704	-76.0#	357	0.00
16		Acenaphthene	0.400	0.412	-3.0	211	-0.01
17		Fluorene	0.400	0.387	3.3	223	0.00
18	I	Phenanthrene-d10	0.400	0.400	0.0	169	0.00
19		4-Bromophenyl-phenylether	0.400	0.401	-0.3	172	-0.01
20		Hexachlorobenzene	0.400	0.405	-1.3	173	0.00
21		Pentachlorophenol	0.400	0.434	-8.5	211	0.00
22		Phenanthrene	0.400	0.394	1.5	172	0.00
23		Anthracene	0.400	0.424	-6.0	192	0.00
24	SURR	Fluoranthene-d10	0.400	0.367	8.3	162	0.00
25		Fluoranthene	0.400	0.376	6.0	164	0.00
26	I	Chrysene-d12	0.400	0.400	0.0	153	-0.01
27		Pyrene	0.400	0.447	-11.7	176	0.00
28	S	Terphenyl-d14	0.400	0.421	-5.2	167	-0.01
29		Benzo(a)anthracene	0.400	0.422	-5.5	165	0.00
30		Chrysene	0.400	0.415	-3.7	165	0.00
31		Indeno(1,2,3-cd)pyrene	0.400	0.429	-7.2	173	0.00
32	I	Perylene-d12	0.400	0.400	0.0	169	0.00
33		Benzo(b)fluoranthene	0.400	0.392	2.0	168	0.00
34		Benzo(k)fluoranthene	0.400	0.382	4.5	161	0.00
35	C	Benzo(a)pyrene	0.400	0.412	-3.0	181	0.00
36		Dibenzo(a,h)anthracene	0.400	0.397	0.8	174	0.00
37		Benzo(g,h,i)perylene	0.400	0.399	0.3	170	0.00

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