Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN081319\
Data File : BN007130.D
Acq On : 13 Aug 2019 14:07
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Aug 14 04:48:06 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	134	0.00
2		n-Nitrosodimethylamine	0.400	0.332	17.0	142	0.00
3	S	2-Fluorophenol	0.400	0.422	-5.5	147	0.00
4	S	Phenol-d6	0.400	0.435	-8.7	157	0.00
5		bis(2-Chloroethyl)ether	0.400	0.404	-1.0	141	0.00
6	I	Naphthalene-d8	0.400	0.400	0.0	143	0.00
7	S	Nitrobenzene-d5	0.400	0.410	-2.5	160	0.00
8		Naphthalene	0.400	0.375	6.3	143	0.00
9		Hexachlorobutadiene	0.400	0.381	4.8	143	-0.01
10	SURR	2-Methylnaphthalene-d10	0.400	0.369	7.8	142	0.00
11		2-Methylnaphthalene	0.400	0.372	7.0	143	0.00
12	I	Acenaphthene-d10	0.400	0.400	0.0	137	0.00
13	S	2,4,6-Tribromophenol	0.400	0.335	16.3	157	0.00
14	S	2-Fluorobiphenyl	0.400	0.230	42.5#	71	0.00
15		Acenaphthylene	0.400	0.450	-12.5	163	0.00
16		Acenaphthene	0.400	0.387	3.3	141	-0.01
17		Fluorene	0.400	0.352	12.0	145	0.00
18	I	Phenanthrene-d10	0.400	0.400	0.0	122	0.00
19		4-Bromophenyl-phenylether	0.400	0.406	-1.5	126	-0.01
20		Hexachlorobenzene	0.400	0.408	-2.0	127	0.00
21		Pentachlorophenol	0.400	0.388	3.0	137	0.00
22		Phenanthrene	0.400	0.387	3.3	123	0.00
23		Anthracene	0.400	0.401	-0.3	132	-0.01
24	SURR	Fluoranthene-d10	0.400	0.358	10.5	115	0.00
25		Fluoranthene	0.400	0.368	8.0	117	0.00
26	I	Chrysene-d12	0.400	0.400	0.0	114	-0.01
27		Pyrene	0.400	0.410	-2.5	121	0.00
28	S	Terphenyl-d14	0.400	0.398	0.5	118	-0.01
29		Benzo(a)anthracene	0.400	0.392	2.0	114	0.00
30		Chrysene	0.400	0.395	1.3	117	-0.01
31		Indeno(1,2,3-cd)pyrene	0.400	0.425	-6.2	128	-0.01
32	I	Perylene-d12	0.400	0.400	0.0	126	0.00
33		Benzo(b)fluoranthene	0.400	0.384	4.0	123	0.00
34		Benzo(k)fluoranthene	0.400	0.372	7.0	117	0.00
35	C	Benzo(a)pyrene	0.400	0.392	2.0	129	-0.01
36		Dibenzo(a,h)anthracene	0.400	0.394	1.5	129	-0.01
37		Benzo(g,h,i)perylene	0.400	0.388	3.0	124	-0.01

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