Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN081219\
Data File : BN007091.D
Acq On : 12 Aug 2019 11:06
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Aug 12 12:13:05 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN081019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Aug 10 21:44:50 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	119	0.00
2		n-Nitrosodimethylamine	0.400	0.298	25.5#	113	0.00
3	S	2-Fluorophenol	0.400	0.348	13.0	108	0.00
4	S	Phenol-d6	0.400	0.344	14.0	110	0.00
5		bis(2-Chloroethyl)ether	0.400	0.383	4.3	119	0.00
6	I	Naphthalene-d8	0.400	0.400	0.0	121	0.00
7	S	Nitrobenzene-d5	0.400	0.342	14.5	113	0.00
8		Naphthalene	0.400	0.377	5.8	122	0.00
9		Hexachlorobutadiene	0.400	0.383	4.3	121	-0.01
10	SURR	2-Methylnaphthalene-d10	0.400	0.367	8.3	119	0.00
11		2-Methylnaphthalene	0.400	0.368	8.0	120	0.00
12	I	Acenaphthene-d10	0.400	0.400	0.0	116	0.00
13	S	2,4,6-Tribromophenol	0.400	0.248	38.0#	99	0.00
14	S	2-Fluorobiphenyl	0.400	0.252	37.0#	66	0.00
15		Acenaphthylene	0.400	0.362	9.5	112	0.00
16		Acenaphthene	0.400	0.377	5.8	117	-0.01
17		Fluorene	0.400	0.326	18.5	114	0.00
18	I	Phenanthrene-d10	0.400	0.400	0.0	110	0.00
19		4-Bromophenyl-phenylether	0.400	0.394	1.5	110	-0.01
20		Hexachlorobenzene	0.400	0.435	-8.7	121	0.00
21		Pentachlorophenol	0.400	0.282	29.5#	89	0.00
22		Phenanthrene	0.400	0.391	2.3	111	0.00
23		Anthracene	0.400	0.352	12.0	104	-0.01
24	SURR	Fluoranthene-d10	0.400	0.340	15.0	98	0.00
25		Fluoranthene	0.400	0.354	11.5	101	0.00
26	I	Chrysene-d12	0.400	0.400	0.0	99	-0.01
27		Pyrene	0.400	0.392	2.0	100	0.00
28	S	Terphenyl-d14	0.400	0.400	0.0	103	0.00
29		Benzo(a)anthracene	0.400	0.349	12.8	88	0.00
30		Chrysene	0.400	0.395	1.3	102	-0.01
31		Indeno(1,2,3-cd)pyrene	0.400	0.359	10.3	94	-0.01
32	I	Perylene-d12	0.400	0.400	0.0	95	0.00
33		Benzo(b)fluoranthene	0.400	0.393	1.8	95	0.00
34		Benzo(k)fluoranthene	0.400	0.402	-0.5	95	0.00
35	C	Benzo(a)pyrene	0.400	0.361	9.8	89	-0.01
36		Dibenzo(a,h)anthracene	0.400	0.388	3.0	96	-0.02
37		Benzo(g,h,i)perylene	0.400	0.389	2.8	94	-0.02

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