Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081019\
Data File : BN007089.D
Acq On : 11 Aug 2019 16:27
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Aug 12 04:44:24 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	139	0.00
2		n-Nitrosodimethylamine	0.400	0.325	18.8	144	0.00
3	S	2-Fluorophenol	0.400	0.359	10.3	130	0.00
4	S	Phenol-d6	0.400	0.363	9.3	136	0.00
5		bis(2-Chloroethyl)ether	0.400	0.379	5.3	138	0.00
6	I	Naphthalene-d8	0.400	0.400	0.0	146	0.00
7	S	Nitrobenzene-d5	0.400	0.346	13.5	138	0.00
8		Naphthalene	0.400	0.372	7.0	146	0.00
9		Hexachlorobutadiene	0.400	0.379	5.3	146	-0.01
10	SURR	2-Methylnaphthalene-d10	0.400	0.368	8.0	146	0.00
11		2-Methylnaphthalene	0.400	0.370	7.5	146	0.00
12	I	Acenaphthene-d10	0.400	0.400	0.0	144	0.00
13	S	2,4,6-Tribromophenol	0.400	0.257	35.8#	127	0.00
14	S	2-Fluorobiphenyl	0.400	0.275	31.3#	90	0.00
15		Acenaphthylene	0.400	0.364	9.0	139	0.00
16		Acenaphthene	0.400	0.376	6.0	145	-0.01
17		Fluorene	0.400	0.323	19.3	140	0.00
18	I	Phenanthrene-d10	0.400	0.400	0.0	134	0.00
19		4-Bromophenyl-phenylether	0.400	0.397	0.8	135	-0.01
20		Hexachlorobenzene	0.400	0.420	-5.0	142	0.00
21		Pentachlorophenol	0.400	0.283	29.3#	109	0.00
22		Phenanthrene	0.400	0.392	2.0	136	0.00
23		Anthracene	0.400	0.365	8.8	131	-0.01
24	SURR	Fluoranthene-d10	0.400	0.347	13.3	122	0.00
25		Fluoranthene	0.400	0.358	10.5	124	0.00
26	I	Chrysene-d12	0.400	0.400	0.0	123	-0.01
27		Pyrene	0.400	0.388	3.0	124	-0.01
28	S	Terphenyl-d14	0.400	0.391	2.3	126	0.00
29		Benzo(a)anthracene	0.400	0.360	10.0	114	-0.01
30		Chrysene	0.400	0.401	-0.3	129	-0.01
31		Indeno(1,2,3-cd)pyrene	0.400	0.386	3.5	126	-0.02
32	I	Perylene-d12	0.400	0.400	0.0	126	-0.02
33		Benzo(b)fluoranthene	0.400	0.387	3.3	124	-0.01
34		Benzo(k)fluoranthene	0.400	0.384	4.0	121	-0.01
35	C	Benzo(a)pyrene	0.400	0.362	9.5	118	-0.02
36		Dibenzo(a,h)anthracene	0.400	0.396	1.0	129	-0.02
37		Benzo(g,h,i)perylene	0.400	0.387	3.3	123	-0.02

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