Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN081219\
Data File : BN007108.D
Acq On : 12 Aug 2019 21:18
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Aug 13 04:17:23 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	135	0.00
2		n-Nitrosodimethylamine	0.400	0.366	8.5	158	0.00
3	S	2-Fluorophenol	0.400	0.409	-2.2	144	0.00
4	S	Phenol-d6	0.400	0.427	-6.7	155	0.00
5		bis(2-Chloroethyl)ether	0.400	0.407	-1.7	144	0.00
6	I	Naphthalene-d8	0.400	0.400	0.0	145	0.00
7	S	Nitrobenzene-d5	0.400	0.397	0.8	157	0.00
8		Naphthalene	0.400	0.375	6.3	146	0.00
9		Hexachlorobutadiene	0.400	0.380	5.0	145	-0.01
10	SURR	2-Methylnaphthalene-d10	0.400	0.372	7.0	146	0.00
11		2-Methylnaphthalene	0.400	0.373	6.8	146	0.00
12	I	Acenaphthene-d10	0.400	0.400	0.0	143	0.00
13	S	2,4,6-Tribromophenol	0.400	0.319	20.3	157	0.00
14	S	2-Fluorobiphenyl	0.400	0.227	43.3#	74	0.00
15		Acenaphthylene	0.400	0.401	-0.3	152	0.00
16		Acenaphthene	0.400	0.377	5.8	144	-0.01
17		Fluorene	0.400	0.332	17.0	143	0.00
18	I	Phenanthrene-d10	0.400	0.400	0.0	133	0.00
19		4-Bromophenyl-phenylether	0.400	0.394	1.5	133	-0.01
20		Hexachlorobenzene	0.400	0.402	-0.5	136	0.00
21		Pentachlorophenol	0.400	0.374	6.5	144	0.00
22		Phenanthrene	0.400	0.385	3.8	133	0.00
23		Anthracene	0.400	0.390	2.5	139	-0.01
24	SURR	Fluoranthene-d10	0.400	0.365	8.8	128	0.00
25		Fluoranthene	0.400	0.368	8.0	127	0.00
26	I	Chrysene-d12	0.400	0.400	0.0	124	-0.01
27		Pyrene	0.400	0.409	-2.2	132	-0.01
28	S	Terphenyl-d14	0.400	0.405	-1.3	131	-0.01
29		Benzo(a)anthracene	0.400	0.407	-1.7	130	-0.01
30		Chrysene	0.400	0.393	1.8	127	-0.01
31		Indeno(1,2,3-cd)pyrene	0.400	0.423	-5.7	139	-0.02
32	I	Perylene-d12	0.400	0.400	0.0	137	-0.01
33		Benzo(b)fluoranthene	0.400	0.385	3.8	134	-0.01
34		Benzo(k)fluoranthene	0.400	0.382	4.5	131	-0.01
35	C	Benzo(a)pyrene	0.400	0.394	1.5	140	-0.02
36		Dibenzo(a,h)anthracene	0.400	0.392	2.0	140	-0.02
37		Benzo(g,h,i)perylene	0.400	0.386	3.5	134	-0.02

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