Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121119\
Data File : BE100841.D
Acq On : 11 Dec 2019 10:20
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 11 11:44:49 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE120419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 04 17:18:11 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	117	-0.01
2		1,4-Dioxane	0.400	0.352	12.0	103	-0.01
3		n-Nitrosodimethylamine	0.400	0.390	2.5	130	-0.01
4	S	2-Fluorophenol	0.400	0.454	-13.5	141	-0.01
5	S	Phenol-d6	0.400	0.480	-20.0	151	0.00
6		bis(2-Chloroethyl)ether	0.400	0.396	1.0	122	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	131	0.00
8	S	Nitrobenzene-d5	0.400	0.516	-29.0#	194	0.00
9		Naphthalene	0.400	0.382	4.5	130	0.00
10		Hexachlorobutadiene	0.400	0.377	5.8	127	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.392	2.0	136	0.00
12		2-Methylnaphthalene	0.400	0.395	1.3	137	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	152	0.00
14	S	2,4,6-Tribromophenol	0.400	1.130	-182.5#	493	0.00
15	S	2-Fluorobiphenyl	0.400	0.335	16.3	133	-0.01
16		Acenaphthylene	0.400	0.376	6.0	166	0.00
17		Acenaphthene	0.400	0.389	2.8	163	0.00
18		Fluorene	0.400	0.487	-21.7	208	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	241	-0.01
20		4-Bromophenyl-phenylether	0.400	0.350	12.5	233	-0.01
21		Hexachlorobenzene	0.400	0.364	9.0	233	0.00
22		Pentachlorophenol	0.400	0.671	-67.8#	2091	0.00
23		Phenanthrene	0.400	0.379	5.3	243	0.00
24		Anthracene	0.400	0.382	4.5	268	0.00
25	SURR	Fluoranthene-d10	0.400	0.357	10.8	239	0.00
26		Fluoranthene	0.400	0.363	9.3	243	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	224	0.00
28		Pyrene	0.400	0.407	-1.7	243	0.00
29	S	Terphenyl-d14	0.400	0.392	2.0	230	0.00
30		Benzo(a)anthracene	0.400	0.421	-5.2	265	-0.01
31		Chrysene	0.400	0.376	6.0	218	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.655	-63.7#	381	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.456	-14.0	289	-0.03
34	I	Perylene-d12	0.400	0.400	0.0	270	-0.01
35		Benzo(b)fluoranthene	0.400	0.347	13.3	247	-0.01
36		Benzo(k)fluoranthene	0.400	0.358	10.5	255	-0.01
37	C	Benzo(a)pyrene	0.400	0.354	11.5	262	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.380	5.0	289	-0.03
39		Benzo(g,h,i)perylene	0.400	0.369	7.8	265	-0.02

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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