Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE070119\
Data File : BE100276.D
Acq On : 1 Jul 2019 21:30
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 02 02:07:11 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE062819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 28 14:06:53 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	252	0.00
2		1,4-Dioxane	0.400	0.258	35.5#	157	0.01
3		n-Nitrosodimethylamine	0.400	0.544	-36.0#	416	-0.02
4	S	2-Fluorophenol	0.400	0.373	6.8	234	0.00
5	S	Phenol-d6	0.400	0.526	-31.5#	333	0.00
6		bis(2-Chloroethyl)ether	0.400	0.596	-49.0#	339	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	343	-0.01
8	S	Nitrobenzene-d5	0.400	0.427	-6.7	372	-0.03
9		Naphthalene	0.400	0.406	-1.5	342	0.00
10		Hexachlorobutadiene	0.400	0.407	-1.7	338	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.437	-9.2	369	-0.01
12		2-Methylnaphthalene	0.400	0.415	-3.7	373	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	365	0.00
14	S	2,4,6-Tribromophenol	0.400	0.441	-10.2	409	-0.01
15	S	2-Fluorobiphenyl	0.400	0.413	-3.2	366	-0.01
16		Acenaphthylene	0.400	0.417	-4.2	367	0.00
17		Acenaphthene	0.400	0.410	-2.5	354	0.00
18		Fluorene	0.400	0.402	-0.5	354	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	318	0.00
20		4-Bromophenyl-phenylether	0.400	0.439	-9.7	335	-0.01
21		Hexachlorobenzene	0.400	0.422	-5.5	321	0.01
22		Pentachlorophenol	0.400	0.509	-27.2#	452	-0.04
23		Phenanthrene	0.400	0.423	-5.7	335	-0.01
24		Anthracene	0.400	0.445	-11.2	358	-0.01
25	SURR	Fluoranthene-d10	0.400	0.393	1.8	309	0.00
26		Fluoranthene	0.400	0.391	2.3	303	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	261	0.00
28		Pyrene	0.400	0.483	-20.7	308	0.00
29	S	Terphenyl-d14	0.400	0.487	-21.7	314	0.00
30		Benzo(a)anthracene	0.400	0.445	-11.2	284	0.01
31		Chrysene	0.400	0.436	-9.0	277	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.422	-5.5	265	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.400	0.0	241	0.01
34	I	Perylene-d12	0.400	0.400	0.0	240	0.00
35		Benzo(b)fluoranthene	0.400	0.429	-7.2	254	0.00
36		Benzo(k)fluoranthene	0.400	0.402	-0.5	232	0.00
37	C	Benzo(a)pyrene	0.400	0.417	-4.2	241	0.00
38		Dibenzo(a,h)anthracene	0.400	0.424	-6.0	252	0.01
39		Benzo(g,h,i)perylene	0.400	0.419	-4.7	242	0.01

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

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