Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE010219\
Data File : BE098600.D
Acq On : 2 Jan 2019 9:44
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jan 02 10:34:04 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE121018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 10 14:59:55 2018
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	63	0.00
2		1,4-Dioxane	0.400	0.454	-13.5	72	0.00
3		n-Nitrosodimethylamine	0.400	0.420	-5.0	71	0.00
4	S	2-Fluorophenol	0.400	0.473	-18.2	76	0.01
5	S	Phenol-d6	0.400	0.524	-31.0#	82	0.01
6		bis(2-Chloroethyl)ether	0.400	0.389	2.8	63	0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	66	0.00
8	S	Nitrobenzene-d5	0.400	0.416	-4.0	74	0.00
9		Naphthalene	0.400	0.420	-5.0	71	0.00
10		Hexachlorobutadiene	0.400	0.455	-13.7	77	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.431	-7.7	72	0.00
12		2-Methylnaphthalene	0.400	0.420	-5.0	71	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	66	0.00
14	S	2,4,6-Tribromophenol	0.400	0.532	-33.0#	92	0.01
15	S	2-Fluorobiphenyl	0.400	0.393	1.8	71	0.02
16		Acenaphthylene	0.400	0.382	4.5	65	0.02
17		Acenaphthene	0.400	0.415	-3.7	69	0.02
18		Fluorene	0.400	0.406	-1.5	67	0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	58	0.00
20		4-Bromophenyl-phenylether	0.400	0.419	-4.7	63	0.01
21		Hexachlorobenzene	0.400	0.448	-12.0	66	0.01
22		Pentachlorophenol	0.400	1.021	-155.2#	183	0.00
23		Phenanthrene	0.400	0.415	-3.7	61	0.01
24		Anthracene	0.400	0.399	0.3	59	0.01
25	SURR	Fluoranthene-d10	0.400	0.415	-3.7	58	0.00
26		Fluoranthene	0.400	0.402	-0.5	56	0.01
27	I	Chrysene-d12	0.400	0.400	0.0	63	0.01
28		Pyrene	0.400	0.379	5.3	58	0.01
29	S	Terphenyl-d14	0.400	0.391	2.3	62	0.00
30		Benzo(a)anthracene	0.400	0.424	-6.0	69	0.00
31		Chrysene	0.400	0.393	1.8	63	0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.491	-22.7	78	0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.413	-3.2	69	-0.06
34	I	Perylene-d12	0.400	0.400	0.0	66	-0.03
35		Benzo(b)fluoranthene	0.400	0.427	-6.7	73	-0.01
36		Benzo(k)fluoranthene	0.400	0.398	0.5	69	-0.02
37	C	Benzo(a)pyrene	0.400	0.408	-2.0	69	-0.03
38		Dibenzo(a,h)anthracene	0.400	0.387	3.3	66	-0.08
39		Benzo(g,h,i)perylene	0.400	0.405	-1.3	69	-0.08

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

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