Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE122118\
Data File : BE098547.D
Acq On : 21 Dec 2018 9:33
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 21 15:42:05 2018
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	47	-0.01
2		1,4-Dioxane	0.400	0.397	0.8	48	-0.01
3		n-Nitrosodimethylamine	0.400	0.411	-2.7	52	0.00
4	S	2-Fluorophenol	0.400	0.429	-7.2	52	0.00
5	S	Phenol-d6	0.400	0.398	0.5	47	-0.01
6		bis(2-Chloroethyl)ether	0.400	0.357	10.8	44	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	55	-0.01
8	S	Nitrobenzene-d5	0.400	0.339	15.3	51	0.00
9		Naphthalene	0.400	0.379	5.3	53	-0.01
10		Hexachlorobutadiene	0.400	0.408	-2.0	57	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.407	-1.7	56	-0.01
12		2-Methylnaphthalene	0.400	0.392	2.0	55	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	66	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.434	-8.5	75	-0.01
15	S	2-Fluorobiphenyl	0.400	0.328	18.0	59	-0.01
16		Acenaphthylene	0.400	0.342	14.5	58	-0.01
17		Acenaphthene	0.400	0.366	8.5	61	-0.01
18		Fluorene	0.400	0.372	7.0	62	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	66	-0.01
20		4-Bromophenyl-phenylether	0.400	0.354	11.5	61	-0.01
21		Hexachlorobenzene	0.400	0.381	4.8	64	-0.01
22		Pentachlorophenol	0.400	0.571	-42.7#	102	-0.01
23		Phenanthrene	0.400	0.363	9.3	61	-0.01
24		Anthracene	0.400	0.335	16.3	56	-0.01
25	SURR	Fluoranthene-d10	0.400	0.407	-1.7	64	-0.01
26		Fluoranthene	0.400	0.395	1.3	62	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	73	-0.01
28		Pyrene	0.400	0.355	11.3	63	-0.01
29	S	Terphenyl-d14	0.400	0.356	11.0	65	-0.01
30		Benzo(a)anthracene	0.400	0.348	13.0	65	-0.02
31		Chrysene	0.400	0.362	9.5	66	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.327	18.3	60	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.335	16.3	64	-0.03
34	I	Perylene-d12	0.400	0.400	0.0	67	-0.02
35		Benzo(b)fluoranthene	0.400	0.366	8.5	64	-0.02
36		Benzo(k)fluoranthene	0.400	0.369	7.8	65	-0.02
37	C	Benzo(a)pyrene	0.400	0.363	9.3	62	-0.03
38		Dibenzo(a,h)anthracene	0.400	0.368	8.0	64	-0.04
39		Benzo(g,h,i)perylene	0.400	0.371	7.3	64	-0.04

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

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