Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121518\
Data File : BE098487.D
Acq On : 16 Dec 2018 1:05
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 16 03:03:45 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	55	0.00
2		1,4-Dioxane	0.400	0.356	11.0	49	0.00
3		n-Nitrosodimethylamine	0.400	0.352	12.0	52	0.01
4	S	2-Fluorophenol	0.400	0.459	-14.8	65	0.01
5	S	Phenol-d6	0.400	0.446	-11.5	61	0.00
6		bis(2-Chloroethyl)ether	0.400	0.322	19.5	45	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	55	-0.01
8	S	Nitrobenzene-d5	0.400	0.378	5.5	57	0.01
9		Naphthalene	0.400	0.417	-4.2	59	0.00
10		Hexachlorobutadiene	0.400	0.479	-19.7	68	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.406	-1.5	56	-0.01
12		2-Methylnaphthalene	0.400	0.382	4.5	54	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	53	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.445	-11.2	62	0.00
15	S	2-Fluorobiphenyl	0.400	0.418	-4.5	60	0.00
16		Acenaphthylene	0.400	0.422	-5.5	57	0.00
17		Acenaphthene	0.400	0.405	-1.3	54	0.00
18		Fluorene	0.400	0.409	-2.2	54	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	50	-0.01
20		4-Bromophenyl-phenylether	0.400	0.376	6.0	49	0.00
21		Hexachlorobenzene	0.400	0.433	-8.2	55	0.00
22		Pentachlorophenol	0.400	0.505	-26.2#	67	0.00
23		Phenanthrene	0.400	0.407	-1.7	52	0.00
24		Anthracene	0.400	0.353	11.8	45	0.00
25	SURR	Fluoranthene-d10	0.400	0.412	-3.0	50	0.00
26		Fluoranthene	0.400	0.417	-4.2	50	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	57	-0.01
28		Pyrene	0.400	0.387	3.3	54	0.00
29	S	Terphenyl-d14	0.400	0.334	16.5	48	0.00
30		Benzo(a)anthracene	0.400	0.350	12.5	51	-0.01
31		Chrysene	0.400	0.413	-3.2	59	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.256	36.0#	37	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.399	0.3	60	0.00
34	I	Perylene-d12	0.400	0.400	0.0	57	-0.01
35		Benzo(b)fluoranthene	0.400	0.427	-6.7	63	0.00
36		Benzo(k)fluoranthene	0.400	0.463	-15.8	69	-0.01
37	C	Benzo(a)pyrene	0.400	0.407	-1.7	59	0.00
38		Dibenzo(a,h)anthracene	0.400	0.370	7.5	54	-0.01
39		Benzo(g,h,i)perylene	0.400	0.422	-5.5	62	-0.01

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

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