Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\Data\BE100317\
Data File : BE094159.D
Acq On : 3 Oct 2017 17:55
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Oct 04 13:06:20 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 29 13:57:50 2017
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	130	0.00
2		1,4-Dioxane	0.400	0.370	7.5	116	0.00
3		n-Nitrosodimethylamine	0.400	0.367	8.3	120	0.00
4	S	2-Fluorophenol	0.400	0.348	13.0	112	0.00
5	S	Phenol-d6	0.400	0.362	9.5	117	0.00
6		bis(2-Chloroethyl)ether	0.400	0.374	6.5	119	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	137	0.00
8	S	Nitrobenzene-d5	0.400	0.322	19.5	117	0.00
9		Naphthalene	0.400	0.399	0.3	138	0.00
10		Hexachlorobutadiene	0.400	0.395	1.3	142	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.401	-0.3	140	0.00
12		2-Methylnaphthalene	0.400	0.402	-0.5	142	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	142	0.00
14	S	2,4,6-Tribromophenol	0.400	0.210	47.5#	85	0.00
15	S	2-Fluorobiphenyl	0.400	0.413	-3.2	140	0.00
16		Acenaphthylene	0.400	0.395	1.3	136	0.00
17		Acenaphthene	0.400	0.428	-7.0	141	0.00
18		Fluorene	0.400	0.408	-2.0	140	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	115	0.00
20		4-Bromophenyl-phenylether	0.400	0.485	-21.2	143	0.00
21		Hexachlorobenzene	0.400	0.626	-56.5#	182	0.00
22		Pentachlorophenol	0.400	0.184	54.0#	54	0.00
23		Phenanthrene	0.400	0.480	-20.0	144	0.00
24		Anthracene	0.400	0.384	4.0	117	0.00
25	SURR	Fluoranthene-d10	0.400	0.376	6.0	113	0.00
26		Fluoranthene	0.400	0.385	3.8	114	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	104	0.00
28		Pyrene	0.400	0.411	-2.7	111	0.00
29	S	Terphenyl-d14	0.400	0.388	3.0	106	0.00
30		Benzo(a)anthracene	0.400	0.378	5.5	98	0.00
31		Chrysene	0.400	0.389	2.8	101	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.088	78.0#	63	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.371	7.3	100	0.00
34	I	Perylene-d12	0.400	0.400	0.0	100	-0.01
35		Benzo(b)fluoranthene	0.400	0.369	7.8	103	0.00
36		Benzo(k)fluoranthene	0.400	0.383	4.3	96	0.00
37	C	Benzo(a)pyrene	0.400	0.385	3.8	100	0.00
38		Dibenzo(a,h)anthracene	0.400	0.372	7.0	98	0.00
39		Benzo(g,h,i)perylene	0.400	0.387	3.3	101	0.00

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

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