Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE110117\
Data File : BE094687.D
Acq On : 1 Nov 2017 12:39
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Nov 01 14:09:30 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 23 22:40:47 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	107	0.00
2		1,4-Dioxane	0.400	0.386	3.5	90	-0.01
3		n-Nitrosodimethylamine	0.400	0.297	25.8#	79	0.01
4	S	2-Fluorophenol	0.400	0.324	19.0	87	0.04
5	S	Phenol-d6	0.400	0.377	5.8	104	0.04
6		bis(2-Chloroethyl)ether	0.400	0.346	13.5	94	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	113	0.00
8	S	Nitrobenzene-d5	0.400	0.338	15.5	99	0.02
9		Naphthalene	0.400	0.354	11.5	97	0.02
10		Hexachlorobutadiene	0.400	0.391	2.3	108	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.344	14.0	95	0.00
12		2-Methylnaphthalene	0.400	0.339	15.3	95	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	101	0.01
14	S	2,4,6-Tribromophenol	0.400	0.334	16.5	88	0.03
15	S	2-Fluorobiphenyl	0.400	0.371	7.3	93	0.01
16		Acenaphthylene	0.400	0.346	13.5	88	0.00
17		Acenaphthene	0.400	0.358	10.5	89	0.00
18		Fluorene	0.400	0.344	14.0	87	0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	95	0.03
20		4-Bromophenyl-phenylether	0.400	0.323	19.3	79	0.03
21		Hexachlorobenzene	0.400	0.583	-45.7#	121	0.00
22		Pentachlorophenol	0.400	0.427	-6.7	102	0.06
23		Phenanthrene	0.400	0.564	-41.0#	119	0.00
24		Anthracene	0.400	0.379	5.3	89	0.00
25	SURR	Fluoranthene-d10	0.400	0.426	-6.5	92	0.00
26		Fluoranthene	0.400	0.414	-3.5	90	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	100	0.00
28		Pyrene	0.400	0.360	10.0	91	0.00
29	S	Terphenyl-d14	0.400	0.353	11.8	90	0.00
30		Benzo(a)anthracene	0.400	0.360	10.0	90	0.01
31		Chrysene	0.400	0.365	8.8	94	0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.352	12.0	96	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.298	25.5#	75	0.05
34	I	Perylene-d12	0.400	0.400	0.0	97	0.02
35		Benzo(b)fluoranthene	0.400	0.347	13.3	87	0.01
36		Benzo(k)fluoranthene	0.400	0.367	8.3	90	0.02
37	C	Benzo(a)pyrene	0.400	0.372	7.0	91	0.02
38		Dibenzo(a,h)anthracene	0.400	0.322	19.5	80	0.03
39		Benzo(g,h,i)perylene	0.400	0.284	29.0#	70	0.06

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

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