Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_E\Data\BE050517\
Data File : BE092957.D
Acq On : 6 May 2017 8:12
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 08 04:29:46 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE050517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 05 14:05:07 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	90	0.00
2		1,4-Dioxane	0.400	0.379	5.3	89	0.00
3		n-Nitrosodimethylamine	0.400	0.472	-18.0	115	0.00
4	S	2-Fluorophenol	0.400	0.573	-43.2#	142	0.01
5	S	Phenol-d6	0.400	0.619	-54.7#	154	0.00
6		bis(2-Chloroethyl)ether	0.400	0.422	-5.5	99	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	104	0.00
8	S	Nitrobenzene-d5	0.400	0.540	-35.0#	155	0.00
9		Naphthalene	0.400	0.376	6.0	104	0.00
10		Hexachlorobutadiene	0.400	0.366	8.5	100	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.391	2.3	109	0.00
12		2-Methylnaphthalene	0.400	0.396	1.0	110	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	118	0.00
14	S	2,4,6-Tribromophenol	0.400	0.802	-100.5#	277	0.01
15	S	2-Fluorobiphenyl	0.400	0.347	13.3	107	0.00
16		Acenaphthylene	0.400	0.458	-14.5	153	0.00
17		Acenaphthene	0.400	0.400	0.0	127	0.00
18		Fluorene	0.400	0.399	0.3	125	0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	111	0.00
20		Hexachlorobenzene	0.400	0.409	-2.2	122	0.00
21		Pentachlorophenol	0.400	0.825	-106.2#	293	0.02
22		Phenanthrene	0.400	0.412	-3.0	119	0.00
23		Anthracene	0.400	0.502	-25.5#	146	0.00
24	SURR	Fluoranthene-d10	0.400	0.498	-24.5	146	0.00
25		Fluoranthene	0.400	0.477	-19.2	139	0.00
26	I	Chrysene-d12	0.400	0.400	0.0	140	0.00
27		Pyrene	0.400	0.396	1.0	143	0.00
28	S	Terphenyl-d14	0.400	0.336	16.0	122	0.00
29		Benzo(a)anthracene	0.400	0.465	-16.3	178	0.02
30		Chrysene	0.400	0.364	9.0	132	0.02
31		Bis(2-ethylhexyl)phthalate	0.400	1.521	-280.3#	546	0.00
32		Indeno(1,2,3-cd)pyrene	0.400	0.403	-0.8	148	0.02
33	I	Perylene-d12	0.400	0.400	0.0	136	0.01
34		Benzo(b)fluoranthene	0.400	0.385	3.8	158	0.01
35		Benzo(k)fluoranthene	0.400	0.338	15.5	129	0.01
36	C	Benzo(a)pyrene	0.400	0.415	-3.7	172	0.00
37		Dibenzo(a,h)anthracene	0.400	0.350	12.5	144	0.00
38		Benzo(g,h,i)perylene	0.400	0.347	13.3	141	0.02

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