Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE050917\
Data File : BE092960.D
Acq On : 9 May 2017 11:59
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 09 17:12:59 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE050517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue May 09 15:34:16 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	88	0.00
2		1,4-Dioxane	0.400	0.416	-4.0	96	0.00
3		n-Nitrosodimethylamine	0.400	0.498	-24.5	119	0.00
4	S	2-Fluorophenol	0.400	0.530	-32.5#	128	0.00
5	S	Phenol-d6	0.400	0.504	-26.0#	123	0.00
6		bis(2-Chloroethyl)ether	0.400	0.440	-10.0	102	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	95	0.00
8	S	Nitrobenzene-d5	0.400	0.452	-13.0	119	0.02
9		Naphthalene	0.400	0.394	1.5	99	0.00
10		Hexachlorobutadiene	0.400	0.398	0.5	100	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.381	4.8	97	0.00
12		2-Methylnaphthalene	0.400	0.392	2.0	99	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	105	0.00
14	S	2,4,6-Tribromophenol	0.400	0.601	-50.2#	184	0.01
15	S	2-Fluorobiphenyl	0.400	0.360	10.0	99	0.01
16		Acenaphthylene	0.400	0.417	-4.2	124	0.00
17		Acenaphthene	0.400	0.398	0.5	112	0.00
18		Fluorene	0.400	0.404	-1.0	112	0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	101	0.00
20		Hexachlorobenzene	0.400	0.438	-9.5	119	0.00
21		Pentachlorophenol	0.400	0.559	-39.8#	161	0.02
22		Phenanthrene	0.400	0.424	-6.0	111	0.00
23		Anthracene	0.400	0.461	-15.3	122	0.00
24	SURR	Fluoranthene-d10	0.400	0.469	-17.2	125	0.00
25		Fluoranthene	0.400	0.445	-11.2	118	0.00
26	I	Chrysene-d12	0.400	0.400	0.0	120	0.00
27		Pyrene	0.400	0.379	5.3	118	0.00
28	S	Terphenyl-d14	0.400	0.342	14.5	106	0.00
29		Benzo(a)anthracene	0.400	0.429	-7.2	141	0.02
30		Chrysene	0.400	0.401	-0.3	124	0.02
31		Bis(2-ethylhexyl)phthalate	0.400	0.650	-62.5#	200	0.00
32		Indeno(1,2,3-cd)pyrene	0.400	0.628	-57.0#	197	0.02
33	I	Perylene-d12	0.400	0.400	0.0	125	0.01
34		Benzo(b)fluoranthene	0.400	0.472	-18.0	177	0.01
35		Benzo(k)fluoranthene	0.400	0.456	-14.0	160	0.01
36	C	Benzo(a)pyrene	0.400	0.452	-13.0	173	0.00
37		Dibenzo(a,h)anthracene	0.400	0.494	-23.5	187	0.02
38		Benzo(g,h,i)perylene	0.400	0.558	-39.5#	209	0.02

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