Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032917\
Data File : BE092861.D
Acq On : 29 Mar 2017 9:37
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 30 01:41:16 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE032817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 29 17:47:56 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	5.000	5.000	0.0	102	-0.02
2		1,4-Dioxane	0.400	0.407	-1.7	101	-0.01
3		n-Nitrosodimethylamine	0.400	0.393	1.8	107	0.00
4	S	2-Fluorophenol	0.400	0.383	4.3	99	0.00
5	S	Phenol-d6	0.400	0.378	5.5	103	0.00
6		bis(2-Chloroethyl)ether	0.400	0.414	-3.5	107	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	103	0.00
8	S	Nitrobenzene-d5	0.400	0.398	0.5	114	0.00
9		Naphthalene	0.400	0.406	-1.5	104	0.00
10		Hexachlorobutadiene	0.400	0.393	1.8	102	0.00
11		2-Methylnaphthalene	0.400	0.384	4.0	101	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	104	0.00
13	S	2,4,6-Tribromophenol	0.400	0.391	2.3	113	0.00
14	S	2-Fluorobiphenyl	0.400	0.403	-0.8	104	-0.01
15		Acenaphthylene	0.400	0.348	13.0	101	-0.02
16		Acenaphthene	0.400	0.381	4.8	103	0.00
17		Fluorene	0.400	0.377	5.8	99	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	101	-0.02
19		Hexachlorobenzene	0.400	0.432	-8.0	112	0.00
20		Pentachlorophenol	0.400	0.511	-27.7#	109	-0.02
21		Phenanthrene	0.400	0.473	-18.2	116	0.00
22		Anthracene	0.400	0.393	1.8	108	0.00
23		Fluoranthene	0.400	0.342	14.5	93	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	101	-0.02
25		Pyrene	0.400	0.346	13.5	93	0.00
26	S	Terphenyl-d14	0.400	0.343	14.2	90	-0.02
27		Benzo(a)anthracene	0.400	0.376	6.0	104	0.00
28		Chrysene	0.400	0.386	3.5	99	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.332	17.0	85	-0.02
30		Indeno(1,2,3-cd)pyrene	0.400	0.335	16.3	91	-0.02
31	I	Perylene-d12	5.000	5.000	0.0	101	-0.01
32		Benzo(b)fluoranthene	0.400	0.338	15.5	88	-0.01
33		Benzo(k)fluoranthene	0.400	0.350	12.5	94	-0.01
34	C	Benzo(a)pyrene	0.400	0.321	19.8	87	-0.01
35		Dibenzo(a,h)anthracene	0.400	0.343	14.2	92	-0.02
36		Benzo(g,h,i)perylene	0.400	0.350	12.5	91	-0.02

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