Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091216\
Data File : BE092369.D
Acq On : 12 Sep 2016 10:20
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Sep 12 14:26:45 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 02 19:34:51 2016
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	112	-0.02
2		1,4-Dioxane	0.400	0.359	10.3	105	-0.01
3		n-Nitrosodimethylamine	0.400	0.373	6.8	120	-0.01
4	S	2-Fluorophenol	0.400	0.335	16.3	110	-0.01
5	S	Phenol-d6	0.400	0.326	18.5	118	0.00
6		bis(2-Chloroethyl)ether	0.400	0.362	9.5	123	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	116	0.00
8	S	Nitrobenzene-d5	0.400	0.366	8.5	135	0.00
9		Naphthalene	0.400	0.403	-0.8	123	-0.02
10		Hexachlorobutadiene	0.400	0.406	-1.5	120	-0.02
11		2-Methylnaphthalene	0.400	0.409	-2.2	128	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	120	0.00
13	S	2,4,6-Tribromophenol	0.400	0.341	14.8	127	-0.01
14	S	2-Fluorobiphenyl	0.400	0.382	4.5	127	0.00
15		Acenaphthylene	0.400	0.389	2.8	130	0.00
16		Acenaphthene	0.400	0.395	1.3	129	0.00
17		Fluorene	0.400	0.391	2.3	130	-0.01
18	I	Phenanthrene-d10	5.000	5.000	0.0	128	0.00
19		Hexachlorobenzene	0.400	0.359	10.3	126	0.00
20		Pentachlorophenol	0.400	0.400	0.0	158	-0.02
21		Phenanthrene	0.400	0.373	6.8	131	0.00
22		Anthracene	0.400	0.367	8.3	142	0.00
23		Fluoranthene	0.400	0.330	17.5	120	-0.02
24	I	Chrysene-d12	5.000	5.000	0.0	106	0.00
25		Pyrene	0.400	0.437	-9.2	122	0.00
26	S	Terphenyl-d14	0.400	0.425	-6.2	120	0.00
27		Benzo(a)anthracene	0.400	0.379	5.3	103	0.00
28		Chrysene	0.400	0.391	2.3	130	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.419	-4.7	142	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.387	3.3	104	0.00
31	I	Perylene-d12	5.000	5.000	0.0	98	0.01
32		Benzo(b)fluoranthene	0.400	0.389	2.8	108	0.01
33		Benzo(k)fluoranthene	0.400	0.392	2.0	100	0.01
34	C	Benzo(a)pyrene	0.400	0.381	4.8	105	0.01
35		Dibenzo(a,h)anthracene	0.400	0.379	5.3	105	0.02
36		Benzo(g,h,i)perylene	0.400	0.396	1.0	105	0.00

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