Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090916\
Data File : BE092356.D
Acq On : 9 Sep 2016 15:57
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Sep 09 17:24:14 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	160	-0.02
2		1,4-Dioxane	0.400	0.335	16.3	142	0.00
3		n-Nitrosodimethylamine	0.400	0.373	6.8	173	0.00
4	S	2-Fluorophenol	0.400	0.310	22.5	146	0.00
5	S	Phenol-d6	0.400	0.329	17.8	171	0.00
6		bis(2-Chloroethyl)ether	0.400	0.351	12.3	169	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	161	0.00
8	S	Nitrobenzene-d5	0.400	0.379	5.3	194	0.00
9		Naphthalene	0.400	0.376	6.0	159	0.00
10		Hexachlorobutadiene	0.400	0.384	4.0	157	0.00
11		2-Methylnaphthalene	0.400	0.363	9.3	158	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	156	0.00
13	S	2,4,6-Tribromophenol	0.400	0.325	18.8	158	0.00
14	S	2-Fluorobiphenyl	0.400	0.375	6.3	162	0.00
15		Acenaphthylene	0.400	0.357	10.8	155	0.00
16		Acenaphthene	0.400	0.368	8.0	156	0.00
17		Fluorene	0.400	0.354	11.5	153	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	155	0.00
19		Hexachlorobenzene	0.400	0.351	12.3	150	0.00
20		Pentachlorophenol	0.400	0.328	18.0	150	0.00
21		Phenanthrene	0.400	0.371	7.3	158	0.00
22		Anthracene	0.400	0.340	15.0	159	0.00
23		Fluoranthene	0.400	0.348	13.0	153	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	161	0.00
25		Pyrene	0.400	0.381	4.8	161	0.00
26	S	Terphenyl-d14	0.400	0.388	3.0	166	0.00
27		Benzo(a)anthracene	0.400	0.341	14.8	140	0.00
28		Chrysene	0.400	0.345	13.8	180	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.311	22.3	146	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.353	11.8	145	0.02
31	I	Perylene-d12	5.000	5.000	0.0	149	0.01
32		Benzo(b)fluoranthene	0.400	0.352	12.0	148	0.01
33		Benzo(k)fluoranthene	0.400	0.363	9.3	140	0.01
34	C	Benzo(a)pyrene	0.400	0.355	11.3	148	0.01
35		Dibenzo(a,h)anthracene	0.400	0.343	14.2	144	0.02
36		Benzo(g,h,i)perylene	0.400	0.350	12.5	140	0.02

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