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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	102	-0.02
2		1,4-Dioxane	0.668	0.680	-1.8	101	-0.01
3		n-Nitrosodimethylamine	0.691	0.678	1.9	107	0.00
4	S	2-Fluorophenol	1.164	1.115	4.2	99	0.00
5	S	Phenol-d6	1.708	1.616	5.4	103	0.00
6		bis(2-Chloroethyl)ether	1.625	1.682	-3.5	107	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
8	S	Nitrobenzene-d5	0.287	0.285	0.7	114	0.00
9		Naphthalene	1.095	1.113	-1.6	104	0.00
10		Hexachlorobutadiene	0.147	0.144	2.0	102	0.00
11		2-Methylnaphthalene	0.695	0.667	4.0	101	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	104	0.00
13	S	2,4,6-Tribromophenol	0.078	0.076	2.6	113	0.00
14	S	2-Fluorobiphenyl	1.648	1.660	-0.7	104	-0.01
15		Acenaphthylene	7.666	6.662	13.1	101	-0.02
16		Acenaphthene	1.367	1.302	4.8	103	0.00
17		Fluorene	1.629	1.535	5.8	99	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	101	-0.02
19		Hexachlorobenzene	0.185	0.200	-8.1	112	0.00
20		Pentachlorophenol	0.048	0.044	8.3	109	-0.02
21		Phenanthrene	1.217	1.437	-18.1	116	0.00
22		Anthracene	1.106	1.087	1.7	108	0.00
23		Fluoranthene	5.597	4.791	14.4	93	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	101	-0.02
25		Pyrene	4.972	4.297	13.6	93	0.00
26	S	Terphenyl-d14	0.864	0.742	14.1	90	-0.02
27		Benzo(a)anthracene	1.185	1.113	6.1	104	0.00
28		Chrysene	1.250	1.207	3.4	99	0.00
29		Bis(2-ethylhexyl)phthalate	0.362	0.242	33.1#	85	-0.02
30		Indeno(1,2,3-cd)pyrene	4.702	3.941	16.2	91	-0.02
31	I	Perylene-d12	1.000	1.000	0.0	101	-0.01
32		Benzo(b)fluoranthene	1.331	1.125	15.5	88	-0.01
33		Benzo(k)fluoranthene	1.192	1.044	12.4	94	-0.01
34	C	Benzo(a)pyrene	1.100	0.882	19.8	87	-0.01
35		Dibenzo(a,h)anthracene	1.170	1.004	14.2	92	-0.02
36		Benzo(g,h,i)perylene	4.851	4.242	12.6	91	-0.02

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