Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\Data\BE012517\
Data File : BE092687.D
Acq On : 25 Jan 2017 17:48
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jan 25 18:31:28 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE011617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 17 10:14:01 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	172#	0.00
2		1,4-Dioxane	0.775	0.674	13.0	142	0.00
3		n-Nitrosodimethylamine	0.902	0.669	25.8#	146	0.01
4	S	2-Fluorophenol	1.116	1.070	4.1	194#	0.00
5	S	Phenol-d6	1.368	1.495	-9.3	229#	-0.02
6		bis(2-Chloroethyl)ether	1.453	1.299	10.6	183#	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	194#	0.00
8	S	Nitrobenzene-d5	0.325	0.286	12.0	205#	-0.02
9		Naphthalene	1.044	1.031	1.2	207#	-0.02
10		Hexachlorobutadiene	0.154	0.153	0.6	203#	-0.02
11		2-Methylnaphthalene	0.648	0.640	1.2	214#	-0.02
12	I	Acenaphthene-d10	1.000	1.000	0.0	194#	0.00
13	S	2,4,6-Tribromophenol	0.126	0.119	5.6	229#	0.01
14	S	2-Fluorobiphenyl	1.230	1.205	2.0	204#	-0.02
15		Acenaphthylene	6.890	6.988	-1.4	221#	-0.02
16		Acenaphthene	1.105	1.121	-1.4	207#	0.00
17		Fluorene	1.387	1.359	2.0	212#	-0.01
18	I	Phenanthrene-d10	1.000	1.000	0.0	190#	0.00
19		Hexachlorobenzene	0.189	0.177	6.3	171#	0.00
20		Pentachlorophenol	0.051	0.041	19.6	165#	0.05
21		Phenanthrene	1.104	1.111	-0.6	198#	0.00
22		Anthracene	1.018	1.088	-6.9	228#	-0.02
23		Fluoranthene	4.895	4.970	-1.5	213#	-0.02
24	I	Chrysene-d12	1.000	1.000	0.0	154#	0.00
25		Pyrene	4.929	5.629	-14.2	211#	-0.02
26	S	Terphenyl-d14	0.707	0.807	-14.1	205#	0.00
27		Benzo(a)anthracene	4.926	4.915	0.2	178#	0.00
28		Chrysene	5.542	5.835	-5.3	175#	0.00
29		Bis(2-ethylhexyl)phthalate	0.913	1.221	-33.7#	288#	0.00
30		Indeno(1,2,3-cd)pyrene	5.987	5.897	1.5	174#	-0.02
31	I	Perylene-d12	1.000	1.000	0.0	149	0.00
32		Benzo(b)fluoranthene	1.210	1.216	-0.5	163#	-0.01
33		Benzo(k)fluoranthene	1.215	1.302	-7.2	186#	-0.01
34	C	Benzo(a)pyrene	1.136	1.202	-5.8	183#	0.00
35		Dibenzo(a,h)anthracene	1.112	1.144	-2.9	173#	-0.02
36		Benzo(g,h,i)perylene	4.699	4.807	-2.3	169#	-0.02

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