Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE012417\
Data File : BE092682.D
Acq On : 25 Jan 2017 1:26
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jan 25 13:48:21 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	147	0.00
2		1,4-Dioxane	0.400	0.343	14.2	124	0.01
3		n-Nitrosodimethylamine	0.400	0.253	36.8#	107	0.03
4	S	2-Fluorophenol	0.400	0.346	13.5	150	0.00
5	S	Phenol-d6	0.400	0.391	2.3	175	0.00
6		bis(2-Chloroethyl)ether	0.400	0.335	16.3	147	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	165	0.00
8	S	Nitrobenzene-d5	0.400	0.343	14.2	171	-0.02
9		Naphthalene	0.400	0.389	2.8	174	-0.02
10		Hexachlorobutadiene	0.400	0.398	0.5	173	-0.02
11		2-Methylnaphthalene	0.400	0.393	1.8	182	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	174	0.00
13	S	2,4,6-Tribromophenol	0.400	0.425	-6.2	191	0.02
14	S	2-Fluorobiphenyl	0.400	0.377	5.8	177	-0.01
15		Acenaphthylene	0.400	0.401	-0.3	197	-0.02
16		Acenaphthene	0.400	0.402	-0.5	184	0.00
17		Fluorene	0.400	0.382	4.5	185	-0.01
18	I	Phenanthrene-d10	5.000	5.000	0.0	172	0.00
19		Hexachlorobenzene	0.400	0.359	10.3	159	0.00
20		Pentachlorophenol	0.400	0.248	38.0#	117	0.14
21		Phenanthrene	0.400	0.401	-0.3	179	0.00
22		Anthracene	0.400	0.414	-3.5	200	-0.02
23		Fluoranthene	0.400	0.374	6.5	177	-0.02
24	I	Chrysene-d12	5.000	5.000	0.0	146	0.00
25		Pyrene	0.400	0.406	-1.5	177	-0.02
26	S	Terphenyl-d14	0.400	0.402	-0.5	170	0.00
27		Benzo(a)anthracene	0.400	0.397	0.8	167	0.00
28		Chrysene	0.400	0.393	1.8	154	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.563	-40.7#	201	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.362	9.5	151	0.00
31	I	Perylene-d12	5.000	5.000	0.0	135	0.00
32		Benzo(b)fluoranthene	0.400	0.391	2.3	143	-0.01
33		Benzo(k)fluoranthene	0.400	0.422	-5.5	166	-0.01
34	C	Benzo(a)pyrene	0.400	0.406	-1.5	158	0.00
35		Dibenzo(a,h)anthracene	0.400	0.398	0.5	151	-0.02
36		Benzo(g,h,i)perylene	0.400	0.391	2.3	146	-0.02

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