Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE082916\
Data File : BE092300.D
Acq On : 29 Aug 2016 21:57
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Aug 30 01:10:02 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE082916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 29 18:19:23 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	107	0.00
2		1,4-Dioxane	0.400	0.385	3.8	101	0.00
3		n-Nitrosodimethylamine	0.400	0.359	10.3	110	-0.01
4	S	2-Fluorophenol	0.400	0.363	9.3	116	0.00
5	S	Phenol-d6	0.400	0.344	14.0	116	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.362	9.5	115	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	114	0.00
8	S	Nitrobenzene-d5	0.400	0.328	18.0	125	0.00
9		Naphthalene	0.400	0.370	7.5	106	0.00
10		Hexachlorobutadiene	0.400	0.378	5.5	107	0.00
11		2-Methylnaphthalene	0.400	0.372	7.0	111	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	121	0.00
13	S	2,4,6-Tribromophenol	0.400	0.377	5.8	126	0.00
14	S	2-Fluorobiphenyl	0.400	0.360	10.0	112	0.00
15		Acenaphthylene	0.400	0.362	9.5	116	0.00
16		Acenaphthene	0.400	0.372	7.0	115	0.00
17		Fluorene	0.400	0.365	8.8	115	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	116	0.00
19		Hexachlorobenzene	0.400	0.379	5.3	117	0.00
20		Pentachlorophenol	0.400	0.440	-10.0	132	0.00
21		Phenanthrene	0.400	0.358	10.5	109	0.00
22		Anthracene	0.400	0.359	10.3	112	0.00
23		Fluoranthene	0.400	0.374	6.5	118	-0.02
24	I	Chrysene-d12	5.000	5.000	0.0	123	0.00
25		Pyrene	0.400	0.372	7.0	121	-0.02
26	S	Terphenyl-d14	0.400	0.363	9.3	114	0.00
27		Benzo(a)anthracene	0.400	0.491	-22.7	136	0.00
28		Chrysene	0.400	0.394	1.5	113	-0.02
29		Bis(2-ethylhexyl)phthalate	0.400	0.462	-15.5	155	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.378	5.5	126	-0.02
31	I	Perylene-d12	5.000	5.000	0.0	119	-0.01
32		Benzo(b)fluoranthene	0.400	0.372	7.0	111	-0.01
33		Benzo(k)fluoranthene	0.400	0.382	4.5	121	-0.01
34	C	Benzo(a)pyrene	0.400	0.379	5.3	122	0.00
35		Dibenzo(a,h)anthracene	0.400	0.380	5.0	129	-0.03
36		Benzo(g,h,i)perylene	0.400	0.375	6.3	120	-0.02

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