Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081916\
Data File : BE092274.D
Acq On : 19 Aug 2016 17:37
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Aug 22 01:03:22 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 10 11:12:32 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	112	0.02
2		1,4-Dioxane	0.400	0.440	-10.0	118	0.00
3		n-Nitrosodimethylamine	0.400	0.481	-20.2	125	0.00
4	S	2-Fluorophenol	0.400	0.386	3.5	119	0.00
5	S	Phenol-d6	0.400	0.342	14.5	122	0.02
6		bis(2-Chloroethyl)ether	0.400	0.394	1.5	122	0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	112	0.00
8	S	Nitrobenzene-d5	0.400	0.403	-0.8	134	0.00
9		Naphthalene	0.400	0.411	-2.7	116	-0.02
10		Hexachlorobutadiene	0.400	0.424	-6.0	120	0.00
11		2-Methylnaphthalene	0.400	0.415	-3.7	119	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	110	0.00
13	S	2,4,6-Tribromophenol	0.400	0.495	-23.7	121	0.00
14	S	2-Fluorobiphenyl	0.400	0.413	-3.2	117	-0.01
15		Acenaphthylene	0.400	0.399	0.3	114	0.00
16		Acenaphthene	0.400	0.414	-3.5	116	0.00
17		Fluorene	0.400	0.412	-3.0	116	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	114	0.02
19		Hexachlorobenzene	0.400	0.406	-1.5	116	0.02
20		Pentachlorophenol	0.400	0.384	4.0	117	0.02
21		Phenanthrene	0.400	0.403	-0.8	119	0.02
22		Anthracene	0.400	0.378	5.5	117	0.02
23		Fluoranthene	0.400	0.379	5.3	114	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	114	0.02
25		Pyrene	0.400	0.400	0.0	119	0.00
26	S	Terphenyl-d14	0.400	0.390	2.5	120	0.00
27		Benzo(a)anthracene	0.400	0.389	2.8	119	0.02
28		Chrysene	0.400	0.404	-1.0	113	0.02
29		Bis(2-ethylhexyl)phthalate	0.400	0.453	-13.2	104	0.02
30		Indeno(1,2,3-cd)pyrene	0.400	0.366	8.5	115	0.00
31	I	Perylene-d12	5.000	5.000	0.0	111	0.00
32		Benzo(b)fluoranthene	0.400	0.381	4.8	115	0.00
33		Benzo(k)fluoranthene	0.400	0.412	-3.0	114	0.00
34	C	Benzo(a)pyrene	0.400	0.380	5.0	115	0.00
35		Dibenzo(a,h)anthracene	0.400	0.369	7.8	114	0.00
36		Benzo(g,h,i)perylene	0.400	0.386	3.5	113	0.00

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