Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE110916\
Data File : BE092519.D
Acq On : 9 Nov 2016 10:58
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Nov 09 14:29:43 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE110816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 08 16:27:22 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	114	-0.02
2		1,4-Dioxane	0.400	0.396	1.0	131	0.00
3		n-Nitrosodimethylamine	0.400	0.373	6.8	96	0.00
4	S	2-Fluorophenol	0.400	0.353	11.8	109	0.00
5	S	Phenol-d6	0.400	0.326	18.5	106	0.00
6		bis(2-Chloroethyl)ether	0.400	0.331	17.3	102	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	114	0.00
8	S	Nitrobenzene-d5	0.400	0.353	11.8	116	0.00
9		Naphthalene	0.400	0.382	4.5	108	0.00
10		Hexachlorobutadiene	0.400	0.393	1.8	109	-0.02
11		2-Methylnaphthalene	0.400	0.375	6.3	111	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	114	0.00
13	S	2,4,6-Tribromophenol	0.400	0.460	-15.0	110	0.00
14	S	2-Fluorobiphenyl	0.400	0.413	-3.2	121	0.00
15		Acenaphthylene	0.400	0.390	2.5	118	-0.02
16		Acenaphthene	0.400	0.395	1.3	116	0.00
17		Fluorene	0.400	0.392	2.0	117	-0.01
18	I	Phenanthrene-d10	5.000	5.000	0.0	117	0.00
19		Hexachlorobenzene	0.400	0.359	10.3	112	0.00
20		Pentachlorophenol	0.400	0.322	19.5	110	0.00
21		Phenanthrene	0.400	0.389	2.8	117	0.00
22		Anthracene	0.400	0.382	4.5	116	0.00
23		Fluoranthene	0.400	0.371	7.3	111	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	114	0.00
25		Pyrene	0.400	0.363	9.3	113	0.00
26	S	Terphenyl-d14	0.400	0.365	8.8	114	0.00
27		Benzo(a)anthracene	0.400	0.359	10.3	111	0.00
28		Chrysene	0.400	0.318	20.5	99	-0.02
29		Bis(2-ethylhexyl)phthalate	0.400	0.321	19.8	85	-0.02
30		Indeno(1,2,3-cd)pyrene	0.400	0.373	6.8	105	0.00
31	I	Perylene-d12	5.000	5.000	0.0	107	0.00
32		Benzo(b)fluoranthene	0.400	0.350	12.5	95	0.00
33		Benzo(k)fluoranthene	0.400	0.340	15.0	97	0.00
34	C	Benzo(a)pyrene	0.400	0.357	10.8	98	0.00
35		Dibenzo(a,h)anthracene	0.400	0.384	4.0	105	-0.02
36		Benzo(g,h,i)perylene	0.400	0.389	2.8	107	0.00

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