Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE103116\
Data File : BE092488.D
Acq On : 31 Oct 2016 20:59
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Nov 01 11:17:38 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE102516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 25 15:25:33 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	137	0.00
2		1,4-Dioxane	0.400	0.316	21.0	123	0.00
3		n-Nitrosodimethylamine	0.400	0.532	-33.0#	172	-0.04
4	S	2-Fluorophenol	0.400	0.620	-55.0#	250	-0.01
5	S	Phenol-d6	0.400	0.631	-57.7#	259	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.504	-26.0#	168	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	141	0.00
8	S	Nitrobenzene-d5	0.400	0.552	-38.0#	191	-0.02
9		Naphthalene	0.400	0.393	1.8	138	0.00
10		Hexachlorobutadiene	0.400	0.408	-2.0	142	0.00
11		2-Methylnaphthalene	0.400	0.393	1.8	144	-0.02
12	I	Acenaphthene-d10	5.000	5.000	0.0	130	0.00
13	S	2,4,6-Tribromophenol	0.400	0.581	-45.2#	223	-0.01
14	S	2-Fluorobiphenyl	0.400	0.424	-6.0	141	-0.01
15		Acenaphthylene	0.400	0.413	-3.2	139	0.00
16		Acenaphthene	0.400	0.404	-1.0	132	0.00
17		Fluorene	0.400	0.415	-3.7	140	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	117	0.00
19		Hexachlorobenzene	0.400	0.425	-6.2	130	0.00
20		Pentachlorophenol	0.400	0.927	-131.8#	382	-0.02
21		Phenanthrene	0.400	0.396	1.0	127	0.00
22		Anthracene	0.400	0.451	-12.7	150	-0.02
23		Fluoranthene	0.400	0.478	-19.5	149	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	122	0.00
25		Pyrene	0.400	0.454	-13.5	149	-0.02
26	S	Terphenyl-d14	0.400	0.455	-13.7	150	-0.02
27		Benzo(a)anthracene	0.400	0.471	-17.7	152	0.00
28		Chrysene	0.400	0.338	15.5	107	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.719	-79.7#	298	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.464	-16.0	153	0.00
31	I	Perylene-d12	5.000	5.000	0.0	126	0.01
32		Benzo(b)fluoranthene	0.400	0.396	1.0	144	0.00
33		Benzo(k)fluoranthene	0.400	0.441	-10.2	134	0.00
34	C	Benzo(a)pyrene	0.400	0.440	-10.0	157	0.00
35		Dibenzo(a,h)anthracene	0.400	0.447	-11.7	155	0.00
36		Benzo(g,h,i)perylene	0.400	0.428	-7.0	142	0.00

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