Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091216\
Data File : BE092378.D
Acq On : 12 Sep 2016 19:38
Operator : UM/SJ
Sample : SSTDCCC0.4EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Sep 13 05:32:55 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 02 19:34:51 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	150	-0.02
2		1,4-Dioxane	0.400	0.326	18.5	131	0.00
3		n-Nitrosodimethylamine	0.400	0.494	-23.5	228	-0.03
4	S	2-Fluorophenol	0.400	0.478	-19.5	211	-0.01
5	S	Phenol-d6	0.400	0.525	-31.3#	256	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.463	-15.8	222	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	157	0.00
8	S	Nitrobenzene-d5	0.400	0.493	-23.2	245	-0.02
9		Naphthalene	0.400	0.402	-0.5	165	-0.02
10		Hexachlorobutadiene	0.400	0.410	-2.5	163	-0.02
11		2-Methylnaphthalene	0.400	0.408	-2.0	173	-0.02
12	I	Acenaphthene-d10	5.000	5.000	0.0	148	0.00
13	S	2,4,6-Tribromophenol	0.400	0.532	-33.0#	245	-0.01
14	S	2-Fluorobiphenyl	0.400	0.392	2.0	161	-0.01
15		Acenaphthylene	0.400	0.417	-4.2	172	0.00
16		Acenaphthene	0.400	0.407	-1.7	164	-0.02
17		Fluorene	0.400	0.403	-0.8	165	-0.01
18	I	Phenanthrene-d10	5.000	5.000	0.0	134	0.00
19		Hexachlorobenzene	0.400	0.395	1.3	146	0.00
20		Pentachlorophenol	0.400	0.862	-115.5#	430	-0.02
21		Phenanthrene	0.400	0.459	-14.8	170	0.00
22		Anthracene	0.400	0.471	-17.7	191	0.00
23		Fluoranthene	0.400	0.621	-55.2#	236	-0.02
24	I	Chrysene-d12	5.000	5.000	0.0	123	-0.02
25		Pyrene	0.400	0.792	-98.0#	255	-0.02
26	S	Terphenyl-d14	0.400	0.533	-33.3#	174	0.00
27		Benzo(a)anthracene	0.400	0.920	-130.0#	288	-0.02
28		Chrysene	0.400	0.763	-90.8#	262	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	-5.000	1350.0#	1525	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.422	-5.5	132	-0.02
31	I	Perylene-d12	5.000	5.000	0.0	113	0.00
32		Benzo(b)fluoranthene	0.400	0.753	-88.3#	240	0.00
33		Benzo(k)fluoranthene	0.400	0.593	-48.2#	174	0.00
34	C	Benzo(a)pyrene	0.400	0.564	-41.0#	179	0.00
35		Dibenzo(a,h)anthracene	0.400	0.420	-5.0	134	0.00
36		Benzo(g,h,i)perylene	0.400	0.412	-3.0	125	-0.02

(#) = Out of Range SPCC's out = 0 CCC's out = 1