Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101716\
Data File : BE092455.D
Acq On : 17 Oct 2016 14:37
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Oct 17 15:39:06 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE092216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 22 18:25:56 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	163	-0.02
2		1,4-Dioxane	0.400	0.368	8.0	161	-0.01
3		n-Nitrosodimethylamine	0.400	0.513	-28.2#	204	-0.02
4	S	2-Fluorophenol	0.400	0.415	-3.7	158	-0.01
5	S	Phenol-d6	0.400	0.369	7.8	163	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.430	-7.5	196	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	153	0.00
8	S	Nitrobenzene-d5	0.400	0.470	-17.5	183	-0.02
9		Naphthalene	0.400	0.435	-8.7	167	-0.02
10		Hexachlorobutadiene	0.400	0.424	-6.0	164	-0.02
11		2-Methylnaphthalene	0.400	0.420	-5.0	163	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	142	0.00
13	S	2,4,6-Tribromophenol	0.400	0.396	1.0	123	0.00
14	S	2-Fluorobiphenyl	0.400	0.457	-14.2	165	0.00
15		Acenaphthylene	0.400	0.410	-2.5	149	0.00
16		Acenaphthene	0.400	0.430	-7.5	148	0.00
17		Fluorene	0.400	0.410	-2.5	148	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	124	0.02
19		Hexachlorobenzene	0.400	0.454	-13.5	151	0.00
20		Pentachlorophenol	0.400	0.392	2.0	127	0.00
21		Phenanthrene	0.400	0.468	-17.0	153	0.00
22		Anthracene	0.400	0.421	-5.2	144	0.02
23		Fluoranthene	0.400	0.415	-3.7	140	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	126	0.00
25		Pyrene	0.400	0.467	-16.8	153	0.00
26	S	Terphenyl-d14	0.400	0.444	-11.0	142	0.02
27		Benzo(a)anthracene	0.400	0.416	-4.0	134	0.00
28		Chrysene	0.400	0.446	-11.5	131	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.384	4.0	120	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.374	6.5	120	0.04
31	I	Perylene-d12	5.000	5.000	0.0	117	0.02
32		Benzo(b)fluoranthene	0.400	0.420	-5.0	123	0.02
33		Benzo(k)fluoranthene	0.400	0.436	-9.0	137	0.02
34	C	Benzo(a)pyrene	0.400	0.418	-4.5	129	0.02
35		Dibenzo(a,h)anthracene	0.400	0.395	1.3	121	0.05
36		Benzo(g,h,i)perylene	0.400	0.398	0.5	120	0.05

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