Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE022417\
Data File : BE092760.D
Acq On : 24 Feb 2017 15:04
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4EC

Quant Time: Feb 24 15:47:53 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE021517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 15 17:05:20 2017
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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	113	-0.02
2		1,4-Dioxane	0.710	0.775	-9.2	125	-0.01
3		n-Nitrosodimethylamine	0.740	0.862	-16.5	151#	-0.05
4	S	2-Fluorophenol	0.973	1.226	-26.0#	149	-0.03
5	S	Phenol-d6	1.194	1.548	-29.6#	152#	-0.05
6		bis(2-Chloroethyl)ether	1.321	1.550	-17.3	144	-0.02
7	I	Naphthalene-d8	1.000	1.000	0.0	104	-0.02
8	S	Nitrobenzene-d5	0.279	0.333	-19.4	129	-0.05
9		Naphthalene	1.082	1.182	-9.2	113	0.00
10		Hexachlorobutadiene	0.159	0.161	-1.3	101	0.00
11		2-Methylnaphthalene	0.680	0.686	-0.9	108	-0.02
12	I	Acenaphthene-d10	1.000	1.000	0.0	118	-0.02
13	S	2,4,6-Tribromophenol	0.106	0.163	-53.8#	186#	-0.03
14	S	2-Fluorobiphenyl	1.219	1.179	3.3	113	-0.02
15		Acenaphthylene	7.544	7.588	-0.6	120	-0.02
16		Acenaphthene	1.096	1.093	0.3	117	-0.02
17		Fluorene	1.407	1.548	-10.0	131	-0.02
18	I	Phenanthrene-d10	1.000	1.000	0.0	126	0.00
19		Hexachlorobenzene	0.193	0.191	1.0	114	-0.02
20		Pentachlorophenol	0.042	0.084	-100.0#	295#	-0.11
21		Phenanthrene	1.188	1.211	-1.9	122	-0.02
22		Anthracene	1.094	1.174	-7.3	139	0.00
23		Fluoranthene	5.096	6.418	-25.9#	159#	-0.02
24	I	Chrysene-d12	1.000	1.000	0.0	152#	-0.02
25		Pyrene	4.761	4.762	-0.0	157#	-0.02
26	S	Terphenyl-d14	0.638	0.641	-0.5	157#	-0.02
27		Benzo(a)anthracene	4.849	5.404	-11.4	173#	-0.02
28		Chrysene	5.752	4.728	17.8	125	0.00
29		Bis(2-ethylhexyl)phthalate	0.674	0.604	10.4	158#	0.00
30		Indeno(1,2,3-cd)pyrene	6.486	5.184	20.1	121	-0.03
31	I	Perylene-d12	1.000	1.000	0.0	138	-0.02
32		Benzo(b)fluoranthene	1.414	1.298	8.2	131	-0.02
33		Benzo(k)fluoranthene	1.495	1.397	6.6	142	-0.02
34	C	Benzo(a)pyrene	1.285	1.230	4.3	141	-0.02
35		Dibenzo(a,h)anthracene	1.355	1.177	13.1	122	-0.03
36		Benzo(g,h,i)perylene	5.728	4.917	14.2	120	-0.03

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