Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE111716\
Data File : BE092557.D
Acq On : 17 Nov 2016 9:16
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Nov 18 13:13:45 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE111616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 16 15:43:18 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	106	-0.02
2		1,4-Dioxane	0.681	0.805	-18.2	130	0.00
3		n-Nitrosodimethylamine	0.654	0.693	-6.0	131	-0.01
4	S	2-Fluorophenol	0.844	0.854	-1.2	107	0.00
5	S	Phenol-d6	1.104	1.205	-9.1	125	0.00
6		bis(2-Chloroethyl)ether	1.341	1.199	10.6	106	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
8	S	Nitrobenzene-d5	0.269	0.270	-0.4	108	-0.02
9		Naphthalene	1.032	1.055	-2.2	108	-0.02
10		Hexachlorobutadiene	0.157	0.159	-1.3	106	0.00
11		2-Methylnaphthalene	0.658	0.655	0.5	111	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	114	0.00
13	S	2,4,6-Tribromophenol	0.112	0.113	-0.9	122	-0.01
14	S	2-Fluorobiphenyl	1.219	1.175	3.6	109	0.00
15		Acenaphthylene	6.962	6.686	4.0	114	0.00
16		Acenaphthene	1.136	1.121	1.3	113	0.00
17		Fluorene	1.408	1.411	-0.2	115	-0.01
18	I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
19		Hexachlorobenzene	0.197	0.205	-4.1	123	0.00
20		Pentachlorophenol	0.036	0.039	-8.3	147	0.00
21		Phenanthrene	1.119	1.150	-2.8	113	-0.02
22		Anthracene	1.053	1.034	1.8	111	0.00
23		Fluoranthene	5.219	5.346	-2.4	116	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	114	0.00
25		Pyrene	3.975	3.839	3.4	118	0.00
26	S	Terphenyl-d14	0.572	0.537	6.1	116	-0.02
27		Benzo(a)anthracene	4.692	4.359	7.1	126	0.00
28		Chrysene	4.250	4.427	-4.2	114	-0.02
29		Bis(2-ethylhexyl)phthalate	0.305	0.342	-12.1	153#	-0.02
30		Indeno(1,2,3-cd)pyrene	4.722	4.307	8.8	109	0.00
31	I	Perylene-d12	1.000	1.000	0.0	106	-0.01
32		Benzo(b)fluoranthene	1.213	1.158	4.5	106	-0.01
33		Benzo(k)fluoranthene	1.270	1.270	0.0	113	-0.01
34	C	Benzo(a)pyrene	1.122	1.094	2.5	112	-0.01
35		Dibenzo(a,h)anthracene	1.111	1.061	4.5	109	-0.02
36		Benzo(g,h,i)perylene	4.702	4.553	3.2	107	-0.02

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