Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE082916\
Data File : BE092300.D
Acq On : 29 Aug 2016 21:57
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Aug 30 01:10:02 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE082916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 29 18:19:23 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	107	0.00
2		1,4-Dioxane	0.711	0.696	2.1	101	0.00
3		n-Nitrosodimethylamine	0.740	0.665	10.1	110	-0.01
4	S	2-Fluorophenol	1.122	1.018	9.3	116	0.00
5	S	Phenol-d6	1.478	1.269	14.1	116	-0.02
6		bis(2-Chloroethyl)ether	1.151	1.041	9.6	115	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	114	0.00
8	S	Nitrobenzene-d5	0.327	0.268	18.0	125	0.00
9		Naphthalene	1.115	1.031	7.5	106	0.00
10		Hexachlorobutadiene	0.169	0.160	5.3	107	0.00
11		2-Methylnaphthalene	0.689	0.641	7.0	111	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	121	0.00
13	S	2,4,6-Tribromophenol	0.135	0.128	5.2	126	0.00
14	S	2-Fluorobiphenyl	1.570	1.412	10.1	112	0.00
15		Acenaphthylene	8.560	7.742	9.6	116	0.00
16		Acenaphthene	1.397	1.299	7.0	115	0.00
17		Fluorene	1.689	1.541	8.8	115	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	116	0.00
19		Hexachlorobenzene	0.206	0.195	5.3	117	0.00
20		Pentachlorophenol	0.049	0.040	18.4	132	0.00
21		Phenanthrene	1.177	1.053	10.5	109	0.00
22		Anthracene	1.080	0.971	10.1	112	0.00
23		Fluoranthene	5.268	4.922	6.6	118	-0.02
24	I	Chrysene-d12	1.000	1.000	0.0	123	0.00
25		Pyrene	4.627	4.304	7.0	121	-0.02
26	S	Terphenyl-d14	0.676	0.614	9.2	114	0.00
27		Benzo(a)anthracene	5.171	4.942	4.4	136	0.00
28		Chrysene	4.699	4.627	1.5	113	-0.02
29		Bis(2-ethylhexyl)phthalate	0.328	0.379	-15.5	155#	0.00
30		Indeno(1,2,3-cd)pyrene	4.773	4.509	5.5	126	-0.02
31	I	Perylene-d12	1.000	1.000	0.0	119	-0.01
32		Benzo(b)fluoranthene	1.185	1.103	6.9	111	-0.01
33		Benzo(k)fluoranthene	1.343	1.284	4.4	121	-0.01
34	C	Benzo(a)pyrene	1.186	1.123	5.3	122	0.00
35		Dibenzo(a,h)anthracene	1.040	0.987	5.1	129	-0.03
36		Benzo(g,h,i)perylene	4.652	4.367	6.1	120	-0.02

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