Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE112316\
Data File : BE092572.D
Acq On : 23 Nov 2016 9:43
Operator : UM/SJ
Sample : SSTDICCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Nov 23 16:01:19 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	104	-0.02
2		1,4-Dioxane	0.681	0.722	-6.0	114	0.00
3		n-Nitrosodimethylamine	0.654	0.557	14.8	103	0.00
4	S	2-Fluorophenol	0.844	0.873	-3.4	108	0.00
5	S	Phenol-d6	1.104	1.084	1.8	111	0.00
6		bis(2-Chloroethyl)ether	1.341	1.120	16.5	98	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
8	S	Nitrobenzene-d5	0.269	0.256	4.8	100	0.00
9		Naphthalene	1.032	1.015	1.6	102	-0.02
10		Hexachlorobutadiene	0.157	0.158	-0.6	102	0.00
11		2-Methylnaphthalene	0.658	0.632	4.0	104	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
13	S	2,4,6-Tribromophenol	0.112	0.108	3.6	110	0.00
14	S	2-Fluorobiphenyl	1.219	1.169	4.1	102	0.00
15		Acenaphthylene	6.962	6.564	5.7	105	0.00
16		Acenaphthene	1.136	1.107	2.6	104	0.00
17		Fluorene	1.408	1.369	2.8	105	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
19		Hexachlorobenzene	0.197	0.187	5.1	107	0.00
20		Pentachlorophenol	0.036	0.035	2.8	125	0.00
21		Phenanthrene	1.119	1.113	0.5	104	-0.02
22		Anthracene	1.053	1.024	2.8	104	0.00
23		Fluoranthene	5.219	5.184	0.7	107	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	103	0.00
25		Pyrene	3.975	3.730	6.2	104	0.00
26	S	Terphenyl-d14	0.572	0.537	6.1	105	0.00
27		Benzo(a)anthracene	4.692	4.217	10.1	110	0.00
28		Chrysene	4.250	4.480	-5.4	105	-0.02
29		Bis(2-ethylhexyl)phthalate	0.305	0.290	4.9	117	-0.02
30		Indeno(1,2,3-cd)pyrene	4.722	4.139	12.3	95	0.00
31	I	Perylene-d12	1.000	1.000	0.0	99	-0.01
32		Benzo(b)fluoranthene	1.213	1.205	0.7	103	0.00
33		Benzo(k)fluoranthene	1.270	1.141	10.2	95	0.00
34	C	Benzo(a)pyrene	1.122	1.033	7.9	99	0.00
35		Dibenzo(a,h)anthracene	1.111	0.994	10.5	95	-0.02
36		Benzo(g,h,i)perylene	4.702	4.272	9.1	94	0.00

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