Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\be032116\
Data File : BE091495.D
Acq On : 21 Mar 2016 9:51
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 21 10:24:01 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 18 11:54:28 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	175	-0.02
2		1,4-Dioxane	0.400	0.349	12.8	160	0.00
3		n-Nitrosodimethylamine	0.400	0.375	6.3	189	-0.02
4	S	2-Fluorophenol	0.400	0.338	15.5	165	0.00
5	S	Phenol-d6	0.400	0.347	13.3	174	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	171	-0.02
7	S	Nitrobenzene-d5	0.400	0.493	-23.2	177	-0.01
8		Naphthalene	0.400	0.386	3.5	177	0.00
9		2-Methylnaphthalene	0.400	0.362	9.5	169	-0.01
10	I	Acenaphthene-d10	5.000	5.000	0.0	163	-0.01
11	S	2,4,6-Tribromophenol	0.400	0.536	-34.0#	194	-0.01
12	S	2-Fluorobiphenyl	0.400	0.377	5.8	167	-0.01
13		Acenaphthylene	0.400	0.434	-8.5	182	-0.01
14		Acenaphthene	0.400	0.376	6.0	160	0.00
15		Fluorene	0.400	0.368	8.0	161	-0.01
16	I	Phenanthrene-d10	5.000	5.000	0.0	157	-0.01
17		Hexachlorobenzene	0.400	0.363	9.3	158	0.00
18		Pentachlorophenol	0.400	0.508	-27.0#	203	-0.02
19		Phenanthrene	0.400	0.378	5.5	160	-0.01
20		Anthracene	0.400	0.359	10.3	162	-0.01
21		Fluoranthene	0.400	0.328	18.0	144	-0.02
22	I	Chrysene-d12	5.000	5.000	0.0	122	0.00
23		Pyrene	0.400	0.471	-17.7	151	-0.02
24	S	Terphenyl-d14	0.400	0.488	-22.0	151	-0.02
25		Benzo(a)anthracene	0.400	0.426	-6.5	139	0.00
26		Chrysene	0.400	0.508	-27.0#	175	-0.02
27		Indeno(1,2,3-cd)pyrene	0.400	0.467	-16.8	156	-0.05
28	I	Perylene-d12	5.000	5.000	0.0	157	-0.02
29		Benzo(b)fluoranthene	0.400	0.332	17.0	145	-0.02
30		Benzo(k)fluoranthene	0.400	0.347	13.3	148	-0.03
31	C	Benzo(a)pyrene	0.400	0.342	14.5	153	-0.02
32		Dibenzo(a,h)anthracene	0.400	0.354	11.5	157	-0.04
33		Benzo(g,h,i)perylene	0.400	0.346	13.5	150	-0.04

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