Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE050316\
Data File : BE091741.D
Acq On : 3 May 2016 12:59
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 04 01:08:44 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE042616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 26 16:09:46 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	122	0.00
2		1,4-Dioxane	0.400	0.318	20.5	96	0.02
3		n-Nitrosodimethylamine	0.400	0.329	17.8	102	0.02
4	S	2-Fluorophenol	0.400	0.343	14.2	108	0.00
5	S	Phenol-d6	0.400	0.328	18.0	106	0.02
6		bis(2-Chloroethyl)ether	0.400	0.374	6.5	113	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	119	0.00
8	S	Nitrobenzene-d5	0.400	0.321	19.8	102	0.01
9		Nitrobenzene	0.400	0.313	21.8	101	0.00
10		Naphthalene	0.400	0.403	-0.8	119	0.00
11		Hexachlorobutadiene	0.400	0.404	-1.0	123	-0.02
12		2-Methylnaphthalene	0.400	0.385	3.8	115	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	115	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.399	0.3	91	0.00
15	S	2-Fluorobiphenyl	0.400	0.396	1.0	116	0.00
16		Acenaphthylene	0.400	0.381	4.8	125	0.00
17		Acenaphthene	0.400	0.383	4.3	111	0.00
18		Fluorene	0.400	0.390	2.5	112	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	115	0.00
20		4-Bromophenyl-phenylether	0.400	0.387	3.3	110	0.00
21		Hexachlorobenzene	0.400	0.405	-1.3	116	0.00
22		Pentachlorophenol	0.400	0.362	9.5	104	0.00
23		Phenanthrene	0.400	0.406	-1.5	114	0.00
24		Anthracene	0.400	0.380	5.0	111	0.00
25		Fluoranthene	0.400	0.385	3.8	109	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	103	0.00
27		Benzidine	0.400	0.462	-15.5	136	0.00
28		Pyrene	0.400	0.423	-5.7	112	0.00
29	S	Terphenyl-d14	0.400	0.418	-4.5	110	0.00
30		Benzo(a)anthracene	0.400	0.384	4.0	101	0.00
31		3,3'-Dichlorobenzidine	0.400	0.347	13.3	103	-0.02
32		Chrysene	0.400	0.476	-19.0	116	0.00
33		Bis(2-ethylhexyl)phthalate	0.400	0.338	15.5	113	-0.02
34		Indeno(1,2,3-cd)pyrene	0.400	0.451	-12.7	114	-0.01
35	I	Perylene-d12	5.000	5.000	0.0	116	0.00
36		Benzo(b)fluoranthene	0.400	0.413	-3.2	123	-0.01
37		Benzo(k)fluoranthene	0.400	0.346	13.5	104	0.00
38	C	Benzo(a)pyrene	0.400	0.373	6.8	114	0.00
39		Dibenzo(a,h)anthracene	0.400	0.377	5.8	113	-0.02
40		Benzo(g,h,i)perylene	0.400	0.373	6.8	112	-0.01

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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