Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE041216\
Data File : BE091645.D
Acq On : 12 Apr 2016 10:29
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Apr 13 07:09:51 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE040816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 08 16:48:14 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	125	-0.02
2		1,4-Dioxane	0.400	0.330	17.5	108	0.00
3		n-Nitrosodimethylamine	0.400	0.372	7.0	125	0.00
4	S	2-Fluorophenol	0.400	0.372	7.0	123	0.00
5	S	Phenol-d6	0.400	0.374	6.5	128	0.00
6		bis(2-Chloroethyl)ether	0.400	0.387	3.3	128	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	128	-0.02
8	S	Nitrobenzene-d5	0.400	0.362	9.5	130	0.00
9		Nitrobenzene	0.400	0.458	-14.5	130	0.00
10		Naphthalene	0.400	0.395	1.3	128	0.00
11		Hexachlorobutadiene	0.400	0.380	5.0	123	0.00
12		2-Methylnaphthalene	0.400	0.386	3.5	128	-0.01
13	I	Acenaphthene-d10	5.000	5.000	0.0	126	0.00
14	S	2,4,6-Tribromophenol	0.400	0.499	-24.7	138	0.00
15	S	2-Fluorobiphenyl	0.400	0.392	2.0	126	0.00
16		Acenaphthylene	0.400	0.373	6.8	145	0.00
17		Acenaphthene	0.400	0.406	-1.5	133	-0.01
18		Fluorene	0.400	0.384	4.0	125	-0.01
19	I	Phenanthrene-d10	5.000	5.000	0.0	131	0.00
20		4-Bromophenyl-phenylether	0.400	0.351	12.3	119	0.00
21		Hexachlorobenzene	0.400	0.371	7.3	122	0.00
22		Pentachlorophenol	0.400	0.455	-13.7	158	0.00
23		Phenanthrene	0.400	0.374	6.5	123	0.00
24		Anthracene	0.400	0.358	10.5	124	0.00
25		Fluoranthene	0.400	0.350	12.5	124	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	131	0.00
27		Benzidine	0.400	0.545	-36.3#	163	-0.02
28		Pyrene	0.400	0.358	10.5	128	-0.02
29	S	Terphenyl-d14	0.400	0.400	0.0	134	0.00
30		Benzo(a)anthracene	0.400	0.414	-3.5	142	0.00
31		3,3'-Dichlorobenzidine	0.400	0.407	-1.7	138	0.00
32		Chrysene	0.400	0.406	-1.5	136	-0.02
33		Bis(2-ethylhexyl)phthalate	0.400	0.407	-1.7	129	0.00
34		Indeno(1,2,3-cd)pyrene	0.400	0.374	6.5	128	-0.01
35	I	Perylene-d12	5.000	5.000	0.0	130	0.00
36		Benzo(b)fluoranthene	0.400	0.411	-2.7	141	-0.01
37		Benzo(k)fluoranthene	0.400	0.371	7.3	127	0.00
38	C	Benzo(a)pyrene	0.400	0.379	5.3	131	0.00
39		Dibenzo(a,h)anthracene	0.400	0.376	6.0	129	-0.01
40		Benzo(g,h,i)perylene	0.400	0.375	6.3	126	-0.01

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

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