Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042016\
Data File : BE091694.D
Acq On : 20 Apr 2016 20:47
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Apr 21 04:24:29 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE042016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 20 18:55:38 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	98	0.00
2		1,4-Dioxane	0.400	0.398	0.5	95	0.00
3		n-Nitrosodimethylamine	0.400	0.394	1.5	101	0.00
4	S	2-Fluorophenol	0.400	0.382	4.5	97	0.00
5	S	Phenol-d6	0.400	0.369	7.8	96	0.00
6		bis(2-Chloroethyl)ether	0.400	0.379	5.3	95	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	100	0.00
8	S	Nitrobenzene-d5	0.400	0.365	8.8	102	0.00
9		Nitrobenzene	0.400	0.369	7.8	102	-0.01
10		Naphthalene	0.400	0.393	1.8	97	0.00
11		Hexachlorobutadiene	0.400	0.404	-1.0	100	0.00
12		2-Methylnaphthalene	0.400	0.388	3.0	98	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	98	0.00
14	S	2,4,6-Tribromophenol	0.400	0.429	-7.2	106	-0.01
15	S	2-Fluorobiphenyl	0.400	0.403	-0.8	99	0.00
16		Acenaphthylene	0.400	0.370	7.5	100	0.00
17		Acenaphthene	0.400	0.393	1.8	98	0.00
18		Fluorene	0.400	0.395	1.3	99	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	100	0.00
20		4-Bromophenyl-phenylether	0.400	0.395	1.3	100	0.00
21		Hexachlorobenzene	0.400	0.400	0.0	98	0.00
22		Pentachlorophenol	0.400	0.388	3.0	96	-0.02
23		Phenanthrene	0.400	0.414	-3.5	99	-0.01
24		Anthracene	0.400	0.382	4.5	100	-0.01
25		Fluoranthene	0.400	0.367	8.3	95	-0.02
26	I	Chrysene-d12	5.000	5.000	0.0	94	0.00
27		Benzidine	0.400	0.432	-8.0	99	0.00
28		Pyrene	0.400	0.395	1.3	99	0.00
29	S	Terphenyl-d14	0.400	0.388	3.0	97	0.00
30		Benzo(a)anthracene	0.400	0.370	7.5	94	0.00
31		3,3'-Dichlorobenzidine	0.400	0.431	-7.7	103	-0.02
32		Chrysene	0.400	0.432	-8.0	104	0.00
33		Bis(2-ethylhexyl)phthalate	0.400	0.433	-8.2	113	-0.02
34		Indeno(1,2,3-cd)pyrene	0.400	0.366	8.5	94	-0.01
35	I	Perylene-d12	5.000	5.000	0.0	95	-0.01
36		Benzo(b)fluoranthene	0.400	0.376	6.0	94	0.00
37		Benzo(k)fluoranthene	0.400	0.375	6.3	94	-0.01
38	C	Benzo(a)pyrene	0.400	0.360	10.0	93	-0.01
39		Dibenzo(a,h)anthracene	0.400	0.359	10.3	93	0.00
40		Benzo(g,h,i)perylene	0.400	0.372	7.0	94	-0.01

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

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