Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\Data\BE051616\
Data File : BE091773.D
Acq On : 16 May 2016 14:41
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 17 04:20:02 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	125	0.00
2		1,4-Dioxane	0.400	0.301	24.8	95	0.02
3		n-Nitrosodimethylamine	0.400	0.289	27.8#	92	0.02
4	S	2-Fluorophenol	0.400	0.342	14.5	110	0.00
5	S	Phenol-d6	0.400	0.339	15.3	113	0.00
6		bis(2-Chloroethyl)ether	0.400	0.362	9.5	112	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	135	0.00
8	S	Nitrobenzene-d5	0.400	0.300	25.0#	109	0.00
9		Nitrobenzene	0.400	0.290	27.5#	106	0.00
10		Naphthalene	0.400	0.390	2.5	130	0.00
11		Hexachlorobutadiene	0.400	0.383	4.3	132	-0.02
12		2-Methylnaphthalene	0.400	0.378	5.5	128	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	139	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.408	-2.0	113	0.00
15	S	2-Fluorobiphenyl	0.400	0.371	7.3	130	0.00
16		Acenaphthylene	0.400	0.366	8.5	144	0.00
17		Acenaphthene	0.400	0.369	7.8	128	0.00
18		Fluorene	0.400	0.363	9.3	126	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	130	0.00
20		4-Bromophenyl-phenylether	0.400	0.364	9.0	117	0.00
21		Hexachlorobenzene	0.400	0.401	-0.3	130	0.00
22		Pentachlorophenol	0.400	0.391	2.3	130	0.00
23		Phenanthrene	0.400	0.386	3.5	122	0.00
24		Anthracene	0.400	0.363	9.3	120	0.00
25		Fluoranthene	0.400	0.349	12.8	112	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	99	0.00
27		Benzidine	0.400	0.450	-12.5	126	0.00
28		Pyrene	0.400	0.437	-9.2	112	0.00
29	S	Terphenyl-d14	0.400	0.422	-5.5	107	0.00
30		Benzo(a)anthracene	0.400	0.372	7.0	94	0.00
31		3,3'-Dichlorobenzidine	0.400	0.334	16.5	95	-0.02
32		Chrysene	0.400	0.437	-9.2	102	0.00
33		Bis(2-ethylhexyl)phthalate	0.400	0.389	2.8	125	-0.02
34		Indeno(1,2,3-cd)pyrene	0.400	0.421	-5.2	102	-0.01
35	I	Perylene-d12	5.000	5.000	0.0	112	0.00
36		Benzo(b)fluoranthene	0.400	0.376	6.0	108	0.00
37		Benzo(k)fluoranthene	0.400	0.343	14.2	100	0.00
38	C	Benzo(a)pyrene	0.400	0.341	14.8	101	0.00
39		Dibenzo(a,h)anthracene	0.400	0.347	13.3	100	-0.01
40		Benzo(g,h,i)perylene	0.400	0.353	11.8	102	-0.02

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

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