Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE052316\
Data File : BE091842.D
Acq On : 23 May 2016 11:38
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 24 02:31:06 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE051916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 19 15:36:49 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	103	-0.02
2		1,4-Dioxane	0.400	0.526	-31.5#	113	-0.02
3		n-Nitrosodimethylamine	0.400	0.413	-3.2	116	0.00
4	S	2-Fluorophenol	0.400	0.462	-15.5	126	0.00
5	S	Phenol-d6	0.400	0.389	2.8	111	0.00
6		bis(2-Chloroethyl)ether	0.400	0.403	-0.8	113	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	92	-0.02
8	S	Nitrobenzene-d5	0.400	0.434	-8.5	112	-0.01
9		Nitrobenzene	0.400	0.418	-4.5	111	-0.01
10		Naphthalene	0.400	0.374	6.5	90	0.00
11		Hexachlorobutadiene	0.400	0.391	2.3	94	-0.02
12		2-Methylnaphthalene	0.400	0.373	6.8	91	-0.01
13	I	Acenaphthene-d10	5.000	5.000	0.0	86	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.565	-41.2#	140	-0.01
15	S	2-Fluorobiphenyl	0.400	0.398	0.5	90	-0.01
16		Acenaphthylene	0.400	0.349	12.8	79	-0.01
17		Acenaphthene	0.400	0.363	9.3	83	-0.01
18		Fluorene	0.400	0.389	2.8	89	-0.01
19	I	Phenanthrene-d10	5.000	5.000	0.0	99	-0.01
20		4-Bromophenyl-phenylether	0.400	0.359	10.3	91	0.00
21		Hexachlorobenzene	0.400	0.344	14.0	86	0.00
22		Pentachlorophenol	0.400	0.275	31.3#	76	0.00
23		Phenanthrene	0.400	0.335	16.3	87	-0.01
24		Anthracene	0.400	0.364	9.0	93	-0.01
25		Fluoranthene	0.400	0.403	-0.8	100	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	108	-0.02
27		Benzidine	0.400	0.525	-31.3#	160	-0.02
28		Pyrene	0.400	0.353	11.8	100	-0.02
29	S	Terphenyl-d14	0.400	0.391	2.3	114	-0.02
30		Benzo(a)anthracene	0.400	0.418	-4.5	121	-0.02
31		3,3'-Dichlorobenzidine	0.400	0.403	-0.8	125	-0.02
32		Chrysene	0.400	0.427	-6.7	106	-0.02
33		Bis(2-ethylhexyl)phthalate	0.400	0.566	-41.5#	146	-0.02
34		Indeno(1,2,3-cd)pyrene	0.400	0.397	0.8	114	-0.03
35	I	Perylene-d12	5.000	5.000	0.0	109	-0.02
36		Benzo(b)fluoranthene	0.400	0.398	0.5	115	-0.02
37		Benzo(k)fluoranthene	0.400	0.390	2.5	112	-0.02
38	C	Benzo(a)pyrene	0.400	0.401	-0.3	116	-0.02
39		Dibenzo(a,h)anthracene	0.400	0.393	1.8	115	-0.05
40		Benzo(g,h,i)perylene	0.400	0.388	3.0	112	-0.05

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

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