Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE060716\
Data File : BE091886.D
Acq On : 7 Jun 2016 20:20
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 08 07:15:16 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 07 17:14:38 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	105	0.00
2		1,4-Dioxane	0.400	0.411	-2.7	99	0.00
3		n-Nitrosodimethylamine	0.400	0.397	0.8	100	0.00
4	S	2-Fluorophenol	0.400	0.416	-4.0	110	0.00
5	S	Phenol-d6	0.400	0.408	-2.0	112	0.00
6		bis(2-Chloroethyl)ether	0.400	0.395	1.3	104	0.00
7		n-Nitroso-di-n-propylamine	0.400	0.385	3.8	104	0.00
8	I	Naphthalene-d8	5.000	5.000	0.0	105	0.00
9	S	Nitrobenzene-d5	0.400	0.401	-0.3	104	-0.02
10		Nitrobenzene	0.400	0.402	-0.5	108	0.00
11		bis(2-Chloroethoxy)methane	0.400	0.398	0.5	111	0.00
12		Naphthalene	0.400	0.423	-5.7	105	0.00
13		Hexachlorobutadiene	0.400	0.424	-6.0	105	0.00
14		2-Methylnaphthalene	0.400	0.415	-3.7	107	0.00
15	I	Acenaphthene-d10	5.000	5.000	0.0	110	0.00
16	S	2,4,6-Tribromophenol	0.400	0.389	2.8	116	-0.02
17	S	2-Fluorobiphenyl	0.400	0.417	-4.2	106	0.00
18		Acenaphthylene	0.400	0.380	5.0	103	0.00
19		2,6-Dinitrotoluene	0.400	0.389	2.8	109	-0.02
20		Acenaphthene	0.400	0.407	-1.7	107	0.00
21		2,4-Dinitrotoluene	0.400	0.425	-6.2	113	-0.02
22		Fluorene	0.400	0.410	-2.5	110	-0.02
23		4-Chlorophenyl-phenylether	0.400	0.425	-6.2	110	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	95	0.00
25		4-Bromophenyl-phenylether	0.400	0.431	-7.7	109	0.00
26		Hexachlorobenzene	0.400	0.442	-10.5	107	0.00
27		Pentachlorophenol	0.400	0.452	-13.0	118	0.00
28		Phenanthrene	0.400	0.445	-11.2	108	-0.01
29		Anthracene	0.400	0.429	-7.2	112	0.00
30		Fluoranthene	0.400	0.431	-7.7	109	-0.02
31	I	Chrysene-d12	5.000	5.000	0.0	109	0.00
32		Benzidine	0.400	0.390	2.5	108	0.00
33		Pyrene	0.400	0.507	-26.7#	115	0.00
34	S	Terphenyl-d14	0.400	0.436	-9.0	110	0.00
35		Benzo(a)anthracene	0.400	0.350	12.5	103	0.00
36		3,3'-Dichlorobenzidine	0.400	0.440	-10.0	121	0.00
37		Chrysene	0.400	0.355	11.3	95	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.428	-7.0	132	0.00
39		Indeno(1,2,3-cd)pyrene	0.400	0.406	-1.5	103	0.00
40	I	Perylene-d12	5.000	5.000	0.0	101	0.00
41		Benzo(b)fluoranthene	0.400	0.419	-4.7	104	0.00

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Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.409	-2.2	104	0.00
43 C	Benzo(a)pyrene	0.400	0.417	-4.2	108	0.00
44	Dibenzo(a,h)anthracene	0.400	0.403	-0.8	104	0.00
45	Benzo(g,h,i)perylene	0.400	0.411	-2.7	102	0.00

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