Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE080116\
Data File : BE092104.D
Acq On : 1 Aug 2016 22:52
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Aug 02 01:28:54 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 01 17:22:55 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	0.00
2		1,4-Dioxane	0.400	0.332	17.0	100	0.00
3		n-Nitrosodimethylamine	0.400	0.351	12.3	96	0.02
4	S	2-Fluorophenol	0.400	0.406	-1.5	102	0.00
5	S	Phenol-d6	0.400	0.502	-25.5#	123	0.00
6		bis(2-Chloroethyl)ether	0.400	0.367	8.3	100	0.00
7		n-Nitroso-di-n-propylamine	0.400	0.374	6.5	103	0.00
8	I	Naphthalene-d8	5.000	5.000	0.0	119	0.00
9	S	Nitrobenzene-d5	0.400	0.368	8.0	106	0.00
10		Nitrobenzene	0.400	0.330	17.5	106	0.00
11		bis(2-Chloroethoxy)methane	0.400	0.343	14.2	103	0.00
12		Naphthalene	0.400	0.394	1.5	116	0.00
13		Hexachlorobutadiene	0.400	0.390	2.5	116	0.00
14		2-Methylnaphthalene	0.400	0.399	0.3	121	0.00
15	I	Acenaphthene-d10	5.000	5.000	0.0	136	0.00
16	S	2,4,6-Tribromophenol	0.400	0.315	21.3	97	0.00
17	S	2-Fluorobiphenyl	0.400	0.390	2.5	135	0.00
18		Acenaphthylene	0.400	0.336	16.0	118	0.00
19		2,6-Dinitrotoluene	0.400	0.276	31.0#	99	0.00
20		Acenaphthene	0.400	0.408	-2.0	138	0.00
21		2,4-Dinitrotoluene	0.400	0.260	35.0#	104	0.00
22		Fluorene	0.400	0.346	13.5	119	0.00
23		4-Chlorophenyl-phenylether	0.400	0.365	8.8	115	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	112	0.00
25		4-Bromophenyl-phenylether	0.400	0.394	1.5	117	0.00
26		Hexachlorobenzene	0.400	0.416	-4.0	119	0.00
27		Pentachlorophenol	0.400	0.349	12.8	70	0.00
28		Phenanthrene	0.400	0.269	32.8#	84	0.00
29		Anthracene	0.400	0.376	6.0	110	0.00
30		Fluoranthene	0.400	0.382	4.5	108	-0.03
31	I	Chrysene-d12	5.000	5.000	0.0	101	0.00
32		Benzidine	0.400	0.519	-29.8#	118	0.00
33		Pyrene	0.400	0.397	0.8	103	0.00
34	S	Terphenyl-d14	0.400	0.170	57.5#	72	0.00
35		Benzo(a)anthracene	0.400	0.402	-0.5	112	0.00
36		3,3'-Dichlorobenzidine	0.400	0.356	11.0	95	0.00
37		Chrysene	0.400	0.405	-1.3	96	-0.02
38		Bis(2-ethylhexyl)phthalate	0.400	0.246	38.5#	77	0.00
39		Indeno(1,2,3-cd)pyrene	0.400	0.368	8.0	96	0.00
40	I	Perylene-d12	5.000	5.000	0.0	95	0.00
41		Benzo(b)fluoranthene	0.400	0.391	2.3	100	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.385	3.8	88	0.00
43 C	Benzo(a)pyrene	0.400	0.388	3.0	95	0.00
44	Dibenzo(a,h)anthracene	0.400	0.386	3.5	95	-0.02
45	Benzo(g,h,i)perylene	0.400	0.397	0.8	96	-0.02

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