Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072216\
Data File : BE092025.D
Acq On : 23 Jul 2016 3:11
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 23 05:03:38 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE072216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 22 15:39:27 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	101	0.00
2		1,4-Dioxane	0.400	0.406	-1.5	97	0.00
3		n-Nitrosodimethylamine	0.400	0.420	-5.0	117	-0.01
4	S	2-Fluorophenol	0.400	0.475	-18.7	132	0.00
5	S	Phenol-d6	0.400	0.474	-18.5	130	0.00
6		bis(2-Chloroethyl)ether	0.400	0.409	-2.2	112	-0.02
7		n-Nitroso-di-n-propylamine	0.400	0.400	0.0	112	0.00
8	I	Naphthalene-d8	5.000	5.000	0.0	103	0.00
9	S	Nitrobenzene-d5	0.400	0.416	-4.0	112	-0.02
10		Nitrobenzene	0.400	0.452	-13.0	113	-0.02
11		bis(2-Chloroethoxy)methane	0.400	0.485	-21.2	135	-0.03
12		Naphthalene	0.400	0.407	-1.7	102	0.00
13		Hexachlorobutadiene	0.400	0.414	-3.5	102	0.00
14		2-Methylnaphthalene	0.400	0.412	-3.0	106	0.00
15	I	Acenaphthene-d10	5.000	5.000	0.0	95	0.00
16	S	2,4,6-Tribromophenol	0.400	0.482	-20.5	135	0.00
17	S	2-Fluorobiphenyl	0.400	0.425	-6.2	105	0.00
18		Acenaphthylene	0.400	0.430	-7.5	124	0.00
19		2,6-Dinitrotoluene	0.400	0.430	-7.5	146	0.00
20		Acenaphthene	0.400	0.404	-1.0	96	0.00
21		2,4-Dinitrotoluene	0.400	0.468	-17.0	127	0.00
22		Fluorene	0.400	0.423	-5.7	108	0.00
23		4-Chlorophenyl-phenylether	0.400	0.425	-6.2	105	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	107	0.00
25		4-Bromophenyl-phenylether	0.400	0.402	-0.5	108	0.00
26		Hexachlorobenzene	0.400	0.415	-3.7	108	0.00
27		Pentachlorophenol	0.400	0.517	-29.3#	153	-0.02
28		Phenanthrene	0.400	0.409	-2.2	107	-0.01
29		Anthracene	0.400	0.404	-1.0	113	-0.01
30		Fluoranthene	0.400	0.407	-1.7	114	-0.02
31	I	Chrysene-d12	5.000	5.000	0.0	82	0.00
32		Benzidine	0.400	0.565	-41.2#	158	0.00
33		Pyrene	0.400	0.534	-33.5#	116	0.00
34	S	Terphenyl-d14	0.400	0.472	-18.0	100	0.00
35		Benzo(a)anthracene	0.400	0.562	-40.5#	128	0.00
36		3,3'-Dichlorobenzidine	0.400	0.658	-64.5#	159	-0.03
37		Chrysene	0.400	0.513	-28.2#	105	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.835	-108.7#	230	0.00
39		Indeno(1,2,3-cd)pyrene	0.400	0.472	-18.0	98	-0.02
40	I	Perylene-d12	5.000	5.000	0.0	97	0.00
41		Benzo(b)fluoranthene	0.400	0.425	-6.2	108	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.388	3.0	93	0.00
43 C	Benzo(a)pyrene	0.400	0.395	1.3	100	0.00
44	Dibenzo(a,h)anthracene	0.400	0.399	0.3	99	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