Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072716\
Data File : BE092070.D
Acq On : 28 Jul 2016 6:59
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 28 07:33:10 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE072716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 27 17:13:26 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	114	-0.02
2		1,4-Dioxane	0.400	0.366	8.5	104	0.00
3		n-Nitrosodimethylamine	0.400	0.410	-2.5	129	0.00
4	S	2-Fluorophenol	0.400	0.492	-23.0	160	0.00
5	S	Phenol-d6	0.400	0.504	-26.0#	162	0.00
6		bis(2-Chloroethyl)ether	0.400	0.394	1.5	124	0.00
7		n-Nitroso-di-n-propylamine	0.400	0.451	-12.7	145	0.00
8	I	Naphthalene-d8	5.000	5.000	0.0	123	0.00
9	S	Nitrobenzene-d5	0.400	0.390	2.5	129	0.00
10		Nitrobenzene	0.400	0.418	-4.5	145	0.00
11		bis(2-Chloroethoxy)methane	0.400	0.397	0.8	140	-0.02
12		Naphthalene	0.400	0.416	-4.0	129	0.00
13		Hexachlorobutadiene	0.400	0.432	-8.0	136	0.00
14		2-Methylnaphthalene	0.400	0.435	-8.7	143	0.00
15	I	Acenaphthene-d10	5.000	5.000	0.0	136	0.00
16	S	2,4,6-Tribromophenol	0.400	0.597	-49.2#	232	0.00
17	S	2-Fluorobiphenyl	0.400	0.408	-2.0	141	0.00
18		Acenaphthylene	0.400	0.437	-9.2	161	0.00
19		2,6-Dinitrotoluene	0.400	0.618	-54.5#	233	0.00
20		Acenaphthene	0.400	0.405	-1.3	141	0.00
21		2,4-Dinitrotoluene	0.400	0.477	-19.2	211	0.00
22		Fluorene	0.400	0.439	-9.7	154	0.00
23		4-Chlorophenyl-phenylether	0.400	0.423	-5.7	145	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	133	0.00
25		4-Bromophenyl-phenylether	0.400	0.422	-5.5	144	0.00
26		Hexachlorobenzene	0.400	0.439	-9.7	149	0.00
27		Pentachlorophenol	0.400	0.715	-78.7#	303	0.00
28		Phenanthrene	0.400	0.437	-9.2	150	0.00
29		Anthracene	0.400	0.452	-13.0	162	0.00
30		Fluoranthene	0.400	0.467	-16.8	177	0.00
31	I	Chrysene-d12	5.000	5.000	0.0	114	0.00
32		Benzidine	0.400	0.947	-136.7#	324	0.00
33		Pyrene	0.400	0.516	-29.0#	161	0.00
34	S	Terphenyl-d14	0.400	0.491	-22.7	158	0.00
35		Benzo(a)anthracene	0.400	0.493	-23.2	138	0.00
36		3,3'-Dichlorobenzidine	0.400	0.728	-82.0#	299	0.00
37		Chrysene	0.400	0.325	18.8	114	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.760	-90.0#	237	0.00
39		Indeno(1,2,3-cd)pyrene	0.400	0.523	-30.8#	160	0.00
40	I	Perylene-d12	5.000	5.000	0.0	124	0.00
41		Benzo(b)fluoranthene	0.400	0.457	-14.2	143	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.455	-13.7	133	0.00
43 C	Benzo(a)pyrene	0.400	0.471	-17.7	151	0.00
44	Dibenzo(a,h)anthracene	0.400	0.485	-21.2	163	0.00
45	Benzo(g,h,i)perylene	0.400	0.477	-19.2	159	0.00

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