Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\Be072216\
Data File : BE092023.D
Acq On : 23 Jul 2016 1:53
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 23 03:02:53 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	106	0.00
2		1,4-Dioxane	0.400	0.403	-0.8	101	0.00
3		n-Nitrosodimethylamine	0.400	0.442	-10.5	130	-0.02
4	S	2-Fluorophenol	0.400	0.509	-27.2#	148	0.00
5	S	Phenol-d6	0.400	0.499	-24.7	146	0.00
6		bis(2-Chloroethyl)ether	0.400	0.419	-4.7	120	-0.02
7		n-Nitroso-di-n-propylamine	0.400	0.437	-9.2	127	0.00
8	I	Naphthalene-d8	5.000	5.000	0.0	111	0.00
9	S	Nitrobenzene-d5	0.400	0.427	-6.7	123	-0.02
10		Nitrobenzene	0.400	0.458	-14.5	124	-0.02
11		bis(2-Chloroethoxy)methane	0.400	0.507	-26.7#	152	-0.02
12		Naphthalene	0.400	0.402	-0.5	108	0.00
13		Hexachlorobutadiene	0.400	0.402	-0.5	107	0.00
14		2-Methylnaphthalene	0.400	0.410	-2.5	113	0.00
15	I	Acenaphthene-d10	5.000	5.000	0.0	103	0.00
16	S	2,4,6-Tribromophenol	0.400	0.511	-27.7#	161	0.00
17	S	2-Fluorobiphenyl	0.400	0.420	-5.0	112	0.00
18		Acenaphthylene	0.400	0.416	-4.0	130	0.00
19		2,6-Dinitrotoluene	0.400	0.435	-8.7	162	0.00
20		Acenaphthene	0.400	0.387	3.3	100	0.00
21		2,4-Dinitrotoluene	0.400	0.812	-103.0#	281	0.00
22		Fluorene	0.400	0.428	-7.0	119	0.00
23		4-Chlorophenyl-phenylether	0.400	0.426	-6.5	114	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	112	0.00
25		4-Bromophenyl-phenylether	0.400	0.408	-2.0	116	0.00
26		Hexachlorobenzene	0.400	0.404	-1.0	110	0.00
27		Pentachlorophenol	0.400	0.589	-47.2#	201	-0.02
28		Phenanthrene	0.400	0.411	-2.7	114	-0.01
29		Anthracene	0.400	0.411	-2.7	121	-0.01
30		Fluoranthene	0.400	0.416	-4.0	123	-0.02
31	I	Chrysene-d12	5.000	5.000	0.0	79	0.00
32		Benzidine	0.400	0.649	-62.2#	188	0.00
33		Pyrene	0.400	0.586	-46.5#	124	0.00
34	S	Terphenyl-d14	0.400	0.544	-36.0#	112	-0.02
35		Benzo(a)anthracene	0.400	0.662	-65.5#	147	0.00
36		3,3'-Dichlorobenzidine	0.400	0.850	-112.5#	211	-0.03
37		Chrysene	0.400	0.513	-28.2#	102	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	1.306	-226.5#	363	0.00
39		Indeno(1,2,3-cd)pyrene	0.400	0.519	-29.8#	105	-0.02
40	I	Perylene-d12	5.000	5.000	0.0	103	0.00
41		Benzo(b)fluoranthene	0.400	0.439	-9.7	119	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.383	4.3	99	0.00
43 C	Benzo(a)pyrene	0.400	0.407	-1.7	110	0.00
44	Dibenzo(a,h)anthracene	0.400	0.406	-1.5	107	0.00
45	Benzo(g,h,i)perylene	0.400	0.410	-2.5	109	0.00

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