Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\Be072216\
Data File : BE092036.D
Acq On : 23 Jul 2016 10:11
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 25 02:16:44 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	122	0.00
2		1,4-Dioxane	0.400	0.368	8.0	105	0.00
3		n-Nitrosodimethylamine	0.400	0.404	-1.0	133	-0.02
4	S	2-Fluorophenol	0.400	0.512	-28.0#	170	0.00
5	S	Phenol-d6	0.400	0.526	-31.5#	179	0.00
6		bis(2-Chloroethyl)ether	0.400	0.431	-7.7	141	-0.02
7		n-Nitroso-di-n-propylamine	0.400	0.450	-12.5	150	0.00
8	I	Naphthalene-d8	5.000	5.000	0.0	134	0.00
9	S	Nitrobenzene-d5	0.400	0.438	-9.5	153	-0.02
10		Nitrobenzene	0.400	0.469	-17.2	155	-0.02
11		bis(2-Chloroethoxy)methane	0.400	0.403	-0.8	146	-0.03
12		Naphthalene	0.400	0.406	-1.5	132	0.00
13		Hexachlorobutadiene	0.400	0.398	0.5	128	0.00
14		2-Methylnaphthalene	0.400	0.413	-3.2	138	0.00
15	I	Acenaphthene-d10	5.000	5.000	0.0	116	0.00
16	S	2,4,6-Tribromophenol	0.400	0.518	-29.5#	184	0.00
17	S	2-Fluorobiphenyl	0.400	0.445	-11.2	133	-0.01
18		Acenaphthylene	0.400	0.491	-22.7	173	0.00
19		2,6-Dinitrotoluene	0.400	0.454	-13.5	195	0.00
20		Acenaphthene	0.400	0.378	5.5	109	0.00
21		2,4-Dinitrotoluene	0.400	0.766	-91.5#	291	-0.03
22		Fluorene	0.400	0.452	-13.0	141	0.00
23		4-Chlorophenyl-phenylether	0.400	0.422	-5.5	127	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	123	0.00
25		4-Bromophenyl-phenylether	0.400	0.422	-5.5	130	0.00
26		Hexachlorobenzene	0.400	0.430	-7.5	128	0.00
27		Pentachlorophenol	0.400	0.561	-40.3#	202	-0.02
28		Phenanthrene	0.400	0.410	-2.5	124	-0.01
29		Anthracene	0.400	0.419	-4.7	134	-0.01
30		Fluoranthene	0.400	0.395	1.3	127	-0.02
31	I	Chrysene-d12	5.000	5.000	0.0	80	0.00
32		Benzidine	0.400	0.772	-93.0#	238	0.00
33		Pyrene	0.400	0.615	-53.7#	131	0.00
34	S	Terphenyl-d14	0.400	0.532	-33.0#	110	-0.02
35		Benzo(a)anthracene	0.400	0.651	-62.7#	145	0.00
36		3,3'-Dichlorobenzidine	0.400	0.980	-145.0#	253	-0.03
37		Chrysene	0.400	0.511	-27.7#	102	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	1.506	-276.5#	424	0.00
39		Indeno(1,2,3-cd)pyrene	0.400	0.569	-42.2#	115	-0.02
40	I	Perylene-d12	5.000	5.000	0.0	102	-0.01
41		Benzo(b)fluoranthene	0.400	0.464	-16.0	124	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.381	4.8	97	0.00
43 C	Benzo(a)pyrene	0.400	0.432	-8.0	115	0.00
44	Dibenzo(a,h)anthracene	0.400	0.452	-13.0	118	-0.02
45	Benzo(g,h,i)perylene	0.400	0.458	-14.5	120	-0.02

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