Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE060716\
Data File : BE091898.D
Acq On : 8 Jun 2016 5:05
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 08 07:58:06 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 07 17:14:38 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	101	0.00
2		1,4-Dioxane	0.781	0.837	-7.2	103	0.00
3		n-Nitrosodimethylamine	0.836	0.811	3.0	95	0.00
4	S	2-Fluorophenol	1.066	1.159	-8.7	111	0.00
5	S	Phenol-d6	1.320	1.418	-7.4	114	0.00
6		bis(2-Chloroethyl)ether	1.431	1.396	2.4	99	0.00
7		n-Nitroso-di-n-propylamine	1.130	1.111	1.7	102	0.00
8	I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
9	S	Nitrobenzene-d5	0.336	0.327	2.7	96	-0.02
10		Nitrobenzene	0.304	0.312	-2.6	108	0.00
11		bis(2-Chloroethoxy)methane	0.507	0.505	0.4	109	0.00
12		Naphthalene	1.079	1.158	-7.3	105	0.00
13		Hexachlorobutadiene	0.161	0.172	-6.8	103	0.00
14		2-Methylnaphthalene	0.696	0.744	-6.9	108	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	111	0.00
16	S	2,4,6-Tribromophenol	0.142	0.155	-9.2	129	-0.02
17	S	2-Fluorobiphenyl	1.517	1.560	-2.8	104	0.00
18		Acenaphthylene	2.263	2.176	3.8	105	0.00
19		2,6-Dinitrotoluene	0.240	0.225	6.2	113	-0.02
20		Acenaphthene	1.383	1.435	-3.8	109	0.00
21		2,4-Dinitrotoluene	0.285	0.228	20.0	107	-0.02
22		Fluorene	1.691	1.731	-2.4	110	-0.02
23		4-Chlorophenyl-phenylether	0.829	0.864	-4.2	108	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	96	0.00
25		4-Bromophenyl-phenylether	0.180	0.190	-5.6	108	0.00
26		Hexachlorobenzene	0.184	0.199	-8.2	106	0.00
27		Pentachlorophenol	0.070	0.070	0.0	134	0.00
28		Phenanthrene	1.158	1.283	-10.8	109	-0.01
29		Anthracene	1.032	1.104	-7.0	113	-0.01
30		Fluoranthene	1.250	1.389	-11.1	114	-0.02
31	I	Chrysene-d12	1.000	1.000	0.0	115	-0.03
32		Benzidine	0.333	0.316	5.1	121	0.00
33		Pyrene	1.262	1.616	-28.1#	123	0.00
34	S	Terphenyl-d14	0.891	1.012	-13.6	107	0.00
35		Benzo(a)anthracene	7.222	8.052	-11.5	108	0.00
36		3,3'-Dichlorobenzidine	0.349	0.378	-8.3	137	0.00
37		Chrysene	5.098	2.822	44.6#	63	-0.03
38		Bis(2-ethylhexyl)phthalate	0.716	0.538	24.9	104	0.00
39		Indeno(1,2,3-cd)pyrene	1.524	1.679	-10.2	104	0.00
40	I	Perylene-d12	1.000	1.000	0.0	102	0.00
41		Benzo(b)fluoranthene	1.222	1.272	-4.1	105	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.268	1.353	-6.7	109	0.00
43 C	Benzo(a)pyrene	1.146	1.214	-5.9	111	0.00
44	Dibenzo(a,h)anthracene	1.068	1.109	-3.8	108	-0.02
45	Benzo(g,h,i)perylene	1.120	1.177	-5.1	105	-0.02

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