Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE060816\
Data File : BE091900.D
Acq On : 8 Jun 2016 14:08
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 09 05:25:32 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 08 16:11:34 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	102	0.00
2		1,4-Dioxane	0.781	0.766	1.9	95	-0.02
3		n-Nitrosodimethylamine	0.836	0.783	6.3	92	0.00
4	S	2-Fluorophenol	1.066	1.005	5.7	96	0.00
5	S	Phenol-d6	1.320	1.242	5.9	100	-0.02
6		bis(2-Chloroethyl)ether	1.431	1.387	3.1	99	0.00
7		n-Nitroso-di-n-propylamine	1.130	1.058	6.4	98	0.00
8	I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
9	S	Nitrobenzene-d5	0.336	0.304	9.5	90	-0.02
10		Nitrobenzene	0.304	0.300	1.3	104	0.00
11		bis(2-Chloroethoxy)methane	0.507	0.476	6.1	103	0.00
12		Naphthalene	1.079	1.155	-7.0	104	0.00
13		Hexachlorobutadiene	0.161	0.172	-6.8	104	0.00
14		2-Methylnaphthalene	0.696	0.727	-4.5	106	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	108	0.00
16	S	2,4,6-Tribromophenol	0.142	0.134	5.6	110	-0.02
17	S	2-Fluorobiphenyl	1.517	1.597	-5.3	105	0.00
18		Acenaphthylene	2.263	2.171	4.1	103	0.00
19		2,6-Dinitrotoluene	0.240	0.217	9.6	106	-0.02
20		Acenaphthene	1.383	1.410	-2.0	105	0.00
21		2,4-Dinitrotoluene	0.285	0.236	17.2	108	-0.02
22		Fluorene	1.691	1.751	-3.5	109	-0.02
23		4-Chlorophenyl-phenylether	0.829	0.874	-5.4	107	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
25		4-Bromophenyl-phenylether	0.180	0.186	-3.3	108	0.00
26		Hexachlorobenzene	0.184	0.209	-13.6	113	0.00
27		Pentachlorophenol	0.070	0.090	-28.6#	176#	0.00
28		Phenanthrene	1.158	1.278	-10.4	111	-0.01
29		Anthracene	1.032	1.075	-4.2	112	0.00
30		Fluoranthene	1.250	1.263	-1.0	105	-0.02
31	I	Chrysene-d12	1.000	1.000	0.0	101	0.00
32		Benzidine	0.333	0.605	-81.7#	205#	0.00
33		Pyrene	1.262	1.676	-32.8#	112	0.00
34	S	Terphenyl-d14	0.891	1.145	-28.5#	106	0.00
35		Benzo(a)anthracene	7.222	8.431	-16.7	100	0.00
36		3,3'-Dichlorobenzidine	0.349	0.326	6.6	104	0.00
37		Chrysene	5.098	3.937	22.8	77	0.00
38		Bis(2-ethylhexyl)phthalate	0.716	0.447	37.6#	76	0.00
39		Indeno(1,2,3-cd)pyrene	1.524	1.735	-13.8	95	0.00
40	I	Perylene-d12	1.000	1.000	0.0	95	0.00
41		Benzo(b)fluoranthene	1.222	1.196	2.1	92	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.268	1.305	-2.9	99	0.00
43 C	Benzo(a)pyrene	1.146	1.103	3.8	95	0.00
44	Dibenzo(a,h)anthracene	1.068	1.059	0.8	97	-0.02
45	Benzo(g,h,i)perylene	1.120	1.125	-0.4	95	-0.02

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