Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE052516\
Data File : BE091874.D
Acq On : 26 May 2016 00:52
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 26 02:04:23 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE052516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 25 19:42:30 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	126	0.00
2		1,4-Dioxane	0.737	0.570	22.7	105	0.00
3		n-Nitrosodimethylamine	0.918	0.767	16.4	120	0.00
4	S	2-Fluorophenol	1.429	1.237	13.4	119	0.00
5	S	Phenol-d6	1.874	1.824	2.7	140	0.00
6		bis(2-Chloroethyl)ether	1.569	1.463	6.8	129	-0.02
7		n-Nitroso-di-n-propylamine	1.452	1.335	8.1	135	0.00
8	I	Naphthalene-d8	1.000	1.000	0.0	141	0.00
9	S	Nitrobenzene-d5	0.406	0.374	7.9	141	0.00
10		Nitrobenzene	0.394	0.360	8.6	146	0.00
11		bis(2-Chloroethoxy)methane	0.474	0.364	23.2	136	0.00
12		Naphthalene	1.080	1.051	2.7	145	0.00
13		Hexachlorobutadiene	0.166	0.159	4.2	139	0.00
14		2-Methylnaphthalene	0.721	0.696	3.5	146	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	142	0.00
16	S	2,4,6-Tribromophenol	0.223	0.176	21.1	126	0.00
17	S	2-Fluorobiphenyl	1.487	1.432	3.7	142	-0.02
18		Acenaphthylene	2.458	2.295	6.6	145	0.00
19		2,6-Dinitrotoluene	0.341	0.262	23.2	130	-0.02
20		Acenaphthene	1.337	1.268	5.2	143	0.00
21		2,4-Dinitrotoluene	0.393	0.276	29.8#	133	0.00
22		Fluorene	1.755	1.613	8.1	139	0.00
23		4-Chlorophenyl-phenylether	0.883	0.786	11.0	133	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	124	0.00
25		4-Bromophenyl-phenylether	0.179	0.176	1.7	131	0.00
26		Hexachlorobenzene	0.170	0.175	-2.9	132	0.00
27		Pentachlorophenol	0.066	0.062	6.1	151#	-0.02
28		Phenanthrene	1.090	1.023	6.1	126	0.00
29		Anthracene	0.999	0.962	3.7	131	0.00
30		Fluoranthene	1.273	1.096	13.9	117	-0.02
31	I	Chrysene-d12	1.000	1.000	0.0	99	0.00
32		Benzidine	0.400	0.438	-9.5	136	0.00
33		Pyrene	1.061	1.170	-10.3	119	0.00
34	S	Terphenyl-d14	0.755	0.789	-4.5	110	0.00
35		Benzo(a)anthracene	1.333	1.393	-4.5	110	0.00
36		3,3'-Dichlorobenzidine	0.732	0.750	-2.5	112	0.00
37		Chrysene	1.222	1.269	-3.8	105	0.00
38		Bis(2-ethylhexyl)phthalate	0.484	0.910	-88.0#	219#	0.00
39		Indeno(1,2,3-cd)pyrene	1.250	1.248	0.2	106	0.00
40	I	Perylene-d12	1.000	1.000	0.0	100	0.00
41		Benzo(b)fluoranthene	1.349	1.230	8.8	104	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.225	1.281	-4.6	102	-0.01
43 C	Benzo(a)pyrene	1.223	1.217	0.5	106	0.00
44	Dibenzo(a,h)anthracene	1.172	1.157	1.3	105	0.00
45	Benzo(g,h,i)perylene	1.189	1.181	0.7	105	0.00

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