Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE052516\
Data File : BE091869.D
Acq On : 25 May 2016 21:44
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 26 13:43:32 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	115	0.00
2		1,4-Dioxane	0.737	0.648	12.1	109	0.00
3		n-Nitrosodimethylamine	0.918	0.782	14.8	112	0.00
4	S	2-Fluorophenol	1.429	1.277	10.6	112	0.00
5	S	Phenol-d6	1.874	1.734	7.5	121	0.00
6		bis(2-Chloroethyl)ether	1.569	1.459	7.0	117	-0.02
7		n-Nitroso-di-n-propylamine	1.452	1.286	11.4	119	0.00
8	I	Naphthalene-d8	1.000	1.000	0.0	126	0.00
9	S	Nitrobenzene-d5	0.406	0.363	10.6	123	0.00
10		Nitrobenzene	0.394	0.359	8.9	131	0.00
11		bis(2-Chloroethoxy)methane	0.474	0.426	10.1	143	0.00
12		Naphthalene	1.080	1.045	3.2	129	0.00
13		Hexachlorobutadiene	0.166	0.158	4.8	123	0.00
14		2-Methylnaphthalene	0.721	0.700	2.9	132	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	130	0.00
16	S	2,4,6-Tribromophenol	0.223	0.178	20.2	117	0.00
17	S	2-Fluorobiphenyl	1.487	1.435	3.5	130	0.00
18		Acenaphthylene	2.458	2.267	7.8	132	0.00
19		2,6-Dinitrotoluene	0.341	0.277	18.8	126	-0.02
20		Acenaphthene	1.337	1.294	3.2	133	0.00
21		2,4-Dinitrotoluene	0.393	0.286	27.2#	126	0.00
22		Fluorene	1.755	1.659	5.5	131	0.00
23		4-Chlorophenyl-phenylether	0.883	0.810	8.3	125	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	119	-0.01
25		4-Bromophenyl-phenylether	0.179	0.172	3.9	123	0.00
26		Hexachlorobenzene	0.170	0.173	-1.8	126	0.00
27		Pentachlorophenol	0.066	0.059	10.6	139	-0.02
28		Phenanthrene	1.090	1.028	5.7	122	0.00
29		Anthracene	0.999	0.950	4.9	124	0.00
30		Fluoranthene	1.273	1.114	12.5	114	-0.02
31	I	Chrysene-d12	1.000	1.000	0.0	99	0.00
32		Benzidine	0.400	0.394	1.5	121	0.00
33		Pyrene	1.061	1.143	-7.7	116	0.00
34	S	Terphenyl-d14	0.755	0.776	-2.8	108	0.00
35		Benzo(a)anthracene	1.333	1.284	3.7	101	0.00
36		3,3'-Dichlorobenzidine	0.732	0.700	4.4	105	0.00
37		Chrysene	1.222	1.215	0.6	101	0.00
38		Bis(2-ethylhexyl)phthalate	0.484	0.481	0.6	116	0.00
39		Indeno(1,2,3-cd)pyrene	1.250	1.220	2.4	104	0.00
40	I	Perylene-d12	1.000	1.000	0.0	101	0.00
41		Benzo(b)fluoranthene	1.349	1.195	11.4	102	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.225	1.262	-3.0	102	-0.01
43 C	Benzo(a)pyrene	1.223	1.158	5.3	103	0.00
44	Dibenzo(a,h)anthracene	1.172	1.117	4.7	103	0.00
45	Benzo(g,h,i)perylene	1.189	1.144	3.8	103	0.00

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