Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE060716\
Data File : BE091886.D
Acq On : 7 Jun 2016 20:20
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 08 07:15:16 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	105	0.00
2		1,4-Dioxane	0.781	0.779	0.3	99	0.00
3		n-Nitrosodimethylamine	0.836	0.831	0.6	100	0.00
4	S	2-Fluorophenol	1.066	1.109	-4.0	110	0.00
5	S	Phenol-d6	1.320	1.348	-2.1	112	0.00
6		bis(2-Chloroethyl)ether	1.431	1.413	1.3	104	0.00
7		n-Nitroso-di-n-propylamine	1.130	1.089	3.6	104	0.00
8	I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
9	S	Nitrobenzene-d5	0.336	0.347	-3.3	104	-0.02
10		Nitrobenzene	0.304	0.305	-0.3	108	0.00
11		bis(2-Chloroethoxy)methane	0.507	0.505	0.4	111	0.00
12		Naphthalene	1.079	1.141	-5.7	105	0.00
13		Hexachlorobutadiene	0.161	0.170	-5.6	105	0.00
14		2-Methylnaphthalene	0.696	0.722	-3.7	107	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
16	S	2,4,6-Tribromophenol	0.142	0.139	2.1	116	-0.02
17	S	2-Fluorobiphenyl	1.517	1.583	-4.4	106	0.00
18		Acenaphthylene	2.263	2.149	5.0	103	0.00
19		2,6-Dinitrotoluene	0.240	0.217	9.6	109	-0.02
20		Acenaphthene	1.383	1.409	-1.9	107	0.00
21		2,4-Dinitrotoluene	0.285	0.242	15.1	113	-0.02
22		Fluorene	1.691	1.732	-2.4	110	-0.02
23		4-Chlorophenyl-phenylether	0.829	0.881	-6.3	110	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
25		4-Bromophenyl-phenylether	0.180	0.194	-7.8	109	0.00
26		Hexachlorobenzene	0.184	0.203	-10.3	107	0.00
27		Pentachlorophenol	0.070	0.062	11.4	118	0.00
28		Phenanthrene	1.158	1.287	-11.1	108	-0.01
29		Anthracene	1.032	1.106	-7.2	112	0.00
30		Fluoranthene	1.250	1.347	-7.8	109	-0.02
31	I	Chrysene-d12	1.000	1.000	0.0	109	0.00
32		Benzidine	0.333	0.297	10.8	108	0.00
33		Pyrene	1.262	1.598	-26.6#	115	0.00
34	S	Terphenyl-d14	0.891	1.099	-23.3	110	0.00
35		Benzo(a)anthracene	7.222	8.032	-11.2	103	0.00
36		3,3'-Dichlorobenzidine	0.349	0.351	-0.6	121	0.00
37		Chrysene	5.098	4.526	11.2	95	0.00
38		Bis(2-ethylhexyl)phthalate	0.716	0.720	-0.6	132	0.00
39		Indeno(1,2,3-cd)pyrene	1.524	1.743	-14.4	103	0.00
40	I	Perylene-d12	1.000	1.000	0.0	101	0.00
41		Benzo(b)fluoranthene	1.222	1.280	-4.7	104	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE060716\
Data File : BE091886.D
Acq On : 7 Jun 2016 20:20
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 08 07:15:16 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)
42	Benzo(k)fluoranthene	1.268	1.296	-2.2	104	0.00
43 C	Benzo(a)pyrene	1.146	1.195	-4.3	108	0.00
44	Dibenzo(a,h)anthracene	1.068	1.076	-0.7	104	0.00
45	Benzo(g,h,i)perylene	1.120	1.152	-2.9	102	0.00

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