Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA E\Data\BE101217\
Data File : BE094311.D
Acq On : 12 Oct 2017 12:02
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Oct 12 22:51:59 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE100617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 06 15:16:13 2017
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	135	0.01
2		1,4-Dioxane	0.725	0.603	16.8	105	0.01
3		n-Nitrosodimethylamine	0.777	0.663	14.7	115	0.03
4	S	2-Fluorophenol	1.333	1.398	-4.9	142	0.03
5	S	Phenol-d6	1.932	2.051	-6.2	145	0.03
6		bis(2-Chloroethyl)ether	1.569	1.562	0.4	137	0.01
7	I	Naphthalene-d8	1.000	1.000	0.0	141	0.02
8	S	Nitrobenzene-d5	0.333	0.347	-4.2	148	0.02
9		Naphthalene	4.505	4.498	0.2	138	0.02
10		Hexachlorobutadiene	0.171	0.178	-4.1	145	0.00
11	SURR	2-Methylnaphthalene-d10	0.616	0.599	2.8	135	0.02
12		2-Methylnaphthalene	0.740	0.716	3.2	135	0.02
13	I	Acenaphthene-d10	1.000	1.000	0.0	126	0.01
14	S	2,4,6-Tribromophenol	0.174	0.165	5.2	128	0.03
15	S	2-Fluorobiphenyl	1.370	1.447	-5.6	132	0.01
16		Acenaphthylene	2.026	2.135	-5.4	132	0.01
17		Acenaphthene	1.338	1.329	0.7	122	0.01
18		Fluorene	1.753	1.659	5.4	118	0.01
19	I	Phenanthrene-d10	1.000	1.000	0.0	64	0.03
20		4-Bromophenyl-phenylether	0.230	0.193	16.1	53	0.03
21		Hexachlorobenzene	0.207	0.377	-82.1#	118	0.03
22		Pentachlorophenol	0.056	0.036	35.7#	41#	0.07
23		Phenanthrene	1.386	1.664	-20.1	76	0.03
24		Anthracene	1.057	1.235	-16.8	77	0.03
25	SURR	Fluoranthene-d10	3.995	6.779	-69.7#	107	0.02
26		Fluoranthene	1.216	2.025	-66.5#	106	0.02
27	I	Chrysene-d12	1.000	1.000	0.0	99	0.02
28		Pyrene	1.376	1.468	-6.7	105	0.02
29	S	Terphenyl-d14	0.755	0.791	-4.8	103	0.01
30		Benzo(a)anthracene	1.314	1.330	-1.2	99	0.03
31		Chrysene	1.297	1.335	-2.9	102	0.03
32		Bis(2-ethylhexyl)phthalate	3.027	3.653	-20.7	126	0.01
33		Indeno(1,2,3-cd)pyrene	1.287	1.078	16.2	82	0.10
34	I	Perylene-d12	1.000	1.000	0.0	101	0.05
35		Benzo(b)fluoranthene	1.345	1.297	3.6	97	0.04
36		Benzo(k)fluoranthene	1.368	1.337	2.3	99	0.04
37	C	Benzo(a)pyrene	1.272	1.240	2.5	99	0.05
38		Dibenzo(a,h)anthracene	1.179	1.014	14.0	87	0.09
39		Benzo(g,h,i)perylene	1.204	0.837	30.5#	71	0.11

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA E\Data\BE101217\
Data File : BE094311.D
Acq On : 12 Oct 2017 12:02
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Oct 12 22:51:59 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE100617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 06 15:16:13 2017
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)

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