Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE032319\
Data File : BE099139.D
Acq On : 24 Mar 2019 8:40
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 24 23:03:18 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE032219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Mar 23 04:40:11 2019
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	148	0.00
2		1,4-Dioxane	0.621	0.602	3.1	165#	-0.01
3		n-Nitrosodimethylamine	0.320	0.368	-15.0	184#	-0.02
4	S	2-Fluorophenol	1.191	1.393	-17.0	194#	0.00
5	S	Phenol-d6	1.715	2.152	-25.5#	212#	0.00
6		bis(2-Chloroethyl)ether	1.394	1.630	-16.9	193#	-0.01
7	I	Naphthalene-d8	1.000	1.000	0.0	170#	-0.01
8	S	Nitrobenzene-d5	0.175	0.396	-126.3#	483#	0.00
9		Naphthalene	4.714	5.175	-9.8	211#	0.00
10		Hexachlorobutadiene	0.231	0.276	-19.5	226#	0.00
11	SURR	2-Methylnaphthalene-d10	0.720	0.835	-16.0	225#	-0.01
12		2-Methylnaphthalene	0.807	0.936	-16.0	228#	-0.01
13	I	Acenaphthene-d10	1.000	1.000	0.0	187#	0.00
14	S	2,4,6-Tribromophenol	0.118	0.250	-111.9#	512#	0.00
15	S	2-Fluorobiphenyl	1.591	1.859	-16.8	238#	-0.01
16		Acenaphthylene	2.135	2.327	-9.0	241#	-0.01
17		Acenaphthene	1.266	1.353	-6.9	234#	0.00
18		Fluorene	1.847	1.893	-2.5	222#	0.00
19	I	Phenanthrene-d10	1.000	1.000	0.0	170#	0.00
20		4-Bromophenyl-phenylether	0.221	0.266	-20.4	240#	0.00
21		Hexachlorobenzene	0.206	0.250	-21.4	246#	0.00
22		Pentachlorophenol	0.061	0.105	-72.1#	426#	-0.01
23		Phenanthrene	1.188	1.236	-4.0	209#	0.00
24		Anthracene	1.124	1.164	-3.6	212#	-0.01
25	SURR	Fluoranthene-d10	5.336	5.222	2.1	192#	0.00
26		Fluoranthene	1.671	1.596	4.5	195#	0.00
27	I	Chrysene-d12	1.000	1.000	0.0	156#	0.00
28		Pyrene	1.357	1.442	-6.3	196#	0.00
29	S	Terphenyl-d14	0.846	1.014	-19.9	212#	0.00
30		Benzo(a)anthracene	1.409	1.489	-5.7	194#	0.00
31		Chrysene	1.408	1.506	-7.0	198#	0.00
32		Bis(2-ethylhexyl)phthalate	1.829	2.609	-42.6#	314#	-0.01
33		Indeno(1,2,3-cd)pyrene	1.514	1.693	-11.8	207#	-0.01
34	I	Perylene-d12	1.000	1.000	0.0	166#	0.00
35		Benzo(b)fluoranthene	1.384	1.395	-0.8	203#	0.00
36		Benzo(k)fluoranthene	1.356	1.373	-1.3	201#	-0.01
37	C	Benzo(a)pyrene	1.279	1.308	-2.3	207#	0.00
38		Dibenzo(a,h)anthracene	1.244	1.301	-4.6	212#	-0.02
39		Benzo(g,h,i)perylene	1.278	1.303	-2.0	203#	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE032319\
Data File : BE099139.D
Acq On : 24 Mar 2019 8:40
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 24 23:03:18 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE032219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Mar 23 04:40:11 2019
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