Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\terastorage\svoasrv\HPCHEM1\BNA_N\Data\BN062819\
Data File : BN006650.D
Acq On : 29 Jun 2019 13:08
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Jul 01 08:16:51 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN062819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 28 13:47:56 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	147	0.00
2		1,4-Dioxane	0.554	0.496	10.5	141	0.00
3		n-Nitrosodimethylamine	0.435	0.434	0.2	166#	0.00
4	S	2-Fluorophenol	0.981	1.076	-9.7	169#	0.00
5	S	Phenol-d6	1.299	1.517	-16.8	182#	0.00
6		bis(2-Chloroethyl)ether	1.158	1.162	-0.3	142	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	149	0.00
8	S	Nitrobenzene-d5	0.269	0.297	-10.4	176#	-0.01
9		Naphthalene	1.051	1.113	-5.9	155#	-0.01
10		Hexachlorobutadiene	0.216	0.233	-7.9	156#	-0.01
11	SURR	2-Methylnaphthalene-d10	0.630	0.647	-2.7	150	0.00
12		2-Methylnaphthalene	0.726	0.759	-4.5	151#	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	132	-0.01
14	S	2,4,6-Tribromophenol	0.188	0.196	-4.3	152#	-0.01
15	S	2-Fluorobiphenyl	1.482	1.629	-9.9	145	-0.01
16		Acenaphthylene	1.907	2.136	-12.0	151#	0.00
17		Acenaphthene	1.256	1.353	-7.7	139	0.00
18		Fluorene	1.585	1.658	-4.6	136	-0.01
19	I	Phenanthrene-d10	1.000	1.000	0.0	119	-0.01
20		4-Bromophenyl-phenylether	0.222	0.249	-12.2	135	-0.01
21		Hexachlorobenzene	0.280	0.304	-8.6	125	0.00
22		Pentachlorophenol	0.045	0.051	-13.3	171#	-0.02
23		Phenanthrene	1.137	1.224	-7.7	127	0.00
24		Anthracene	0.987	1.076	-9.0	134	0.00
25	SURR	Fluoranthene-d10	1.175	1.106	5.9	111	0.00
26		Fluoranthene	1.382	1.314	4.9	110	0.00
27	I	Chrysene-d12	1.000	1.000	0.0	77	-0.01
28		Pyrene	1.400	1.957	-39.8#	105	0.00
29	S	Terphenyl-d14	0.838	1.195	-42.6#	109	0.00
30		Benzo(a)anthracene	1.308	1.495	-14.3	89	-0.01
31		Chrysene	1.410	1.515	-7.4	78	-0.01
32		Bis(2-ethylhexyl)phthalate	0.658	1.257	-91.0#	164#	-0.01
33		Indeno(1,2,3-cd)pyrene	1.427	1.595	-11.8	74	-0.02
34	I	Perylene-d12	1.000	1.000	0.0	67	-0.02
35		Benzo(b)fluoranthene	1.361	1.420	-4.3	69	-0.02
36		Benzo(k)fluoranthene	1.450	1.453	-0.2	65	-0.02
37	C	Benzo(a)pyrene	1.272	1.372	-7.9	69	-0.02
38		Dibenzo(a,h)anthracene	1.111	1.303	-17.3	73	-0.02
39		Benzo(g,h,i)perylene	1.109	1.337	-20.6	75	-0.02

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\terastorage\svoasrv\HPCHEM1\BNA_N\Data\BN062819\
Data File : BN006650.D
Acq On : 29 Jun 2019 13:08
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_N
LabSampleId :
SSTDCCC0.4

Quant Time: Jul 01 08:16:51 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN062819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 28 13:47:56 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