Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN032219\
Data File : BN004917.D
Acq On : 23 Mar 2019 10:19
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Mar 24 00:55:42 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN032019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 20 16:06:20 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	282	0.00
2		1,4-Dioxane	0.400	0.383	4.3	268	0.00
3		n-Nitrosodimethylamine	0.400	0.390	2.5	297	0.00
4	S	2-Fluorophenol	0.400	0.445	-11.2	308	0.00
5	S	Phenol-d6	0.400	0.474	-18.5	336	0.00
6		bis(2-Chloroethyl)ether	0.400	0.426	-6.5	308	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	289	-0.01
8	S	Nitrobenzene-d5	0.400	0.441	-10.2	321	-0.01
9		Naphthalene	0.400	0.412	-3.0	294	-0.03
10		Hexachlorobutadiene	0.400	0.394	1.5	282	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.393	1.8	284	-0.01
12		2-Methylnaphthalene	0.400	0.391	2.3	279	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	254	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.374	6.5	253	-0.01
15	S	2-Fluorobiphenyl	0.400	0.421	-5.2	262	-0.01
16		Acenaphthylene	0.400	0.434	-8.5	289	-0.01
17		Acenaphthene	0.400	0.408	-2.0	256	-0.01
18		Fluorene	0.400	0.381	4.8	237	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	226	-0.01
20		4-Bromophenyl-phenylether	0.400	0.399	0.3	227	-0.01
21		Hexachlorobenzene	0.400	0.383	4.3	214	-0.01
22		Pentachlorophenol	0.400	0.688	-72.0#	450	-0.01
23		Phenanthrene	0.400	0.397	0.8	222	-0.01
24		Anthracene	0.400	0.409	-2.2	236	-0.01
25	SURR	Fluoranthene-d10	0.400	0.391	2.3	223	0.00
26		Fluoranthene	0.400	0.372	7.0	209	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	169	-0.01
28		Pyrene	0.400	0.499	-24.7	207	0.00
29	S	Terphenyl-d14	0.400	0.493	-23.2	206	0.00
30		Benzo(a)anthracene	0.400	0.423	-5.7	181	-0.01
31		Chrysene	0.400	0.396	1.0	162	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.703	-75.7#	322	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.274	31.5#	112	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	134	-0.02
35		Benzo(b)fluoranthene	0.400	0.401	-0.3	132	-0.01
36		Benzo(k)fluoranthene	0.400	0.382	4.5	130	-0.01
37	C	Benzo(a)pyrene	0.400	0.385	3.8	130	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.335	16.3	111	-0.01
39		Benzo(g,h,i)perylene	0.400	0.328	18.0	108	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN032219\
Data File : BN004917.D
Acq On : 23 Mar 2019 10:19
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Mar 24 00:55:42 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN032019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 20 16:06:20 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	Amount	Calc.	%Dev	Area%	Dev(min)

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