Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA N\Data\BN030118\
Data File : BN000207.D
Acq On : 01 Mar 2018 08:44
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Mar 01 16:55:36 2018
Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-SIM-BN022718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 27 15:07:57 2018
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	68	0.00
2		1,4-Dioxane	0.400	0.411	-2.7	75	0.00
3		n-Nitrosodimethylamine	0.400	0.360	10.0	61	0.00
4	S	2-Fluorophenol	0.400	0.437	-9.2	73	0.00
5	S	Phenol-d6	0.400	0.427	-6.7	74	0.00
6		bis(2-Chloroethyl)ether	0.400	0.415	-3.7	68	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	70	0.00
8	S	Nitrobenzene-d5	0.400	0.396	1.0	69	0.00
9		Naphthalene	0.400	0.405	-1.3	69	0.00
10		Hexachlorobutadiene	0.400	0.407	-1.7	69	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.409	-2.2	70	0.00
12		2-Methylnaphthalene	0.400	0.405	-1.3	70	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	75	0.00
14	S	2,4,6-Tribromophenol	0.400	0.408	-2.0	78	0.00
15	S	2-Fluorobiphenyl	0.400	0.387	3.3	71	0.00
16		Acenaphthylene	0.400	0.355	11.3	68	0.00
17		Acenaphthene	0.400	0.381	4.8	71	-0.01
18		Fluorene	0.400	0.379	5.3	72	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	78	0.00
20		4-Bromophenyl-phenylether	0.400	0.373	6.8	73	0.00
21		Hexachlorobenzene	0.400	0.375	6.3	73	0.00
22		Pentachlorophenol	0.400	0.418	-4.5	84	0.00
23		Phenanthrene	0.400	0.375	6.3	73	0.00
24		Anthracene	0.400	0.357	10.8	71	0.00
25	SURR	Fluoranthene-d10	0.400	0.369	7.8	73	0.00
26		Fluoranthene	0.400	0.363	9.3	72	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	79	0.00
28		Pyrene	0.400	0.372	7.0	73	0.00
29	S	Terphenyl-d14	0.400	0.377	5.8	73	0.00
30		Benzo(a)anthracene	0.400	0.392	2.0	77	0.00
31		Chrysene	0.400	0.386	3.5	74	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.366	8.5	78	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.366	8.5	69	0.00
34	I	Perylene-d12	0.400	0.400	0.0	76	0.00
35		Benzo(b)fluoranthene	0.400	0.366	8.5	70	0.00
36		Benzo(k)fluoranthene	0.400	0.383	4.3	72	0.00
37	C	Benzo(a)pyrene	0.400	0.366	8.5	70	0.00
38		Dibenzo(a,h)anthracene	0.400	0.363	9.3	70	0.00
39		Benzo(g,h,i)perylene	0.400	0.367	8.3	69	0.00

Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA N\Data\BN030118\
Data File : BN000207.D
Acq On : 01 Mar 2018 08:44
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Mar 01 16:55:36 2018
Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-SIM-BN022718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 27 15:07:57 2018
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