Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE070217\
Data File : BE093418.D
Acq On : 2 Jul 2017 17:44
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 03 06:49:13 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE070217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Jul 02 13:43:38 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	98	0.00
2		1,4-Dioxane	0.400	0.383	4.3	94	0.00
3		n-Nitrosodimethylamine	0.400	0.379	5.3	109	0.00
4	S	2-Fluorophenol	0.400	0.378	5.5	95	0.00
5	S	Phenol-d6	0.400	0.384	4.0	100	0.00
6		bis(2-Chloroethyl)ether	0.400	0.387	3.3	97	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	99	0.00
8	S	Nitrobenzene-d5	0.400	0.383	4.3	102	0.00
9		Naphthalene	0.400	0.390	2.5	98	0.00
10		Hexachlorobutadiene	0.400	0.393	1.8	98	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.390	2.5	99	0.00
12		2-Methylnaphthalene	0.400	0.389	2.8	99	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	101	0.00
14	S	2,4,6-Tribromophenol	0.400	0.324	19.0	90	0.00
15	S	2-Fluorobiphenyl	0.400	0.385	3.8	101	0.00
16		Acenaphthylene	0.400	0.362	9.5	97	0.00
17		Acenaphthene	0.400	0.381	4.8	100	0.00
18		Fluorene	0.400	0.384	4.0	100	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	91	0.00
20		4-Bromophenyl-phenylether	0.400	0.405	-1.3	100	0.00
21		Hexachlorobenzene	0.400	0.391	2.3	100	0.00
22		Pentachlorophenol	0.400	0.332	17.0	82	0.00
23		Phenanthrene	0.400	0.400	0.0	100	0.00
24		Anthracene	0.400	0.390	2.5	95	0.00
25	SURR	Fluoranthene-d10	0.400	0.388	3.0	97	0.00
26		Fluoranthene	0.400	0.395	1.3	99	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	97	0.00
28		Pyrene	0.400	0.393	1.8	97	0.00
29	S	Terphenyl-d14	0.400	0.392	2.0	97	0.00
30		Benzo(a)anthracene	0.400	0.382	4.5	97	0.00
31		Chrysene	0.400	0.388	3.0	95	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.332	17.0	98	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.390	2.5	97	0.00
34	I	Perylene-d12	0.400	0.400	0.0	97	0.00
35		Benzo(b)fluoranthene	0.400	0.381	4.8	95	0.00
36		Benzo(k)fluoranthene	0.400	0.399	0.3	101	0.00
37	C	Benzo(a)pyrene	0.400	0.378	5.5	97	0.00
38		Dibenzo(a,h)anthracene	0.400	0.390	2.5	98	0.00
39		Benzo(g,h,i)perylene	0.400	0.390	2.5	99	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE070217\
Data File : BE093418.D
Acq On : 2 Jul 2017 17:44
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 03 06:49:13 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE070217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Jul 02 13:43:38 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	Amount	Calc.	%Dev	Area%	Dev(min)

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