Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE111617\
Data File : BE094818.D
Acq On : 15 Nov 2017 17:09
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Nov 16 02:01:26 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 23 22:40:47 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	66	0.01
2		1,4-Dioxane	0.400	0.453	-13.2	65	-0.03
3		n-Nitrosodimethylamine	0.400	0.357	10.8	59	0.00
4	S	2-Fluorophenol	0.400	0.344	14.0	57	0.02
5	S	Phenol-d6	0.400	0.390	2.5	66	0.06
6		bis(2-Chloroethyl)ether	0.400	0.392	2.0	66	0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	69	0.02
8	S	Nitrobenzene-d5	0.400	0.358	10.5	64	0.05
9		Naphthalene	0.400	0.358	10.5	60	0.03
10		Hexachlorobutadiene	0.400	0.412	-3.0	69	-0.02
11	SURR	2-Methylnaphthalene-d10	0.400	0.359	10.3	60	0.03
12		2-Methylnaphthalene	0.400	0.355	11.3	60	0.03
13	I	Acenaphthene-d10	0.400	0.400	0.0	73	0.02
14	S	2,4,6-Tribromophenol	0.400	0.513	-28.2#	97	0.03
15	S	2-Fluorobiphenyl	0.400	0.347	13.3	62	0.02
16		Acenaphthylene	0.400	0.352	12.0	65	0.00
17		Acenaphthene	0.400	0.360	10.0	65	0.00
18		Fluorene	0.400	0.385	3.8	70	0.02
19	I	Phenanthrene-d10	0.400	0.400	0.0	94	0.03
20		4-Bromophenyl-phenylether	0.400	0.298	25.5#	71	0.03
21		Hexachlorobenzene	0.400	0.490	-22.5	100	0.00
22		Pentachlorophenol	0.400	0.426	-6.5	100	0.03
23		Phenanthrene	0.400	0.516	-29.0#	108	0.00
24		Anthracene	0.400	0.451	-12.7	103	0.00
25	SURR	Fluoranthene-d10	0.400	0.431	-7.7	91	0.01
26		Fluoranthene	0.400	0.422	-5.5	90	0.01
27	I	Chrysene-d12	0.400	0.400	0.0	99	0.01
28		Pyrene	0.400	0.363	9.3	91	0.01
29	S	Terphenyl-d14	0.400	0.364	9.0	91	0.00
30		Benzo(a)anthracene	0.400	0.372	7.0	92	0.01
31		Chrysene	0.400	0.397	0.8	101	0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.360	10.0	97	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.392	2.0	98	0.07
34	I	Perylene-d12	0.400	0.400	0.0	98	0.03
35		Benzo(b)fluoranthene	0.400	0.363	9.3	92	0.02
36		Benzo(k)fluoranthene	0.400	0.374	6.5	93	0.03
37	C	Benzo(a)pyrene	0.400	0.383	4.3	94	0.04
38		Dibenzo(a,h)anthracene	0.400	0.394	1.5	99	0.05
39		Benzo(g,h,i)perylene	0.400	0.418	-4.5	104	0.09

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE111617\
Data File : BE094818.D
Acq On : 15 Nov 2017 17:09
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Nov 16 02:01:26 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 23 22:40:47 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