Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE051616\
Data File : BE091788.D
Acq On : 17 May 2016 00:40
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 17 05:11:43 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE042616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 26 16:09:46 2016
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	5.000	5.000	0.0	125	0.00
2		1,4-Dioxane	0.400	0.297	25.8#	94	0.00
3		n-Nitrosodimethylamine	0.400	0.291	27.3#	93	0.02
4	S	2-Fluorophenol	0.400	0.358	10.5	115	0.00
5	S	Phenol-d6	0.400	0.359	10.3	119	0.00
6		bis(2-Chloroethyl)ether	0.400	0.361	9.8	111	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	136	0.00
8	S	Nitrobenzene-d5	0.400	0.324	19.0	118	0.00
9		Nitrobenzene	0.400	0.308	23.0	114	0.00
10		Naphthalene	0.400	0.383	4.3	129	0.00
11		Hexachlorobutadiene	0.400	0.383	4.3	133	-0.02
12		2-Methylnaphthalene	0.400	0.384	4.0	131	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	140	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.421	-5.2	121	0.00
15	S	2-Fluorobiphenyl	0.400	0.368	8.0	131	0.00
16		Acenaphthylene	0.400	0.377	5.8	150	0.00
17		Acenaphthene	0.400	0.371	7.3	131	0.00
18		Fluorene	0.400	0.366	8.5	128	-0.01
19	I	Phenanthrene-d10	5.000	5.000	0.0	129	0.00
20		4-Bromophenyl-phenylether	0.400	0.382	4.5	122	0.00
21		Hexachlorobenzene	0.400	0.405	-1.3	130	0.00
22		Pentachlorophenol	0.400	0.353	11.8	113	0.00
23		Phenanthrene	0.400	0.389	2.8	122	0.00
24		Anthracene	0.400	0.370	7.5	121	0.00
25		Fluoranthene	0.400	0.380	5.0	121	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	104	-0.02
27		Benzidine	0.400	0.414	-3.5	115	0.00
28		Pyrene	0.400	0.428	-7.0	115	0.00
29	S	Terphenyl-d14	0.400	0.402	-0.5	106	0.00
30		Benzo(a)anthracene	0.400	0.389	2.8	103	0.00
31		3,3'-Dichlorobenzidine	0.400	0.418	-4.5	125	-0.02
32		Chrysene	0.400	0.394	1.5	96	0.00
33		Bis(2-ethylhexyl)phthalate	0.400	0.559	-39.8#	188	-0.02
34		Indeno(1,2,3-cd)pyrene	0.400	0.431	-7.7	109	-0.01
35	I	Perylene-d12	5.000	5.000	0.0	115	-0.01
36		Benzo(b)fluoranthene	0.400	0.361	9.8	107	-0.01
37		Benzo(k)fluoranthene	0.400	0.373	6.8	112	-0.01
38	C	Benzo(a)pyrene	0.400	0.364	9.0	110	-0.01
39		Dibenzo(a,h)anthracene	0.400	0.370	7.5	109	-0.02
40		Benzo(g,h,i)perylene	0.400	0.376	6.0	111	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE051616\
Data File : BE091788.D
Acq On : 17 May 2016 00:40
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 17 05:11:43 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE042616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 26 16:09:46 2016
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