Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE050316\
Data File : BE091750.D
Acq On : 3 May 2016 19:18
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 04 01:40:36 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	151	0.00
2		1,4-Dioxane	0.400	0.298	25.5#	114	0.00
3		n-Nitrosodimethylamine	0.400	0.338	15.5	130	0.02
4	S	2-Fluorophenol	0.400	0.385	3.8	150	0.00
5	S	Phenol-d6	0.400	0.401	-0.3	161	0.00
6		bis(2-Chloroethyl)ether	0.400	0.395	1.3	148	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	157	0.00
8	S	Nitrobenzene-d5	0.400	0.392	2.0	165	0.00
9		Nitrobenzene	0.400	0.377	5.8	161	0.00
10		Naphthalene	0.400	0.414	-3.5	161	0.00
11		Hexachlorobutadiene	0.400	0.436	-9.0	175	-0.02
12		2-Methylnaphthalene	0.400	0.404	-1.0	159	-0.01
13	I	Acenaphthene-d10	5.000	5.000	0.0	153	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.475	-18.7	160	0.00
15	S	2-Fluorobiphenyl	0.400	0.399	0.3	155	0.00
16		Acenaphthylene	0.400	0.427	-6.7	186	0.00
17		Acenaphthene	0.400	0.397	0.8	153	0.00
18		Fluorene	0.400	0.404	-1.0	154	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	140	0.00
20		4-Bromophenyl-phenylether	0.400	0.426	-6.5	148	0.00
21		Hexachlorobenzene	0.400	0.430	-7.5	151	0.00
22		Pentachlorophenol	0.400	0.438	-9.5	164	0.00
23		Phenanthrene	0.400	0.411	-2.7	141	0.00
24		Anthracene	0.400	0.436	-9.0	156	0.00
25		Fluoranthene	0.400	0.462	-15.5	160	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	122	0.00
27		Benzidine	0.400	0.671	-67.8#	273	0.00
28		Pyrene	0.400	0.506	-26.5#	159	0.00
29	S	Terphenyl-d14	0.400	0.479	-19.7	149	0.00
30		Benzo(a)anthracene	0.400	0.434	-8.5	135	0.00
31		3,3'-Dichlorobenzidine	0.400	0.559	-39.8#	196	-0.02
32		Chrysene	0.400	0.462	-15.5	133	0.00
33		Bis(2-ethylhexyl)phthalate	0.400	0.979	-144.7#	386	-0.02
34		Indeno(1,2,3-cd)pyrene	0.400	0.475	-18.7	142	0.00
35	I	Perylene-d12	5.000	5.000	0.0	127	0.00
36		Benzo(b)fluoranthene	0.400	0.453	-13.2	148	0.00
37		Benzo(k)fluoranthene	0.400	0.398	0.5	132	0.00
38	C	Benzo(a)pyrene	0.400	0.436	-9.0	146	0.00
39		Dibenzo(a,h)anthracene	0.400	0.427	-6.7	139	-0.01
40		Benzo(g,h,i)perylene	0.400	0.431	-7.7	141	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE050316\
Data File : BE091750.D
Acq On : 3 May 2016 19:18
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 04 01:40:36 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