Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE050316\
Data File : BE091741.D
Acq On : 3 May 2016 12:59
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 04 01:08:44 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	122	0.00
2		1,4-Dioxane	0.752	0.600	20.2	96	0.02
3		n-Nitrosodimethylamine	0.908	0.747	17.7	102	0.02
4	S	2-Fluorophenol	1.228	1.053	14.3	108	0.00
5	S	Phenol-d6	1.629	1.337	17.9	106	0.02
6		bis(2-Chloroethyl)ether	1.433	1.341	6.4	113	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	119	0.00
8	S	Nitrobenzene-d5	0.323	0.259	19.8	102	0.01
9		Nitrobenzene	0.338	0.264	21.9	101	0.00
10		Naphthalene	1.010	1.018	-0.8	119	0.00
11		Hexachlorobutadiene	0.149	0.151	-1.3	123	-0.02
12		2-Methylnaphthalene	0.649	0.624	3.9	115	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	115	-0.01
14	S	2,4,6-Tribromophenol	0.157	0.113	28.0#	91	0.00
15	S	2-Fluorobiphenyl	1.414	1.400	1.0	116	0.00
16		Acenaphthylene	1.923	1.830	4.8	125	0.00
17		Acenaphthene	1.247	1.193	4.3	111	0.00
18		Fluorene	1.550	1.512	2.5	112	0.00
19	I	Phenanthrene-d10	1.000	1.000	0.0	115	0.00
20		4-Bromophenyl-phenylether	0.180	0.174	3.3	110	0.00
21		Hexachlorobenzene	0.180	0.182	-1.1	116	0.00
22		Pentachlorophenol	0.085	0.065	23.5	104	0.00
23		Phenanthrene	1.136	1.154	-1.6	114	0.00
24		Anthracene	1.025	0.972	5.2	111	0.00
25		Fluoranthene	1.188	1.143	3.8	109	0.00
26	I	Chrysene-d12	1.000	1.000	0.0	103	0.00
27		Benzidine	0.286	0.306	-7.0	136	0.00
28		Pyrene	1.130	1.194	-5.7	112	0.00
29	S	Terphenyl-d14	0.780	0.815	-4.5	110	0.00
30		Benzo(a)anthracene	1.221	1.172	4.0	101	0.00
31		3,3'-Dichlorobenzidine	0.623	0.541	13.2	103	-0.02
32		Chrysene	1.269	1.511	-19.1	116	0.00
33		Bis(2-ethylhexyl)phthalate	0.370	0.313	15.4	113	-0.02
34		Indeno(1,2,3-cd)pyrene	1.130	1.275	-12.8	114	-0.01
35	I	Perylene-d12	1.000	1.000	0.0	116	0.00
36		Benzo(b)fluoranthene	1.172	1.209	-3.2	123	-0.01
37		Benzo(k)fluoranthene	1.191	1.029	13.6	104	0.00
38	C	Benzo(a)pyrene	1.103	1.030	6.6	114	0.00
39		Dibenzo(a,h)anthracene	1.045	0.986	5.6	113	-0.02
40		Benzo(g,h,i)perylene	1.062	0.990	6.8	112	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE050316\
Data File : BE091741.D
Acq On : 3 May 2016 12:59
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 04 01:08:44 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	AvgRF	CCRF	%Dev	Area%	Dev(min)

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