Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072716\
Data File : BE092070.D
Acq On : 28 Jul 2016 6:59
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 28 07:33:10 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE072716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 27 17:13:26 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	114	-0.02
2		1,4-Dioxane	0.687	0.629	8.4	104	0.00
3		n-Nitrosodimethylamine	0.819	0.840	-2.6	129	0.00
4	S	2-Fluorophenol	1.114	1.370	-23.0	160#	0.00
5	S	Phenol-d6	1.487	1.874	-26.0#	162#	0.00
6		bis(2-Chloroethyl)ether	1.437	1.417	1.4	124	0.00
7		n-Nitroso-di-n-propylamine	1.187	1.339	-12.8	145	0.00
8	I	Naphthalene-d8	1.000	1.000	0.0	123	0.00
9	S	Nitrobenzene-d5	0.373	0.364	2.4	129	0.00
10		Nitrobenzene	0.362	0.378	-4.4	145	0.00
11		bis(2-Chloroethoxy)methane	0.417	0.414	0.7	140	-0.02
12		Naphthalene	1.095	1.140	-4.1	129	0.00
13		Hexachlorobutadiene	0.164	0.177	-7.9	136	0.00
14		2-Methylnaphthalene	0.707	0.769	-8.8	143	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	136	0.00
16	S	2,4,6-Tribromophenol	0.175	0.246	-40.6#	232#	0.00
17	S	2-Fluorobiphenyl	2.003	2.041	-1.9	141	0.00
18		Acenaphthylene	12.962	14.152	-9.2	161#	0.00
19		2,6-Dinitrotoluene	1.331	1.992	-49.7#	233#	0.00
20		Acenaphthene	1.366	1.382	-1.2	141	0.00
21		2,4-Dinitrotoluene	1.684	1.612	4.3	211#	0.00
22		Fluorene	2.145	2.352	-9.7	154#	0.00
23		4-Chlorophenyl-phenylether	1.055	1.116	-5.8	145	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	133	0.00
25		4-Bromophenyl-phenylether	0.206	0.217	-5.3	144	0.00
26		Hexachlorobenzene	0.213	0.234	-9.9	149	0.00
27		Pentachlorophenol	0.060	0.073	-21.7	303#	0.00
28		Phenanthrene	1.271	1.388	-9.2	150#	0.00
29		Anthracene	1.156	1.308	-13.1	162#	0.00
30		Fluoranthene	5.652	6.605	-16.9	177#	0.00
31	I	Chrysene-d12	1.000	1.000	0.0	114	0.00
32		Benzidine	0.246	0.501	-103.7#	324#	0.00
33		Pyrene	6.442	8.311	-29.0#	161#	0.00
34	S	Terphenyl-d14	0.931	1.144	-22.9	158#	0.00
35		Benzo(a)anthracene	2.391	2.694	-12.7	138	0.00
36		3,3'-Dichlorobenzidine	0.461	0.804	-74.4#	299#	0.00
37		Chrysene	2.026	2.069	-2.1	114	0.00
38		Bis(2-ethylhexyl)phthalate	4.801	8.435	-75.7#	237#	0.00
39		Indeno(1,2,3-cd)pyrene	5.859	7.655	-30.7#	160#	0.00
40	I	Perylene-d12	1.000	1.000	0.0	124	0.00
41		Benzo(b)fluoranthene	1.515	1.676	-10.6	143	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072716\
Data File : BE092070.D
Acq On : 28 Jul 2016 6:59
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 28 07:33:10 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE072716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 27 17:13:26 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)
42	Benzo(k)fluoranthene	1.567	1.699	-8.4	133	0.00
43 C	Benzo(a)pyrene	1.273	1.500	-17.8	151#	0.00
44	Dibenzo(a,h)anthracene	1.111	1.347	-21.2	163#	0.00
45	Benzo(g,h,i)perylene	4.638	5.533	-19.3	159#	0.00

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