Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE080116\
Data File : BE092096.D
Acq On : 1 Aug 2016 16:51
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE080116

Quant Time: Aug 01 20:08:10 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 01 17:22:55 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	113	0.00
2		1,4-Dioxane	0.400	0.396	1.0	124	0.00
3		n-Nitrosodimethylamine	0.400	0.418	-4.5	124	0.00
4	S	2-Fluorophenol	0.400	0.412	-3.0	113	0.00
5	S	Phenol-d6	0.400	0.494	-23.5	132	0.00
6		bis(2-Chloroethyl)ether	0.400	0.427	-6.7	128	0.00
7		n-Nitroso-di-n-propylamine	0.400	0.419	-4.7	126	0.00
8	I	Naphthalene-d8	5.000	5.000	0.0	113	0.00
9	S	Nitrobenzene-d5	0.400	0.478	-19.5	131	0.00
10		Nitrobenzene	0.400	0.425	-6.2	130	0.00
11		bis(2-Chloroethoxy)methane	0.400	0.482	-20.5	126	0.00
12		Naphthalene	0.400	0.457	-14.2	128	0.00
13		Hexachlorobutadiene	0.400	0.468	-17.0	132	0.00
14		2-Methylnaphthalene	0.400	0.461	-15.3	133	0.00
15	I	Acenaphthene-d10	5.000	5.000	0.0	120	0.00
16	S	2,4,6-Tribromophenol	0.400	0.404	-1.0	111	0.00
17	S	2-Fluorobiphenyl	0.400	0.448	-12.0	138	0.00
18		Acenaphthylene	0.400	0.463	-15.8	145	0.00
19		2,6-Dinitrotoluene	0.400	0.413	-3.2	131	0.00
20		Acenaphthene	0.400	0.463	-15.8	140	0.00
21		2,4-Dinitrotoluene	0.400	0.362	9.5	128	0.00
22		Fluorene	0.400	0.448	-12.0	137	0.00
23		4-Chlorophenyl-phenylether	0.400	0.435	-8.7	121	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	112	0.00
25		4-Bromophenyl-phenylether	0.400	0.469	-17.2	139	0.00
26		Hexachlorobenzene	0.400	0.487	-21.7	139	0.00
27		Pentachlorophenol	0.400	1.312	-228.0#	394	-0.02
28		Phenanthrene	0.400	0.468	-17.0	133	0.00
29		Anthracene	0.400	0.463	-15.8	135	0.00
30		Fluoranthene	0.400	0.486	-21.5	132	0.00
31	I	Chrysene-d12	5.000	5.000	0.0	116	0.00
32		Benzidine	0.400	0.557	-39.3#	166	0.00
33		Pyrene	0.400	0.433	-8.2	129	0.00
34	S	Terphenyl-d14	0.400	0.391	2.3	114	0.00
35		Benzo(a)anthracene	0.400	0.427	-6.7	135	0.00
36		3,3'-Dichlorobenzidine	0.400	0.423	-5.7	129	0.00
37		Chrysene	0.400	0.474	-18.5	128	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.358	10.5	114	0.00
39		Indeno(1,2,3-cd)pyrene	0.400	0.445	-11.2	132	0.00
40	I	Perylene-d12	5.000	5.000	0.0	111	0.00
41		Benzo(b)fluoranthene	0.400	0.453	-13.2	134	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE080116\
Data File : BE092096.D
Acq On : 1 Aug 2016 16:51
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE080116

Quant Time: Aug 01 20:08:10 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 01 17:22:55 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)
42	Benzo(k)fluoranthene	0.400	0.478	-19.5	128	0.00
43 C	Benzo(a)pyrene	0.400	0.461	-15.3	132	0.00
44	Dibenzo(a,h)anthracene	0.400	0.449	-12.2	130	0.00
45	Benzo(g,h,i)perylene	0.400	0.467	-16.8	132	-0.02

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