Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE011617\
Data File : BE092656.D
Acq On : 16 Jan 2017 17:42
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE011617

Quant Time: Jan 17 10:21:48 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE011617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 17 10:14:01 2017
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	108	0.00
2		1,4-Dioxane	0.400	0.401	-0.3	104	0.00
3		n-Nitrosodimethylamine	0.400	0.346	13.5	107	0.01
4	S	2-Fluorophenol	0.400	0.323	19.3	103	0.00
5	S	Phenol-d6	0.400	0.328	18.0	108	0.00
6		bis(2-Chloroethyl)ether	0.400	0.335	16.3	108	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	105	0.00
8	S	Nitrobenzene-d5	0.400	0.331	17.3	105	0.00
9		Naphthalene	0.400	0.383	4.3	109	0.00
10		Hexachlorobutadiene	0.400	0.383	4.3	106	0.00
11		2-Methylnaphthalene	0.400	0.358	10.5	106	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	108	0.00
13	S	2,4,6-Tribromophenol	0.400	0.413	-3.2	114	0.00
14	S	2-Fluorobiphenyl	0.400	0.370	7.5	108	0.00
15		Acenaphthylene	0.400	0.357	10.8	109	0.00
16		Acenaphthene	0.400	0.378	5.5	108	0.00
17		Fluorene	0.400	0.362	9.5	109	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	112	0.00
19		Hexachlorobenzene	0.400	0.350	12.5	101	0.00
20		Pentachlorophenol	0.400	0.339	15.3	104	0.00
21		Phenanthrene	0.400	0.388	3.0	112	0.00
22		Anthracene	0.400	0.363	9.3	114	0.00
23		Fluoranthene	0.400	0.362	9.5	111	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	109	0.00
25		Pyrene	0.400	0.339	15.3	111	0.00
26	S	Terphenyl-d14	0.400	0.337	15.8	107	0.00
27		Benzo(a)anthracene	0.400	0.339	15.3	108	0.00
28		Chrysene	0.400	0.381	4.8	112	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.418	-4.5	90	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.329	17.8	103	0.00
31	I	Perylene-d12	5.000	5.000	0.0	106	0.00
32		Benzo(b)fluoranthene	0.400	0.329	17.8	95	-0.01
33		Benzo(k)fluoranthene	0.400	0.388	3.0	120	-0.01
34	C	Benzo(a)pyrene	0.400	0.350	12.5	107	-0.01
35		Dibenzo(a,h)anthracene	0.400	0.340	15.0	102	-0.02
36		Benzo(g,h,i)perylene	0.400	0.346	13.5	102	0.00

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