Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE040417\
Data File : BE092886.D
Acq On : 4 Apr 2017 15:58
Operator : SJ/MA
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE040417

Quant Time: Apr 04 16:37:50 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE040417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 04 16:34:01 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	107	-0.02
2		1,4-Dioxane	0.400	0.408	-2.0	109	0.00
3		n-Nitrosodimethylamine	0.400	0.394	1.5	112	0.00
4	S	2-Fluorophenol	0.400	0.400	0.0	117	0.00
5	S	Phenol-d6	0.400	0.386	3.5	117	0.00
6		bis(2-Chloroethyl)ether	0.400	0.391	2.3	107	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	109	0.00
8	S	Nitrobenzene-d5	0.400	0.416	-4.0	126	-0.02
9		Naphthalene	0.400	0.401	-0.3	109	0.00
10		Hexachlorobutadiene	0.400	0.397	0.8	108	0.00
11		2-Methylnaphthalene	0.400	0.393	1.8	110	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	113	-0.02
13	S	2,4,6-Tribromophenol	0.400	0.369	7.8	123	-0.01
14	S	2-Fluorobiphenyl	0.400	0.394	1.5	109	-0.01
15		Acenaphthylene	0.400	0.369	7.8	117	-0.02
16		Acenaphthene	0.400	0.385	3.8	112	0.00
17		Fluorene	0.400	0.378	5.5	112	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	113	-0.02
19		Hexachlorobenzene	0.400	0.383	4.3	109	0.00
20		Pentachlorophenol	0.400	0.430	-7.5	116	-0.02
21		Phenanthrene	0.400	0.389	2.8	113	0.00
22		Anthracene	0.400	0.357	10.8	114	0.00
23		Fluoranthene	0.400	0.368	8.0	111	-0.02
24	I	Chrysene-d12	5.000	5.000	0.0	110	-0.02
25		Pyrene	0.400	0.380	5.0	113	0.00
26	S	Terphenyl-d14	0.400	0.385	3.8	111	-0.02
27		Benzo(a)anthracene	0.400	0.352	12.0	114	0.00
28		Chrysene	0.400	0.410	-2.5	108	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.442	-10.5	121	-0.02
30		Indeno(1,2,3-cd)pyrene	0.400	0.356	11.0	101	-0.02
31	I	Perylene-d12	5.000	5.000	0.0	105	-0.01
32		Benzo(b)fluoranthene	0.400	0.369	7.8	104	-0.01
33		Benzo(k)fluoranthene	0.400	0.362	9.5	107	-0.01
34	C	Benzo(a)pyrene	0.400	0.346	13.5	104	-0.01
35		Dibenzo(a,h)anthracene	0.400	0.356	11.0	102	-0.03
36		Benzo(g,h,i)perylene	0.400	0.367	8.3	102	-0.02

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