Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE090116\
Data File : BE092333.D
Acq On : 1 Sep 2016 17:47
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE090116

Quant Time: Sep 02 10:55:34 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE090116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 02 03:36:16 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	92	0.00
2		1,4-Dioxane	0.400	0.451	-12.7	106	0.00
3		n-Nitrosodimethylamine	0.400	0.384	4.0	94	0.00
4	S	2-Fluorophenol	0.400	0.387	3.3	98	0.00
5	S	Phenol-d6	0.400	0.336	16.0	98	0.00
6		bis(2-Chloroethyl)ether	0.400	0.440	-10.0	101	0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	91	0.00
8	S	Nitrobenzene-d5	0.400	0.328	18.0	97	0.00
9		Naphthalene	0.400	0.424	-6.0	104	0.00
10		Hexachlorobutadiene	0.400	0.440	-10.0	108	0.00
11		2-Methylnaphthalene	0.400	0.412	-3.0	103	0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	89	-0.02
13	S	2,4,6-Tribromophenol	0.400	0.350	12.5	91	0.00
14	S	2-Fluorobiphenyl	0.400	0.410	-2.5	99	0.00
15		Acenaphthylene	0.400	0.422	-5.5	103	0.00
16		Acenaphthene	0.400	0.431	-7.7	102	0.00
17		Fluorene	0.400	0.428	-7.0	102	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	91	0.00
19		Hexachlorobenzene	0.400	0.429	-7.2	104	0.00
20		Pentachlorophenol	0.400	0.374	6.5	89	0.00
21		Phenanthrene	0.400	0.422	-5.5	100	0.00
22		Anthracene	0.400	0.405	-1.3	100	0.00
23		Fluoranthene	0.400	0.414	-3.5	98	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	92	0.00
25		Pyrene	0.400	0.400	0.0	97	0.00
26	S	Terphenyl-d14	0.400	0.388	3.0	95	0.00
27		Benzo(a)anthracene	0.400	0.418	-4.5	102	0.00
28		Chrysene	0.400	0.447	-11.7	102	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.361	9.8	87	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.420	-5.0	100	0.00
31	I	Perylene-d12	5.000	5.000	0.0	89	0.00
32		Benzo(b)fluoranthene	0.400	0.393	1.8	103	0.00
33		Benzo(k)fluoranthene	0.400	0.444	-11.0	100	-0.01
34	C	Benzo(a)pyrene	0.400	0.415	-3.7	101	0.00
35		Dibenzo(a,h)anthracene	0.400	0.408	-2.0	100	-0.02
36		Benzo(g,h,i)perylene	0.400	0.416	-4.0	100	0.00

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