Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102516\
Data File : BE092479.D
Acq On : 25 Oct 2016 14:43
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE102516

Quant Time: Oct 25 15:30:55 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE102516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 25 15:25:33 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	89	0.00
2		1,4-Dioxane	0.400	0.370	7.5	90	0.00
3		n-Nitrosodimethylamine	0.400	0.446	-11.5	82	0.00
4	S	2-Fluorophenol	0.400	0.425	-6.2	103	0.00
5	S	Phenol-d6	0.400	0.418	-4.5	100	0.00
6		bis(2-Chloroethyl)ether	0.400	0.430	-7.5	89	0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	94	0.00
8	S	Nitrobenzene-d5	0.400	0.469	-17.2	99	0.00
9		Naphthalene	0.400	0.381	4.8	89	0.00
10		Hexachlorobutadiene	0.400	0.392	2.0	91	0.00
11		2-Methylnaphthalene	0.400	0.382	4.5	93	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	95	0.00
13	S	2,4,6-Tribromophenol	0.400	0.405	-1.3	100	0.00
14	S	2-Fluorobiphenyl	0.400	0.416	-4.0	101	0.00
15		Acenaphthylene	0.400	0.396	1.0	98	0.00
16		Acenaphthene	0.400	0.406	-1.5	98	0.00
17		Fluorene	0.400	0.403	-0.8	100	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	95	0.00
19		Hexachlorobenzene	0.400	0.410	-2.5	101	0.00
20		Pentachlorophenol	0.400	0.431	-7.7	83	0.00
21		Phenanthrene	0.400	0.383	4.3	100	0.00
22		Anthracene	0.400	0.369	7.8	100	0.00
23		Fluoranthene	0.400	0.390	2.5	99	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	94	0.00
25		Pyrene	0.400	0.386	3.5	97	-0.02
26	S	Terphenyl-d14	0.400	0.397	0.8	101	0.00
27		Benzo(a)anthracene	0.400	0.373	6.8	93	0.00
28		Chrysene	0.400	0.419	-4.7	102	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.448	-12.0	103	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.376	6.0	96	0.00
31	I	Perylene-d12	5.000	5.000	0.0	95	0.00
32		Benzo(b)fluoranthene	0.400	0.341	14.8	94	0.00
33		Benzo(k)fluoranthene	0.400	0.406	-1.5	94	0.00
34	C	Benzo(a)pyrene	0.400	0.347	13.3	94	0.00
35		Dibenzo(a,h)anthracene	0.400	0.368	8.0	96	0.00
36		Benzo(g,h,i)perylene	0.400	0.378	5.5	95	0.00

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