Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE092216\
Data File : BE092400.D
Acq On : 22 Sep 2016 17:49
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE092216

Quant Time: Sep 22 18:39:58 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE092216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 22 18:25:56 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	103	-0.02
2		1,4-Dioxane	0.400	0.347	13.3	97	-0.01
3		n-Nitrosodimethylamine	0.400	0.440	-10.0	99	-0.01
4	S	2-Fluorophenol	0.400	0.391	2.3	92	0.00
5	S	Phenol-d6	0.400	0.331	17.3	89	0.00
6		bis(2-Chloroethyl)ether	0.400	0.340	15.0	91	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	100	0.00
8	S	Nitrobenzene-d5	0.400	0.401	-0.3	94	0.00
9		Naphthalene	0.400	0.411	-2.7	103	0.00
10		Hexachlorobutadiene	0.400	0.402	-0.5	101	0.00
11		2-Methylnaphthalene	0.400	0.400	0.0	101	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	103	0.00
13	S	2,4,6-Tribromophenol	0.400	0.413	-3.2	95	0.01
14	S	2-Fluorobiphenyl	0.400	0.374	6.5	97	0.01
15		Acenaphthylene	0.400	0.380	5.0	100	0.02
16		Acenaphthene	0.400	0.408	-2.0	101	0.00
17		Fluorene	0.400	0.383	4.3	100	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	100	0.02
19		Hexachlorobenzene	0.400	0.367	8.3	99	0.00
20		Pentachlorophenol	0.400	0.380	5.0	97	0.00
21		Phenanthrene	0.400	0.377	5.8	100	0.02
22		Anthracene	0.400	0.361	9.8	100	0.02
23		Fluoranthene	0.400	0.370	7.5	101	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	101	0.00
25		Pyrene	0.400	0.379	5.3	99	0.02
26	S	Terphenyl-d14	0.400	0.376	6.0	96	0.02
27		Benzo(a)anthracene	0.400	0.379	5.3	97	0.00
28		Chrysene	0.400	0.417	-4.2	98	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.322	19.5	80	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.364	9.0	93	0.05
31	I	Perylene-d12	5.000	5.000	0.0	98	0.03
32		Benzo(b)fluoranthene	0.400	0.399	0.3	98	0.03
33		Benzo(k)fluoranthene	0.400	0.372	7.0	98	0.03
34	C	Benzo(a)pyrene	0.400	0.379	5.3	98	0.02
35		Dibenzo(a,h)anthracene	0.400	0.375	6.3	96	0.05
36		Benzo(g,h,i)perylene	0.400	0.387	3.3	98	0.05

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