Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE080916\
Data File : BE092243.D
Acq On : 10 Aug 2016 00:16
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE080916

Quant Time: Aug 10 11:32:18 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 10 11:12:32 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	107	0.00
2		1,4-Dioxane	0.400	0.368	8.0	98	0.00
3		n-Nitrosodimethylamine	0.400	0.420	-5.0	98	0.00
4	S	2-Fluorophenol	0.400	0.333	16.8	99	0.00
5	S	Phenol-d6	0.400	0.329	17.8	113	0.00
6		bis(2-Chloroethyl)ether	0.400	0.328	18.0	98	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	108	0.00
8	S	Nitrobenzene-d5	0.400	0.334	16.5	107	-0.02
9		Naphthalene	0.400	0.368	8.0	100	0.00
10		Hexachlorobutadiene	0.400	0.376	6.0	102	0.00
11		2-Methylnaphthalene	0.400	0.366	8.5	101	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	107	0.00
13	S	2,4,6-Tribromophenol	0.400	0.453	-13.2	98	0.00
14	S	2-Fluorobiphenyl	0.400	0.361	9.8	99	0.00
15		Acenaphthylene	0.400	0.360	10.0	100	0.00
16		Acenaphthene	0.400	0.371	7.3	101	0.00
17		Fluorene	0.400	0.363	9.3	99	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	107	0.00
19		Hexachlorobenzene	0.400	0.380	5.0	101	0.00
20		Pentachlorophenol	0.400	0.360	10.0	98	0.02
21		Phenanthrene	0.400	0.361	9.8	100	0.00
22		Anthracene	0.400	0.338	15.5	97	0.00
23		Fluoranthene	0.400	0.348	13.0	98	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	111	-0.02
25		Pyrene	0.400	0.346	13.5	100	0.00
26	S	Terphenyl-d14	0.400	0.343	14.2	102	0.00
27		Benzo(a)anthracene	0.400	0.380	5.0	113	0.00
28		Chrysene	0.400	0.345	13.8	94	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.417	-4.2	84	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.324	19.0	99	-0.02
31	I	Perylene-d12	5.000	5.000	0.0	109	0.00
32		Benzo(b)fluoranthene	0.400	0.339	15.3	100	0.00
33		Benzo(k)fluoranthene	0.400	0.376	6.0	101	0.00
34	C	Benzo(a)pyrene	0.400	0.340	15.0	101	0.00
35		Dibenzo(a,h)anthracene	0.400	0.331	17.3	100	0.00
36		Benzo(g,h,i)perylene	0.400	0.349	12.8	100	0.00

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