Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\Be072216\
Data File : BE092007.D
Acq On : 22 Jul 2016 15:29
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE072216

Quant Time: Jul 22 16:31:32 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE072216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 22 15:39:27 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	94	0.00
2		1,4-Dioxane	0.400	0.432	-8.0	96	-0.02
3		n-Nitrosodimethylamine	0.400	0.419	-4.7	108	-0.02
4	S	2-Fluorophenol	0.400	0.427	-6.7	110	0.00
5	S	Phenol-d6	0.400	0.435	-8.7	108	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.411	-2.7	105	-0.02
7		n-Nitroso-di-n-propylamine	0.400	0.385	3.8	100	0.00
8	I	Naphthalene-d8	5.000	5.000	0.0	95	0.00
9	S	Nitrobenzene-d5	0.400	0.409	-2.2	101	-0.02
10		Nitrobenzene	0.400	0.439	-9.7	100	-0.02
11		bis(2-Chloroethoxy)methane	0.400	0.393	1.8	101	-0.03
12		Naphthalene	0.400	0.398	0.5	92	0.00
13		Hexachlorobutadiene	0.400	0.403	-0.8	92	0.00
14		2-Methylnaphthalene	0.400	0.399	0.3	95	0.00
15	I	Acenaphthene-d10	5.000	5.000	0.0	81	0.00
16	S	2,4,6-Tribromophenol	0.400	0.460	-15.0	107	0.00
17	S	2-Fluorobiphenyl	0.400	0.449	-12.2	95	0.00
18		Acenaphthylene	0.400	0.455	-13.7	113	0.00
19		2,6-Dinitrotoluene	0.400	0.430	-7.5	125	0.00
20		Acenaphthene	0.400	0.388	3.0	79	0.00
21		2,4-Dinitrotoluene	0.400	0.454	-13.5	105	0.00
22		Fluorene	0.400	0.451	-12.7	99	0.00
23		4-Chlorophenyl-phenylether	0.400	0.466	-16.5	98	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	99	0.00
25		4-Bromophenyl-phenylether	0.400	0.404	-1.0	100	0.00
26		Hexachlorobenzene	0.400	0.406	-1.5	97	0.00
27		Pentachlorophenol	0.400	0.460	-15.0	115	0.00
28		Phenanthrene	0.400	0.409	-2.2	99	-0.01
29		Anthracene	0.400	0.404	-1.0	104	-0.01
30		Fluoranthene	0.400	0.385	3.8	100	0.00
31	I	Chrysene-d12	5.000	5.000	0.0	87	0.00
32		Benzidine	0.400	0.468	-17.0	125	0.00
33		Pyrene	0.400	0.408	-2.0	95	0.00
34	S	Terphenyl-d14	0.400	0.451	-12.7	102	0.00
35		Benzo(a)anthracene	0.400	0.440	-10.0	107	0.00
36		3,3'-Dichlorobenzidine	0.400	0.462	-15.5	112	0.00
37		Chrysene	0.400	0.458	-14.5	100	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.442	-10.5	119	0.00
39		Indeno(1,2,3-cd)pyrene	0.400	0.395	1.3	87	0.00
40	I	Perylene-d12	5.000	5.000	0.0	93	0.00
41		Benzo(b)fluoranthene	0.400	0.384	4.0	94	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.390	2.5	90	0.00
43 C	Benzo(a)pyrene	0.400	0.373	6.8	90	0.00
44	Dibenzo(a,h)anthracene	0.400	0.370	7.5	88	0.00
45	Benzo(g,h,i)perylene	0.400	0.385	3.8	92	0.00

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