Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE060716\
Data File : BE091884.D
Acq On : 7 Jun 2016 16:33
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE060716

Quant Time: Jun 07 19:38:48 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 07 17:14:38 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	100	0.00
2		1,4-Dioxane	0.781	0.847	-8.5	103	-0.02
3		n-Nitrosodimethylamine	0.836	0.872	-4.3	100	0.00
4	S	2-Fluorophenol	1.066	1.094	-2.6	103	0.00
5	S	Phenol-d6	1.320	1.363	-3.3	107	-0.02
6		bis(2-Chloroethyl)ether	1.431	1.425	0.4	100	0.00
7		n-Nitroso-di-n-propylamine	1.130	1.081	4.3	98	0.00
8	I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
9	S	Nitrobenzene-d5	0.336	0.324	3.6	93	-0.02
10		Nitrobenzene	0.304	0.308	-1.3	103	0.00
11		bis(2-Chloroethoxy)methane	0.507	0.551	-8.7	115	0.00
12		Naphthalene	1.079	1.153	-6.9	101	0.00
13		Hexachlorobutadiene	0.161	0.173	-7.5	101	0.00
14		2-Methylnaphthalene	0.696	0.729	-4.7	102	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
16	S	2,4,6-Tribromophenol	0.142	0.141	0.7	109	-0.02
17	S	2-Fluorobiphenyl	1.517	1.609	-6.1	100	0.00
18		Acenaphthylene	2.263	2.195	3.0	98	0.00
19		2,6-Dinitrotoluene	0.240	0.221	7.9	103	-0.02
20		Acenaphthene	1.383	1.465	-5.9	104	0.00
21		2,4-Dinitrotoluene	0.285	0.241	15.4	105	-0.02
22		Fluorene	1.691	1.734	-2.5	102	-0.02
23		4-Chlorophenyl-phenylether	0.829	0.884	-6.6	102	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	104	0.00
25		4-Bromophenyl-phenylether	0.180	0.170	5.6	105	0.00
26		Hexachlorobenzene	0.184	0.176	4.3	102	0.00
27		Pentachlorophenol	0.070	0.054	22.9	114	0.00
28		Phenanthrene	1.158	1.112	4.0	103	0.00
29		Anthracene	1.032	0.955	7.5	106	0.00
30		Fluoranthene	1.250	1.137	9.0	101	0.00
31	I	Chrysene-d12	1.000	1.000	0.0	103	0.00
32		Benzidine	0.333	0.144	56.8#	49#	0.00
33		Pyrene	1.262	1.526	-20.9	104	0.00
34	S	Terphenyl-d14	0.891	1.090	-22.3	103	0.00
35		Benzo(a)anthracene	7.222	8.429	-16.7	102	0.00
36		3,3'-Dichlorobenzidine	0.349	0.299	14.3	97	0.00
37		Chrysene	5.098	5.275	-3.5	105	0.00
38		Bis(2-ethylhexyl)phthalate	0.716	0.655	8.5	113	0.00
39		Indeno(1,2,3-cd)pyrene	1.524	1.836	-20.5	102	0.00
40	I	Perylene-d12	1.000	1.000	0.0	95	0.00
41		Benzo(b)fluoranthene	1.222	1.270	-3.9	98	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.268	1.273	-0.4	97	0.00
43 C	Benzo(a)pyrene	1.146	1.152	-0.5	99	0.00
44	Dibenzo(a,h)anthracene	1.068	1.118	-4.7	102	-0.02
45	Benzo(g,h,i)perylene	1.120	1.180	-5.4	99	-0.02

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