Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE052516\
Data File : BE091869.D
Acq On : 25 May 2016 21:44
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 26 13:43:32 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE052516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 25 19:42:30 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	115	0.00
2		1,4-Dioxane	0.400	0.419	-4.7	109	0.00
3		n-Nitrosodimethylamine	0.400	0.341	14.8	112	0.00
4	S	2-Fluorophenol	0.400	0.357	10.8	112	0.00
5	S	Phenol-d6	0.400	0.370	7.5	121	0.00
6		bis(2-Chloroethyl)ether	0.400	0.372	7.0	117	-0.02
7		n-Nitroso-di-n-propylamine	0.400	0.354	11.5	119	0.00
8	I	Naphthalene-d8	5.000	5.000	0.0	126	0.00
9	S	Nitrobenzene-d5	0.400	0.357	10.8	123	0.00
10		Nitrobenzene	0.400	0.365	8.8	131	0.00
11		bis(2-Chloroethoxy)methane	0.400	0.360	10.0	143	0.00
12		Naphthalene	0.400	0.387	3.3	129	0.00
13		Hexachlorobutadiene	0.400	0.380	5.0	123	0.00
14		2-Methylnaphthalene	0.400	0.388	3.0	132	0.00
15	I	Acenaphthene-d10	5.000	5.000	0.0	130	0.00
16	S	2,4,6-Tribromophenol	0.400	0.319	20.3	117	0.00
17	S	2-Fluorobiphenyl	0.400	0.386	3.5	130	0.00
18		Acenaphthylene	0.400	0.387	3.3	132	0.00
19		2,6-Dinitrotoluene	0.400	0.325	18.8	126	-0.02
20		Acenaphthene	0.400	0.399	0.3	133	0.00
21		2,4-Dinitrotoluene	0.400	0.441	-10.2	126	0.00
22		Fluorene	0.400	0.378	5.5	131	0.00
23		4-Chlorophenyl-phenylether	0.400	0.344	14.0	125	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	119	-0.01
25		4-Bromophenyl-phenylether	0.400	0.384	4.0	123	0.00
26		Hexachlorobenzene	0.400	0.406	-1.5	126	0.00
27		Pentachlorophenol	0.400	0.361	9.8	139	-0.02
28		Phenanthrene	0.400	0.367	8.3	122	0.00
29		Anthracene	0.400	0.380	5.0	124	0.00
30		Fluoranthene	0.400	0.334	16.5	114	-0.02
31	I	Chrysene-d12	5.000	5.000	0.0	99	0.00
32		Benzidine	0.400	0.448	-12.0	121	0.00
33		Pyrene	0.400	0.431	-7.7	116	0.00
34	S	Terphenyl-d14	0.400	0.411	-2.7	108	0.00
35		Benzo(a)anthracene	0.400	0.386	3.5	101	0.00
36		3,3'-Dichlorobenzidine	0.400	0.383	4.3	105	0.00
37		Chrysene	0.400	0.323	19.3	101	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.398	0.5	116	0.00
39		Indeno(1,2,3-cd)pyrene	0.400	0.390	2.5	104	0.00
40	I	Perylene-d12	5.000	5.000	0.0	101	0.00
41		Benzo(b)fluoranthene	0.400	0.354	11.5	102	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.412	-3.0	102	-0.01
43 C	Benzo(a)pyrene	0.400	0.379	5.3	103	0.00
44	Dibenzo(a,h)anthracene	0.400	0.381	4.8	103	0.00
45	Benzo(g,h,i)perylene	0.400	0.385	3.8	103	0.00

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