Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE061516\
Data File : BE091933.D
Acq On : 15 Jun 2016 20:14
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 16 01:00:28 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE061416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 15 12:00:38 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	82	0.02
2		1,4-Dioxane	0.400	0.400	0.0	80	0.00
3		n-Nitrosodimethylamine	0.400	0.316	21.0	64	0.04
4	S	2-Fluorophenol	0.400	0.349	12.8	72	0.02
5	S	Phenol-d6	0.400	0.404	-1.0	74	0.02
6		bis(2-Chloroethyl)ether	0.400	0.344	14.0	71	0.02
7		n-Nitroso-di-n-propylamine	0.400	0.411	-2.7	73	0.02
8	I	Naphthalene-d8	5.000	5.000	0.0	84	0.03
9	S	Nitrobenzene-d5	0.400	0.364	9.0	76	0.02
10		Nitrobenzene	0.400	0.355	11.3	72	0.02
11		bis(2-Chloroethoxy)methane	0.400	0.340	15.0	66	0.03
12		Naphthalene	0.400	0.386	3.5	76	0.00
13		Hexachlorobutadiene	0.400	0.384	4.0	75	0.00
14		2-Methylnaphthalene	0.400	0.368	8.0	75	0.01
15	I	Acenaphthene-d10	5.000	5.000	0.0	79	0.00
16	S	2,4,6-Tribromophenol	0.400	0.331	17.3	68	0.02
17	S	2-Fluorobiphenyl	0.400	0.402	-0.5	74	0.00
18		Acenaphthylene	0.400	0.358	10.5	71	0.03
19		2,6-Dinitrotoluene	0.400	0.420	-5.0	69	0.00
20		Acenaphthene	0.400	0.373	6.8	74	0.00
21		2,4-Dinitrotoluene	0.400	0.387	3.3	74	0.03
22		Fluorene	0.400	0.365	8.8	72	0.02
23		4-Chlorophenyl-phenylether	0.400	0.371	7.3	70	0.02
24	I	Phenanthrene-d10	5.000	5.000	0.0	78	0.00
25		4-Bromophenyl-phenylether	0.400	0.365	8.8	70	0.02
26		Hexachlorobenzene	0.400	0.391	2.3	73	0.00
27		Pentachlorophenol	0.400	0.295	26.3#	55	0.02
28		Phenanthrene	0.400	0.386	3.5	72	0.00
29		Anthracene	0.400	0.358	10.5	71	0.00
30		Fluoranthene	0.400	0.335	16.3	64	0.00
31	I	Chrysene-d12	5.000	5.000	0.0	79	0.00
32		Benzidine	0.400	0.316	21.0	70	0.02
33		Pyrene	0.400	0.333	16.8	67	0.00
34	S	Terphenyl-d14	0.400	0.319	20.3	61	0.00
35		Benzo(a)anthracene	0.400	0.395	1.3	80	0.00
36		3,3'-Dichlorobenzidine	0.400	0.282	29.5#	67	0.03
37		Chrysene	0.400	0.413	-3.2	74	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.339	15.3	71	-0.03
39		Indeno(1,2,3-cd)pyrene	0.400	0.327	18.3	66	0.00
40	I	Perylene-d12	5.000	5.000	0.0	75	-0.01
41		Benzo(b)fluoranthene	0.400	0.382	4.5	69	-0.01

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.341	14.8	64	0.00
43 C	Benzo(a)pyrene	0.400	0.347	13.3	66	0.00
44	Dibenzo(a,h)anthracene	0.400	0.349	12.8	65	-0.02
45	Benzo(g,h,i)perylene	0.400	0.351	12.3	65	-0.02

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