Evaluate Continuing Calibration Report

Data Path : V:\HPCHEM1\BNA E\DATA\BE070916\
Data File : BE091996.D
Acq On : 9 Jul 2016 13:27
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 11 04:51:47 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE070516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 05 16:14:58 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	97	0.00
2		1,4-Dioxane	0.400	0.391	2.3	91	0.00
3		n-Nitrosodimethylamine	0.400	0.349	12.8	93	-0.02
4	S	2-Fluorophenol	0.400	0.349	12.8	89	0.00
5	S	Phenol-d6	0.400	0.336	16.0	90	0.00
6		bis(2-Chloroethyl)ether	0.400	0.378	5.5	98	-0.02
7		n-Nitroso-di-n-propylamine	0.400	0.332	17.0	86	-0.02
8	I	Naphthalene-d8	5.000	5.000	0.0	99	0.00
9	S	Nitrobenzene-d5	0.400	0.365	8.8	92	-0.02
10		Nitrobenzene	0.400	0.351	12.3	93	0.00
11		bis(2-Chloroethoxy)methane	0.400	0.410	-2.5	88	0.00
12		Naphthalene	0.400	0.388	3.0	94	0.00
13		Hexachlorobutadiene	0.400	0.373	6.8	86	0.00
14		2-Methylnaphthalene	0.400	0.376	6.0	91	-0.01
15	I	Acenaphthene-d10	5.000	5.000	0.0	90	0.00
16	S	2,4,6-Tribromophenol	0.400	0.426	-6.5	96	0.00
17	S	2-Fluorobiphenyl	0.400	0.425	-6.2	93	0.00
18		Acenaphthylene	0.400	0.405	-1.3	94	0.00
19		2,6-Dinitrotoluene	0.400	0.393	1.8	112	0.00
20		Acenaphthene	0.400	0.380	5.0	84	0.00
21		2,4-Dinitrotoluene	0.400	0.539	-34.8#	165	0.00
22		Fluorene	0.400	0.408	-2.0	90	0.00
23		4-Chlorophenyl-phenylether	0.400	0.410	-2.5	88	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	101	0.00
25		4-Bromophenyl-phenylether	0.400	0.356	11.0	88	0.00
26		Hexachlorobenzene	0.400	0.377	5.8	94	0.00
27		Pentachlorophenol	0.400	0.361	9.8	105	0.00
28		Phenanthrene	0.400	0.384	4.0	93	0.00
29		Anthracene	0.400	0.364	9.0	93	0.00
30		Fluoranthene	0.400	0.372	7.0	94	0.00
31	I	Chrysene-d12	5.000	5.000	0.0	120	0.00
32		Benzidine	0.400	0.360	10.0	63	0.00
33		Pyrene	0.400	0.326	18.5	97	0.00
34	S	Terphenyl-d14	0.400	0.358	10.5	106	0.00
35		Benzo(a)anthracene	0.400	0.399	0.3	121	0.00
36		3,3'-Dichlorobenzidine	0.400	0.213	46.8#	70	0.00
37		Chrysene	0.400	0.428	-7.0	129	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.254	36.5#	83	-0.03
39		Indeno(1,2,3-cd)pyrene	0.400	0.329	17.8	96	-0.02
40	I	Perylene-d12	5.000	5.000	0.0	110	0.00
41		Benzo(b)fluoranthene	0.400	0.396	1.0	110	-0.01

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.346	13.5	96	0.00
43 C	Benzo(a)pyrene	0.400	0.358	10.5	99	0.00
44	Dibenzo(a,h)anthracene	0.400	0.341	14.8	96	0.00
45	Benzo(g,h,i)perylene	0.400	0.343	14.2	95	0.00

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