Evaluate Continuing Calibration Report

Data Path : V:\HPCHEM1\BNA E\DATA\BE070916\
Data File : BE091991.D
Acq On : 9 Jul 2016 10:19
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 11 04:38:01 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	105	0.00
2		1,4-Dioxane	0.400	0.385	3.8	97	0.00
3		n-Nitrosodimethylamine	0.400	0.356	11.0	103	0.00
4	S	2-Fluorophenol	0.400	0.326	18.5	91	0.00
5	S	Phenol-d6	0.400	0.327	18.3	95	0.00
6		bis(2-Chloroethyl)ether	0.400	0.386	3.5	109	-0.02
7		n-Nitroso-di-n-propylamine	0.400	0.333	16.8	94	-0.02
8	I	Naphthalene-d8	5.000	5.000	0.0	107	0.00
9	S	Nitrobenzene-d5	0.400	0.350	12.5	95	-0.02
10		Nitrobenzene	0.400	0.347	13.3	99	0.00
11		bis(2-Chloroethoxy)methane	0.400	0.361	9.8	84	-0.03
12		Naphthalene	0.400	0.389	2.8	101	0.00
13		Hexachlorobutadiene	0.400	0.380	5.0	95	0.00
14		2-Methylnaphthalene	0.400	0.374	6.5	98	-0.01
15	I	Acenaphthene-d10	5.000	5.000	0.0	98	0.00
16	S	2,4,6-Tribromophenol	0.400	0.385	3.8	90	0.00
17	S	2-Fluorobiphenyl	0.400	0.407	-1.7	97	0.00
18		Acenaphthylene	0.400	0.399	0.3	101	0.00
19		2,6-Dinitrotoluene	0.400	0.390	2.5	120	0.00
20		Acenaphthene	0.400	0.372	7.0	89	0.00
21		2,4-Dinitrotoluene	0.400	0.519	-29.8#	167	0.00
22		Fluorene	0.400	0.398	0.5	95	0.00
23		4-Chlorophenyl-phenylether	0.400	0.404	-1.0	94	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	105	0.00
25		4-Bromophenyl-phenylether	0.400	0.360	10.0	93	0.00
26		Hexachlorobenzene	0.400	0.394	1.5	102	0.00
27		Pentachlorophenol	0.400	0.449	-12.2	150	0.00
28		Phenanthrene	0.400	0.391	2.3	99	0.00
29		Anthracene	0.400	0.368	8.0	98	0.00
30		Fluoranthene	0.400	0.374	6.5	99	0.00
31	I	Chrysene-d12	5.000	5.000	0.0	122	0.00
32		Benzidine	0.400	0.464	-16.0	110	0.00
33		Pyrene	0.400	0.330	17.5	100	0.00
34	S	Terphenyl-d14	0.400	0.351	12.3	106	0.00
35		Benzo(a)anthracene	0.400	0.368	8.0	113	0.00
36		3,3'-Dichlorobenzidine	0.400	0.222	44.5#	73	0.00
37		Chrysene	0.400	0.399	0.3	122	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.242	39.5#	80	-0.03
39		Indeno(1,2,3-cd)pyrene	0.400	0.317	20.8	94	-0.02
40	I	Perylene-d12	5.000	5.000	0.0	107	0.00
41		Benzo(b)fluoranthene	0.400	0.386	3.5	105	0.00

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Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.353	11.8	96	0.00
43 C	Benzo(a)pyrene	0.400	0.355	11.3	96	0.00
44	Dibenzo(a,h)anthracene	0.400	0.347	13.3	95	0.00
45	Benzo(g,h,i)perylene	0.400	0.344	14.0	93	0.00

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