Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061217\
Data File : BE093106.D
Acq On : 12 Jun 2017 17:53
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 13 01:10:52 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE061217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 12 14:07:54 2017
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	0.400	0.400	0.0	83	0.00
2		1,4-Dioxane	0.400	0.422	-5.5	104	0.00
3		n-Nitrosodimethylamine	0.400	0.421	-5.2	96	0.00
4	S	2-Fluorophenol	0.400	0.411	-2.7	100	0.00
5	S	Phenol-d6	0.400	0.402	-0.5	95	0.00
6		bis(2-Chloroethyl)ether	0.400	0.426	-6.5	99	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	83	0.00
8	S	Nitrobenzene-d5	0.400	0.417	-4.2	100	0.00
9		Naphthalene	0.400	0.428	-7.0	100	0.00
10		Hexachlorobutadiene	0.400	0.426	-6.5	102	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.431	-7.7	104	0.00
12		2-Methylnaphthalene	0.400	0.420	-5.0	99	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	87	0.00
14	S	2,4,6-Tribromophenol	0.400	0.411	-2.7	107	0.00
15	S	2-Fluorobiphenyl	0.400	0.430	-7.5	104	0.00
16		Acenaphthylene	0.400	0.394	1.5	102	0.00
17		Acenaphthene	0.400	0.423	-5.7	105	0.00
18		Fluorene	0.400	0.427	-6.7	106	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	77	0.00
20		4-Bromophenyl-phenylether	0.400	0.416	-4.0	93	0.00
21		Hexachlorobenzene	0.400	0.508	-27.0#	127	0.00
22		Pentachlorophenol	0.400	0.395	1.3	96	0.00
23		Phenanthrene	0.400	0.427	-6.7	97	0.00
24		Anthracene	0.400	0.417	-4.2	103	0.00
25	SURR	Fluoranthene-d10	0.400	0.428	-7.0	98	0.00
26		Fluoranthene	0.400	0.428	-7.0	101	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	72	0.00
28		Pyrene	0.400	0.464	-16.0	91	0.00
29	S	Terphenyl-d14	0.400	0.449	-12.2	89	0.00
30		Benzo(a)anthracene	0.400	0.396	1.0	80	0.00
31		Chrysene	0.400	0.443	-10.7	89	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.304	24.0	59	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.372	7.0	76	0.00
34	I	Perylene-d12	0.400	0.400	0.0	63	0.00
35		Benzo(b)fluoranthene	0.400	0.453	-13.2	84	0.00
36		Benzo(k)fluoranthene	0.400	0.419	-4.7	77	0.00
37	C	Benzo(a)pyrene	0.400	0.415	-3.7	77	0.00
38		Dibenzo(a,h)anthracene	0.400	0.415	-3.7	77	0.00
39		Benzo(g,h,i)perylene	0.400	0.424	-6.0	79	0.00

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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