Zeaxanthin in Soybean oil: impact of oxidative stability, degradation pattern and product analysis

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Summary of Figures and tables in Supporting Information:

Table S1. Pearson correlation and regression models coefficients relating the stability indexes, Y=peroxide value (mmol/kg), to the zeaxanthin content, X (μ g/g), in soybean oil.

Table S2. Reaction rate constant (k) and correlation of coefficient (R^2) of three kinetic models of all-E-zeaxanthin in soybean oil at isothermal condition (110°C).

 Table S3. Zeaxanthin-thermodegraded compounds that tentatively identified in soybean oil.

Figure S1. Possible routes for the formation of 6-methyl-3,5-heptadien-2-one, 3-hydroxy-5,6-epoxy-7,8-dihydro- β -ionone, 3-hydroxy- β -ionone and 3-hydroxy- β -cy-clocitral, from the Zeaxanthin.

Table S	1.
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Sample	10 (µg/g)	50 (µg/g)	100 (μg/g)	200 (µg/g)	400 (µg/g)
Pearson					
correlation	-0.954	-0.981	-0.988	-0.960	-0.995
coefficient					
p-value	0.003	0.012	0.024	0.005	0.028
Democrien medal	Y = -4.3886*X	Y= -0.9936*X +	Y = -0.5726*X +	Y= -0.2368*X +	Y= -0.1137*X +
Regression model	+ 47.315	43.217	56.95	50.543	48.923
\mathbb{R}^2	0.922	0.964	0.980	0.918	0.990

	Zero-order reaction ^a		The 1st-order reaction ^b		The 2nd-order reaction ^c	
Zeaxanthin			k/(h ⁻	(R^2)		
	k	R ²	k	R ²	k	R ²
10 µg/g	0.7033	0.9275	0.1460	0.9383	0.0372	0.8273
50 µg/g	3.1206	0.9319	0.1404	0.9482	0.0077	0.8633
100 µg/g	6.7028	0.9453	0.1325	0.9609	0.0031	0.8686
200 µg/g	13.569	0.9521	0.1451	0.9609	0.0019	0.8603
400 µg/g	27.9400	0.9805	0.1470	0.9814	0.0009	0.8910

Table S2.

 $a \quad C_A - C_{A,0} = -kt$

 $b \quad \ln \left(C_A / C_{A,0} \right) = -kt$ $c \quad 1 \quad 1$

$$\frac{1}{C_A-C_{A,0}}=kt$$

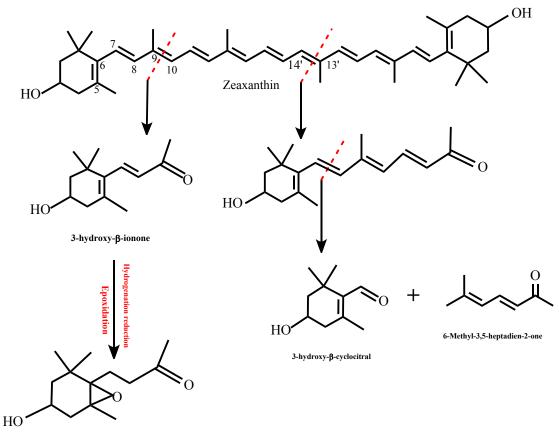
where C_A was the zeaxanthin concentration at time t and $C_{A,0}$ was the initial zeaxanthin concentration, and the reaction rate constants k were based on linear regression.

Table	S3 .
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peak	compounds	relative percentage (%) ^a
1	6-methyl- 3,5-heptadien-2-one	1.77
2	3-hydroxy-5,6-epoxy-7,8-dihydro-β-ionone	0.49
3	3-hydroxy-β-ionone	0.35
4	3-hydroxy-β-cyclocitral	1.02

^{*a*} Data are expressed as percentage values with respect to the total volatile profile.

Figure S1.



³⁻hydroxy-5,6-epoxy-7,8-dihydro-β-ionone