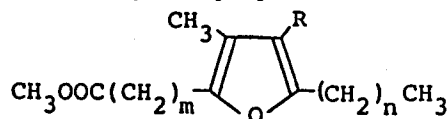


## ERRATA

In the paper "The Composition of Furan Fatty Acids in the Crayfish," by Kazuo Ishii et al., Vol. 23, No. 7, pp. 694-700, there was an omission of one scheme in each of Tables 1 and 2.

TABLE 1

Furan Fatty Acid Composition of Sterol Esters<sup>a</sup> from Crayfish Hepatopancreas



Peak no.	ECL <sup>b</sup>	CCL <sup>c</sup>	m	n	R	Wt. % of total fatty acids	Peak no.	ECL	CCL	m	n	R	Wt. % of total fatty acids
1*	15.09	12	2	4	H	0.02	18*	21.38	19	8	5	H	0.03
2	15.42	12	2	4	CH <sub>3</sub>	0.04	19 (F <sub>4</sub> )	21.51	18	10	2	CH <sub>3</sub>	3.45
3*	17.32	14	4	4	CH <sub>3</sub>	trace <sup>d</sup>	20*	21.97	19	8	5	CH <sub>3</sub>	trace
4*	17.79	14	6	2	CH <sub>3</sub>	trace	21*	22.12	19	9	4	CH <sub>3</sub>	0.06
5*	18.33	16	4	6	H	0.09	22*	22.36	19	10	3	CH <sub>3</sub>	0.11
6*	18.46	16	6	4	H	0.09	23 (F <sub>3</sub> )	22.51	20	10	4	H	0.93
7	18.80	16	8	2	H	0.03	24 (F <sub>4</sub> )	23.20	20	10	4	CH <sub>3</sub>	11.91
8 (F <sub>0</sub> )	19.07	16	6	4	CH <sub>3</sub>	1.76	25	23.66	20	12	2	CH <sub>3</sub>	0.09
9 (F <sub>1</sub> )	19.34	16	8	2	CH <sub>3</sub>	0.98	26*	23.93	21	10	5	CH <sub>3</sub>	0.02
10*	19.56	17	8	3	H	0.02	27*	24.06	21	11	4	CH <sub>3</sub>	0.02
11*	19.95	17	6	5	CH <sub>3</sub>	0.02	28*	24.40	22	10	6	H	0.04
12*	20.06	17	7	4	CH <sub>3</sub>	0.04	29	24.60	22	12	4	H	0.02
13*	20.23	17	8	3	CH <sub>3</sub>	0.10	30	25.26	22	12	4	CH <sub>3</sub>	0.11
14 (F <sub>2</sub> )	20.44	18	8	4	H	1.64	32	24.24				olefinic F <sub>3</sub> <sup>e</sup>	0.02
15	20.72	18	10	2	H	0.07	33	24.34				olefinic F <sub>4</sub> <sup>e</sup>	0.02
16*	20.98	18	6	6	CH <sub>3</sub>	0.02	34	25.61				H	0.11
17 (F <sub>3</sub> )	21.09	18	8	4	CH <sub>3</sub>	6.78	35	25.96				CH <sub>3</sub>	0.13

<sup>a</sup>Represents 18.9% of the total lipids of hepatopancreas.

<sup>b</sup>Equivalent chain length.

<sup>c</sup>Carbon chain length.

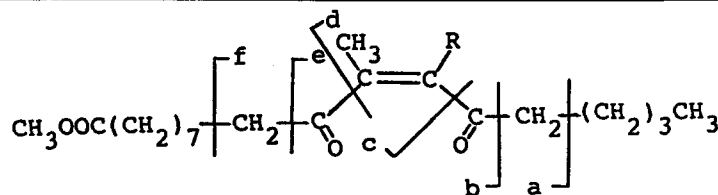
<sup>d</sup>Less than 0.01%.

<sup>e</sup>F<sub>3</sub> or F<sub>4</sub> methyl ester with one additional double bond, conjugated with a furan ring.

\*The acid is unknown.

TABLE 2

Comparison of Major Mass Spectral Fragments of Dimethyldiketo-ene (36) with Those of Monomethyldiketo-ene (40)



Fragments	36 (R=CH <sub>3</sub> ) m/z (rel. int.) <sup>a</sup>	40 (R=H) m/z (rel. int.) <sup>a</sup>
M*	352 (48)	338 (9)
M* - H <sub>2</sub> O	334 (68)	320 (20)
M* - OCH <sub>3</sub>	321 (28)	307 (22)
a	295 (55)	281 (3)
a - CH <sub>3</sub> OH	263 (100)	249 (10)
b	281 (4)	267 (0)
b - CH <sub>3</sub> OH	249 (15)	235 (40)
c	252 (1)	239 (29)
c - CH <sub>3</sub> OH	221 (3)	207 (12)
d	153 (4)	139 (23)
e	181 (25)	167 (100)
f	195 (90)	181 (42)
g <sup>b</sup>	177 (29)	163 (18)
h <sup>c</sup>	205 (28)	191 (20)

<sup>a</sup>Relative intensity.

<sup>b</sup>The fragments corresponding to base peaks of olefinic F acids with a double bond in the alkyl chain (Fig. 2B).

<sup>c</sup>The fragments corresponding to base peaks of olefinic F acids with a double bond in the alkylcarboxyl chain (Fig. 2C).