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## LETTERS TO THE EDITOR

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### THE TAUTOMERISM OF 5-MERCAPTO-2-ACETYLFURAN

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By acidifying the sodium salt of 5-mercapto-2-acetylfuran (I) [1] with formic acid, we obtained for the first time 5-mercapto-2-acetylfuran (II), yellow crystals, mp 121° (ex hexane). Found: C 50.78, 51.04; H 3.83, 4.27; S 22.13%. Calculated for  $C_6H_6O_2S$ : C 50.68; H 4.25; S 22.55%. The IV spectrum of II was determined in octane, dioxane, and methanol. For comparison purposes, the spectrum of 5-methylmercapto-2-acetylfuran (III) was observed in the same solvents.

The spectrum of III exhibits a band with  $\lambda_{\max}$  302 m $\mu$ , lg  $\epsilon$  4.09 (in octane). With the other solvents there is a bathochromic shift of 5-10 m $\mu$ . The spectrum of II has two absorption zones, an intense band with  $\lambda_{\max}$  270 m $\mu$  and a less intense wide band in the 300 m $\mu$  region. Both bands vanish in alcoholic alkali, and one with  $\lambda_{\max}$  380 m $\mu$  appears. A similar band is found in the spectrum of the sodium salt of I (obtained by crystallization from water, after evaporating an aqueous solution under reduced pressure; yellow crystals with a violet reflex, mp 180°, decomp.)

These results make it possible to assume that, in solution, II exists as two tautomers A and B



with tautomer B, which absorbs in the 270 m $\mu$  region, prevailing. We assign the band in the 300 m $\mu$  region to absorption by the thiol form A. In alkali both tautomers are converted into the sodium salt I.

We are continuing our research on the subject.

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