## 5-(3,4-Dimethoxybenzylidene)-1,3-diethyl-2-thioxodihydropyrimidine-4,6(1H,5H)-dione

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Received: 13 September 2003 / Accepted: 4 February 2004 / Published: 24 February 2004


5-(3,4-Dimethoxybenzylidene)-1,3-diethyl-2-thioxodihydropyrimidine-4,6(1H,5H)-dione 3 was prepared by Knoevenagel condensation of 3,4di-methoxybenzaldehyde 1 and N,N-diethylthiobarbituric acid 2 in ethanol using piperidine as a base [1,2].
$\mathrm{N}, \mathrm{N}$-diethylthiobarbituric acid $2(6.0 \mathrm{~g}, 0.03 \mathrm{~mol})$ and 3,4-dimethoxy- benzaldehyde $1(5.0 \mathrm{~g}, 0.03 \mathrm{~mol})$ in ethanol ( 50 mL ) was heated under reflux for ten minutes. Piperidine $(1.5 \mathrm{~mL})$ was added in one portion and the reflux was continued for further three hours. The reaction mixture was cooled to room temperature and the solid formed was filtered, washed with cooled ethanol ( $2 \times 50 \mathrm{~mL}$ ) and dried. 5-(3,4-
Dimethoxybenzylidene)-1,3-diethyl-2-thioxodihydropyrimidine-4,6( $1 H, 5 H$ )-dione 3 was recrystallized from ethanol as deep yellow crystals ( $9.85 \mathrm{~g}, 94 \%$ ).
M.p. $169{ }^{\circ} \mathrm{C}$ (EtOH, uncorrected).

UV 1 max $(\mathrm{nm} ; E t O H) / \mathrm{e}\left(\mathrm{dm}^{3} . \mathrm{mol}^{-1} . \mathrm{cm}^{-1}\right)$ 260/2009, 370/2875 and 220/ 2475.

IR nmax ( $\mathrm{cm}^{-1}$; KBr Disk) $1694(\mathrm{C}=\mathrm{O}), 1670(\mathrm{~N}-\mathrm{CO}-\mathrm{N}), 1632(\mathrm{C}=\mathrm{C})$.
${ }^{1} \mathrm{H}-\mathrm{NMR}\left(400 \mathrm{MHz} ; \mathrm{CDCl}_{3} ; \mathrm{Me}_{4} \mathrm{Si}\right) \mathrm{dH} 8.48(1 \mathrm{H}, \mathrm{s}$, olefinic Proton), $8.35(1 \mathrm{H}), 7.84(1 \mathrm{H}, \mathrm{d}, J=8.5 \mathrm{~Hz})$, $6.96(1 \mathrm{H}, \mathrm{d}, J=8.5 \mathrm{~Hz}), 4.51,4.50\left(4 \mathrm{H}, \mathrm{t}, 2 \mathrm{xCH}_{2}\right), 4.00,3.98(5 \mathrm{H}, \mathrm{s}, 2 \mathrm{xMeO}), 1.33,1.28\left(6 \mathrm{H}, \mathrm{q}, 2 \mathrm{xCH}_{3}\right)$.
${ }^{13} \mathrm{C}-\mathrm{NMR}\left(100 \mathrm{MHz} ; \mathrm{CDCl}_{3} ; \mathrm{Me}_{4} \mathrm{Si}\right)$ d C $178.8(\mathrm{C}=\mathrm{S}), 161.4,160.9(2 \mathrm{xC}=\mathrm{O}), 154.74,148.16(\mathrm{CH}$ olefinic), $133.6,131.97,122.5,120.8,117.38,109.9,55.4,55.02\left(2 \mathrm{x} \mathrm{CH}_{3} \mathrm{O}\right), 44.1,43.5\left(2 \mathrm{xCH}_{2}\right), 13.0,12.7\left(2 \mathrm{xCH}_{3}\right)$.

Anal.Calc. for C17H20N2O4S ( 348.418): C 58.60, H 5.79 , N 8.04; found : C 58.45, H 5.89, N 7.95.

## References

1. G. Jones, Org. React., 1967, 15, 204.
2. K. Tanaka, X. Chen and F. Yoneda, Tetrahedron, 1988, 44, 3241.
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