REACTION OF 2-AMINOMETHYLETHYLENEIMINE WITH SOME KETO AND

 α - AND β -DICARBONYL COMPOUNDS*

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Continuing our research on the reactivity of 2-aminomethylethyleneimine (I) in reactions with carbonyl compounds, as a result of which we obtained a new heterocyclic system – 1,3-diazabicyclo[3.1.0]hexane – we present the results of an investigation of the reactions of I with several keto and α - and β -dicarbonyl compounds (with acetone, glyceraldehyde, diacetyl, and acetylacetone).

We found that 2-substituted-1,3-diazabicyclo[3.1.0]hexanes IIa,b are formed as the chief reaction products in the reaction of I with acetone and glyceraldehyde.

EXPERIMENTAL

 $\frac{2-\text{Methyl}-2-(2-\text{hydroxyethyl})-1, 3-\text{diazabicyclo}[3.1.0]\text{hexane (IIa).} \text{ This compound had}}{\text{bp 95-97° (2 mm) and nD^{22} 1.4635. Found: C 58.7; H 10.5; N 19.5%. C_7H_14N_2O. Calculated:} C 58.7; H 10.5; N 19.7%. IR spectrum, cm⁻¹: 3010, 3070 <math>\left(\gamma \times \bigvee_{CH_2}^{CH_2}\right)$; 3340-3400 (OH, NH); 1220 $\left(\delta \times \bigvee_{CH_2}^{CH_2}\right)$.

 $\frac{2-(1,2-\text{Dihydroxyethyl})-1,3-\text{diazabicyclo}[3.1.0]\text{hexane (IIb).}}{(2 \text{ mm}) \text{ and } n_D^{-2} 1.5324.} \text{ Found: C 50.2; H 8.7; N 19.3%. C_6H_{12}N_2O_2.} \text{ Calculated: C 50.0; H 8.3; N 19.4. IR spectrum, cm^{-1}: 3010, 3060 (vN (); 3340-3400 (v OH, NH); 1240 (<math>\delta N$ ().

Condensation products of the IIIa,b type were obtained by reaction of I with the simplest $\alpha-$ and $\beta-dicarbonyl compounds:$

R H H $R = COCH_3;$ R H H $R = COCH_3;$ $R = CH_2COCH_3$

2-Methyl-2-acetyl-1,3-diazabicyclo[3.1.0]hexane (IIIa). This compound had mp 42° (from hexane). Found: C 60.4; H 8.4; N 19.8%. $C_7H_{12}N_2O$. Calculated: C 60.3; H 8.6; N 20.0%. IR spectrum, v, cm⁻¹: 3290 (NH); 3012, 3080 (N \checkmark); 1690 (C=0).

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 $\frac{2-\text{Methyl}-2-(\text{acetonyl})-1,3-\text{diazabicyclo}[3.1.0]\text{hexane (IIIb).}}{(3 \text{ mm}) \text{ and } n_D^{-2} 1.5460.}$ Found: C 62.1; H 9.2; N 18.0%. C₈H₁₄N₂O. Calculated: C 62.0 H 9.1; H 18.2%. IR spectrum, v, cm⁻¹: 3280 (NH); 3010, 3080 (N \checkmark); 1700 (C=0).

LITERATURE CITED

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