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Note

## Investigation of IR-Spectra of *N*-Benzoyl-*N*-phenylhydroxylamine in Carbon Tetrachloride

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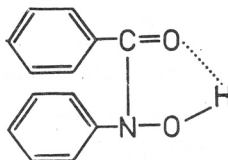
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Spectra of carbon tetrachloride solutions of *N*-benzoyl-*N*-phenylhydroxylamine in concentrations between  $1 \times 10^{-4}$  and  $25 \times 10^{-4}$  M, in the spectral range from  $2800 \text{ cm}^{-1}$  to  $4000 \text{ cm}^{-1}$  were measured. Appearance of the absorption band at  $3070 \text{ cm}^{-1}$  is ascribed to C—H stretching, and the band at  $3250 \text{ cm}^{-1}$  to H...O—H bond. The validity of Beer's law for the band at  $3250 \text{ cm}^{-1}$  and the fact that no bands between  $3400$  and  $4000 \text{ cm}^{-1}$  were detected indicate the presence of an intramolecular H bond.

Recently *N*-benzoyl-*N*-phenylhydroxylamine (BPH) has been applied for the solvent extraction of some elements<sup>1,2</sup>.

In our determination<sup>3</sup> of equilibrium constants for complexes of vanadium with BPH, we used nonpolar organic solvents. In this respect it was important to know the behaviour of BPH in such solvents.



BPH is a polar molecule capable of forming hydrogen bonds. This may lead to strong selfassociation in non-polar organic solvents. IR spectroscopy was used to show the type of hydrogen bonding.

Spectra of solutions of BPH in  $\text{CCl}_4$  at concentrations from  $1 \times 10^{-4}$  to  $25 \times 10^{-4}$  M were measured. A 102.6 mm cell with glass windows about 0.8 mm thick was used. The spectra were obtained with a VEB Carl Zeiss, Jena, double beam IR spectrophotometer Model UR-10. The spectral slit width ranged from 2 to  $4 \text{ cm}^{-1}$ .

In the spectral range from  $2800 \text{ cm}^{-1}$  to  $4000 \text{ cm}^{-1}$  two absorption bands were detected, at  $3070 \text{ cm}^{-1}$  and  $3250 \text{ cm}^{-1}$  (Fig. 1). From the analogy of these bands and literature data<sup>4-7</sup>, we ascribed the band at  $3070 \text{ cm}^{-1}$  to C—H stretching and the band at  $3250 \text{ cm}^{-1}$  to the intramolecularly bonded OH group. The fact that no bands between  $3400$  and  $4000 \text{ cm}^{-1}$  were detected indicates the absence of free OH groups and hence intramolecular bonding.

This is further substantiated by the fact that the plot of the absorption intensity at  $3250 \text{ cm}^{-1}$  vs. BPH concentration in  $\text{CCl}_4$  gives a straight line (Fig. 2).

Hence, it is concluded that at these concentrations of BPH no association takes place.

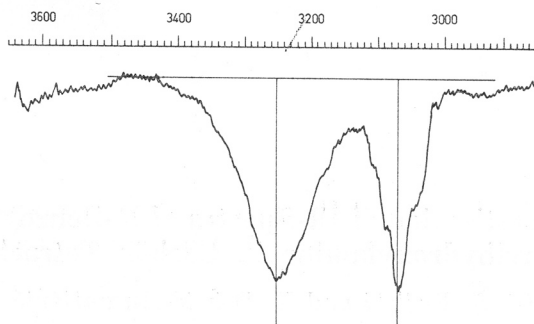


Fig. 1. IR absorption spectrum of  $8 \times 10^{-4}$  M BPH solution in  $\text{CCl}_4$ .

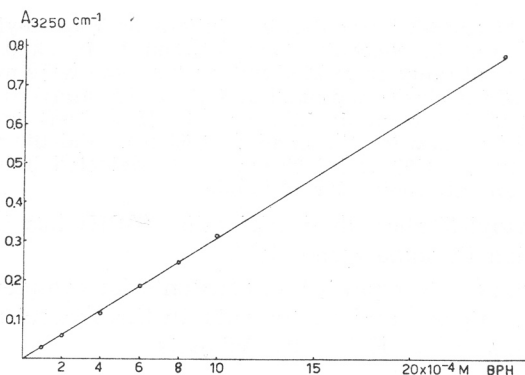


Fig. 2. Calibration curve for the BPH solutions in  $\text{CCl}_4$  at  $3250 \text{ cm}^{-1}$ .

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#### IZVOD

##### Ispitivanje IR spektra *N*-benzoil-*N*-fenilhidroksilamina

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Snimljeni su IR spektri *N*-benzoil-*N*-fenilhidroksilamina u ugljen tetrahloridu za koncentracije od  $1 \times 10^{-4}$  do  $25 \times 10^{-4}$  M, i u spektralnoj oblasti od 2800 do  $4000 \text{ cm}^{-1}$ . Pojava trake na  $3070 \text{ cm}^{-1}$  pripisuje se C—H istežućoj vibraciji, a trake na  $3250 \text{ cm}^{-1}$  O...H...O vezi. Utvrđeno je da za traku na  $3250 \text{ cm}^{-1}$  važi Beer-ov zakon što, pored odsustva traka u oblasti od  $3400$  do  $3700 \text{ cm}^{-1}$ , ukazuje na postojanje intramolekulske vodonične veze.

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