

Synthesis, characterisation and biocidal studies of Cu(II) and Co(II) complexes with drug alprazolam [8-chloro-1-methyl-6-phenyl-4H-1,2,4-triazolo (4,3-a) (1,4) benzodiazepine]

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ABSTRACT

In the present study, Cu(II) and Co(II) complexes have been prepared by reacting metal Nitrate with the ligand (Alprazolam), which have the general structural composition $[(L) M(NO_3)_2]$, where $M=Cu(II)$ and $Co(II)$ and $L=Alprazolam$ [8-chloro-1-methyl-6-phenyl-4H-1,2,4-triazolo (4,3-a) (1,4) benzodiazepine]. The complexes were characterised on the basis of elemental analysis, molecular weight and IR data. The shift in the characteristic infrared frequencies of the free ligand bands confirm the coordination. The IR data of the complexes reveal the bidentate nature with [N(1) and N(4)] sites of the ligand. All the complexes, ligand and metal salts have been screened for antibacterial and antifungal activities against selected microbes. Biocidal studies show that the Cu(II) complexes is more effective than Co(II) complexes.

Key words: Metal complexes, alprazolam and biocidal study.

INTRODUCTION

The complexes of 1,4-benzodiazepine with metal ions have also been reported in literature¹⁻⁵. Alprazolam a benzodiazepine drug is included in the category of tranquillizer drugs. It is used to treat anxiety disorders and as an adjunctive treatment for depression. It has been reported that complexes of 1,4-benzodiazepine also possess anticancer^{6,7} and pharmacological⁸ properties. Metal complexes of 1,4-benzodiazepines possessing biological activity may be even more active than free ligand⁹. Copper (II) complex of Alprazolam appears to be quite active, having a rapid onset of action and it also prolongs the duration as compared to that of Alprazolam itself. Different coordination modes have been including coordination with the neutral ligand viz. N(4), However, none of the modes has been ascertained. Neutral complexes of 1,4-

benzodiazepines such as Nitrazepam¹¹ where the ligands act as anions through labialization of proton at N(1) atom giving (imido) complexes, have been reported. The structure of the ligand Alprazolam (AZ) is given in figure (1).

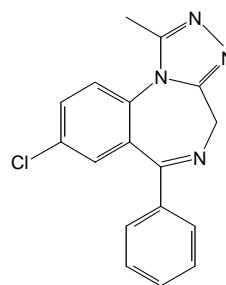


Fig. 1: Alprazolam (AZ)

The proposed work has been taken up to investigate the mode of coordination of the metal ion with (AZ) drug of increased biological activity of the drug on complexation as reported in the literature¹². The results will throw light on the mode of coordination of the metal ion with the AZ drug as well as on the complex formation of other benzodiazepines in general.

EXPERIMENTAL

All the chemical used were of AR grade and their solution were prepared in double distilled water. The complexes were prepared by mixing molar solutions of $\text{Cu}(\text{NO}_3)_2 \cdot 3\text{H}_2\text{O}$, $\text{Co}(\text{NO}_3)_2 \cdot 6\text{H}_2\text{O}$ (dissolved in double distilled water) and the ligand Alprazolam (dissolved in ethanol). The pH of the mixture was adjusted at 7.5-8.5.

The stoichiometry of the complexes of the drug (0.025M) with Cu(II) and Co(II) (0.025M) metal ions was found by carrying out potentiometric titration against standard (0.1M) NaOH solution in ethanol-water mixture. The pH changes observed during the titration were plotted against moles of alkali(m) added per moles of metal ion and ligand as depicted in figs. 3 and 4. The stoichiometric ratio is confirmed by *Job's method*¹³. The coloured precipitates were filtered, washed several times with hot water followed by ethanol to free it from the soluble impurities. The complex were finally dried in an oven at 100°C and stored in desiccator. The purity of the complexes was checked by TLC. The complexes were dissolved in benzene: Acetic Acid (2:1) and TLC was carried out in ethanol: Benzene (80:20) system. The Retention factor (R_f value) was calculated by following formula.

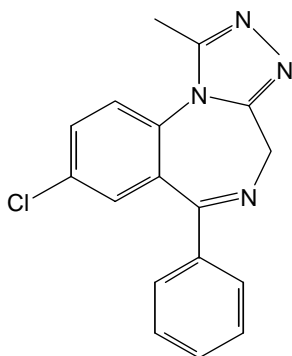


Fig. 2:

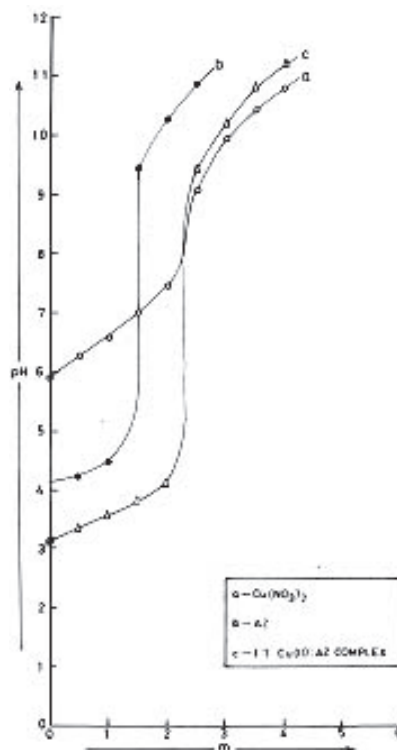


Fig. 3: System Cu(II): AZ

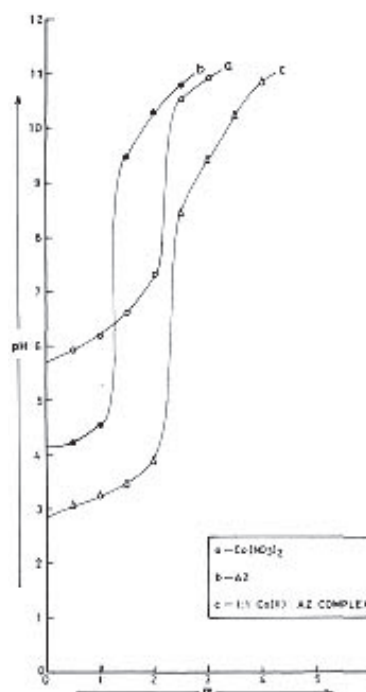


Fig. 4: System Co(II): AZ

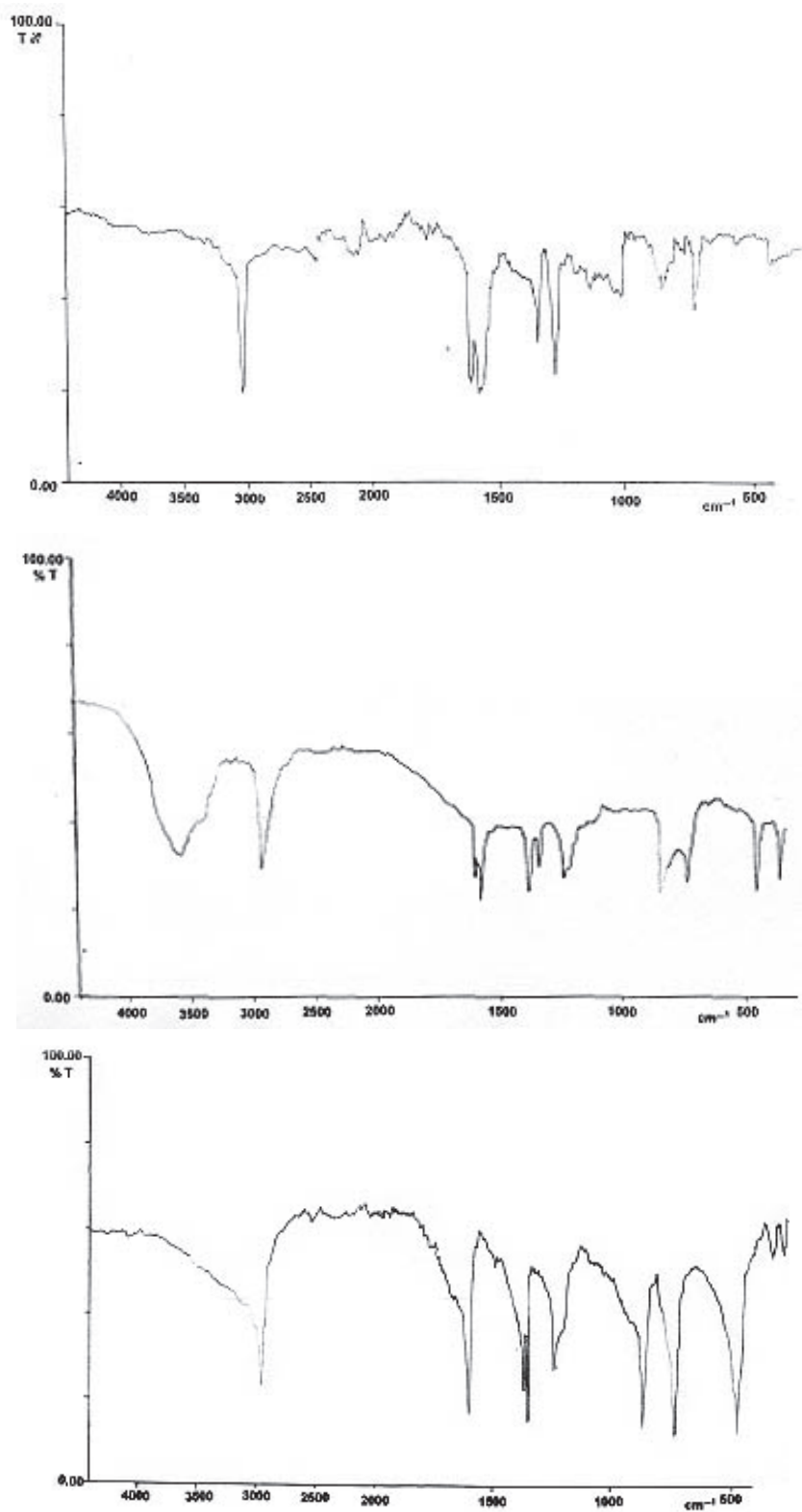


Fig. 5: IR Spectrum of AZ Complexes

suspension of the test organism separately. Standard drug was used to check and compare the activity of the complexes. It indicated that metal complexes of the drug had a significant activity at a lower concentration and bacteria were incubated at $24 \pm 1^\circ\text{C}$ for 20 hours. The data and results are presented in Table 3.

The IR spectra in KBr matrix were recorded on Perkin-Elmer 842-Spectrometer. Elemental analysis of C,H,N were carried out at CDRI, Lucknow.

RESULTS AND DISCUSSION

The molecular formula of the Cu(II) and Co(II) complexes correspond to $[(C_{17}H_{13}Cl N_4) \cdot Cu(NO_3)_2]$ and $[(C_{17}H_{13}Cl N_4) \cdot Co(NO_3)_2]$. The molecular weights 496.34 and 491.73 respectively for Cu(II), Co(II) complexes were determined on the basis of elemental analysis. The physical and analytical data of ligand Cu(II) and Co(II) complexes are given in table 2.

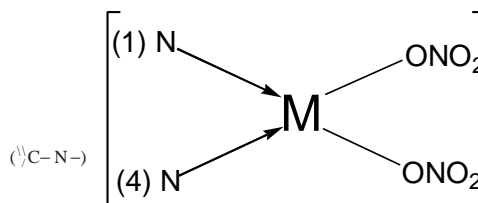
The IR spectrum of the ligand exhibits bands in the region 1628 cm^{-1} , 1280 cm^{-1} , 1600 cm^{-1} , 740 cm^{-1} , 2960 cm^{-1} and 1355 cm^{-1} which may be assigned to $\nu(C=N)$, $\nu(C-N)$, $\nu(C_6H_5)$, $\nu(Cl)$, $\nu(CH_2)$ and $\nu(-CH_3)$ respectively. The ligand band observed in the range of 1628 cm^{-1} undergoes the lower shifting to 1612 cm^{-1} and 1613 cm^{-1} in Cu(II) and Co(II) complexes respectively, indicating azomethine nitrogen N(4) atom of Benzodiazepine ring in coordination to the metal ion complexes. The Cu(II) and Co(II) complexes the band attributed to the vibration mode ν appear at 1263 cm^{-1} and 1265 cm^{-1} showing lower shifting as compared to the ligand (1280 cm^{-1}), indicating nitrogen (1) participating in complexation. In IR spectrum of the ligand the bands attributed to vibrational mode $\nu(-C_6H_5)$, $\nu(CH_3)$, $\nu(-Cl)$, $\nu(CH_2)$ appears at 1600

cm^{-1} , 1355 cm^{-1} , 740 cm^{-1} and 2960 cm^{-1} showing small positive shift in complexation.

The presence of new bands at 480 cm^{-1} and 340 cm^{-1} in Cu (II) complex and 492 cm^{-1} and 340 cm^{-1} in Co (II) are attributed to ν (M-N) linkage¹⁶. Other strong bands at 1383 cm^{-1} and 861 cm^{-1} in Cu(II) complexes and 1370 cm^{-1} and 850 cm^{-1} in Co (II) complex suggest monodentate nitrate group in complexes¹⁷. The representative structure of the complex may be shown as

Biocidal Activity

The biological activities of Cu(II) and Co(II) complexes, metal nitrates and ligand were screened against bacteria – *E. coli*, *S. typhi*, *B. subtilis* and *S. aureus* and fungi – *A. flavous*, *A. niger*, *P. triticena* and *F. species*. The zone of inhibition of Cu(II) and



Probable structure of metal complexes
Where M=Cu (II), Co (II)

Co(II) complexes on bacteria and fungal species are presented in Table 3.

The above data indicate that the zone of inhibition at 500 ppm concentration is best as compared to the standard drug. The results show that the Cu(II) complex of Alprazolam is more effective towards all fungi and bacteria as compared to Co(II) complex of Alprazolam and drug Alprazolam itself.

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