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Synthesis and Characterization of Mixed Ligand Complexes of Transition Metal Chelates of 1-Nitroso-2-naphthol and 8-Hydroxyquinoline with Quinaldinic Acid

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ABSTRACT

Mixed ligand transition metal complexes having general formula ML, HL', where M = Cu(II) or Pd(II); L deprotonated 1-nitroso-2-naphthol or 8-hydroxyquinoline; HL' = Quinaldinic acid have been synthesised and characterised on the basis of Infrared, electronic absorption spectral data, magnetic and conductance measurements. The above study reveals the octahedral geometry of the complexes. Microbiological studies revealed that all prepared mixed ligand complexes are relatively potential inhibitor of the growth of both the test bacteria and one fungi.

Complexes, Quinaldinic Key words: Mixed Ligand acid, 8-Hydroxyquinoline, 1-Nitroso-2-naphthol.

INTRODUCTION

Quinaldinic acid has been extensively investigated for their possible roles in analytical determination of metal ions¹⁻ ⁴. They have replaceable hydrogen atom as well as two donor atoms suitably placed to form a five membered chelating ring with metal ion. The versatile chelating ability of quinaldinic acid with various metals is well established⁵⁻¹⁰. In the present communication, we report the synthesis and characterization of a number of mixed ligand complexes of Cu(II) or Pd(II) with above mentioned ligands.

EXPERIMENTAL

Quinaldinic acid(HQuinA), 1-Nitroso-2-naphthol(1N2N) and 8-Hydroxyquinoline(8HQ) of AnalaR grade were used as such.

Preparation of the complexes

To the suspension of 0.01 mole transition metal salt of organic acid(1N2N or 8HQ) in absolute ethanol, 0.01 mole quinaldinic acid(HQuinA) was added. The whole reaction mixture was refluxed with constant stirring on a hot plate of magnetic stirrer for 1-2 hours and cooled to give characteristic colour solid adduct. The adduct was filtered, washed with absolute ethanol and dried in an electric oven at 100°C.

Some physical properties and analytical data of the ligand(HQuinA) and the new mixed ligand complexes obtained are listed in Table-1.

The mixed ligand complexes are generally coloured. They are appreciably soluble in most polar solvents such as methanol, DMF etc. but are insoluble in non-polar solvents such as benzene, toluene, ether etc. The complexes are found to be stable when stored under dry conditions. Melting/decomposition temperatures of the complexes have been found to be higher than those of the corresponding ligand, indicating thereby their greater stability.

Molar conductance

Molar conductance values of all these complexes were measured in methanol at 23° C at a concentration of 10^{-3} M. The values are given in Table-1. Low values of molar conductance (5.2 - 6.2 ohm⁻¹cm²mol⁻¹) show that these complexes are non-electrolyte in nature.

Infrared spectra

Infrared spectra of the ligand(Quinaldinic acid) and their mixed ligand Cu(II) or Pd(II) complexes were recorded in KCl phase between 4000-400 cm⁻¹ with the help of JASCO FTIR spectrophotometer model -5300. Selected IR absorption bands are shown in Table - 2.

IR spectra of quinaldinic acid & its complexes

The broad band at 3400 cm⁻¹ and multiple bands in the region 2700-1800 cm⁻¹ in the spectrum of ligand(HQuinA) indicate strong intramolecular hydrogen bonding involving the carboxyl hydrogen atom and nitrogen atom of quinoline ring. In the spectra of mixed ligand complexes, these bands have disappeared. However, the spectra of the complexes exhibit new broad bands of weak to medium intensities in the region 3400-3000 cm⁻¹, which may be attributed

O-H...O/N...H-O absorption. This suggests hydrogen bonding to be a dominant factor in stabilizing these complexes.

The medium intensity band at 1680 cm⁻¹ in the spectra of ligand, all probability, be assigned to the antisymmetric -COOH has shifted lower stretching frequency of the group, to frequencies by 5-30 cm⁻¹ in almost all the complexes. The 1620 cm⁻¹ and 1560 cm⁻¹ bands of the ligand too, in almost all the complexes, have shifted to lower frequencies by 10-40 cm⁻¹. The 1535 cm⁻¹, which appear as a medium intensity in the spectrum of the ligand, shows an appreciable shift of about 20-35 cm⁻¹ in the spectra of these complexes. These features are suggestive of coordination of ligand with Cu(II) or Pd(II) through oxygen atom of carboxylate(COO⁻) moiety.

The 1580 cm⁻¹ (assigned to C=N) band of the ligand has shifted to lower frequencies by 10-15 cm⁻¹ in almost all the complexes. These features suggest the coordination of ligand with Cu(II) or Pd(II) metal through nitrogen atom of quinoline ring.

Electronic absorption spectra

Electronic absorption spectra of the ligand(Quinaldinic acid) and their mixed ligand Cu(II) or Pd(II) complexes were recorded on PERKIN ELMER LAMBDA-15 UV-VIS spectro-photometer in paraffin solvent. The bands observed are given in Table -2.

A comparative look of electronic absorption spectral data of the ligand and their complexes indicate that n-p and p- p^* transition of the ligand have shifted to higher frequencies.

The electronic absorption spectra of mixed ligand Cu(II) complexes show broad band at 640 nm with quinaldinic acid.

The strong absorption in the region 570-575 nm in all the complexes attributed to ligand absorption and charge transfer transition, showing coordination of ligand with metal chelates by L-M interaction. However from the position of spectral bands and their low intensities it is supposed that the complexes are in octahedral geometry.

Magnetic moment

Magnetic moment of mixed ligand transition metal complexes have been measured by Can Faraday magnetic susceptibility balance at 30°C. The magnetic moment values are shown in Table-1.

The spin only value of magnetic moment of Cu(II) octahedral complexes is 1.73 BM corresponding to one unpaired electron. The magnetic moment(μ_{eff}) values for Cu(II) complexes are in the range of 1.96 - 2.00 BM, suggest the octahedral geometry of the complexes. The complexes of Pd(II) are diamagnetic.

Microbiological Studies

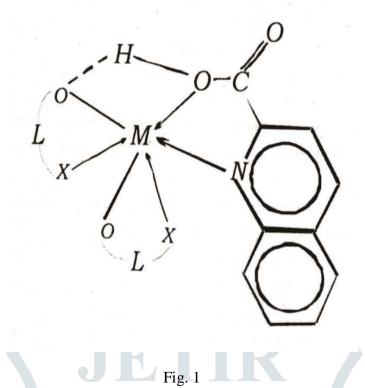
Minimum inhibitory concentration(MIC) values (μ gml⁻¹) of mixed ligand transition metal complexes have been examinated by Serial dilute method for activity against some bacteria¹¹, viz., E. coli. S. aureus and fungi¹² viz. C. albicans. Althought a few of the mixed ligand complexes exhibited significant inhibition whereas other did not show any activity. The results are summarized in Table – 3.

The results of antibacterial and antifungal activity evaluation revealed that three of the mixed ligand complexes, $Cu(8HQ)_2HQuinA$, $Cu(1N2N)_2.HQuinA$ and $Pd(1N2N)_2.HQuinA$ were relatively potential inhibitor of the growth of both the test bacteria and one fungi.

The MIC value of Cu(1N2N)₂.HQuinA & Pd(1N2N)₂.HQuinA at concentration of 40 μ gml⁻¹ while Cu(8HQ)₂.HQuinA at concentration of 20 μ gml⁻¹ for E. Coli. The MIC values of Cu(8HQ)₂.HQuinA, Cu(1N2N)₂.HQuinA and Pd(1N2N)₂.HQuinA at concentration 40 μ gml⁻¹ and 20 μ gml⁻¹ respectively for fungi C. albicans. The MIC values of all three oxygen bridge transition metal complexes at the same concentration 40 μ gml⁻¹ for S. aureus.

Structure & bonding

Based on the analytical and spectral(infrared & electronic absorption) studies, the structure and bonding of the newly prepared mixed ligand Cu(II) or Pd(II) complexes involving some chelating organic acids and quinaldinic acid may tentatively be proposed as shown in Fig. 1.



[Where M = Cu(II) or Pd(II); L = deprotonated 8-hydroxyquinoline or 1-nitroso-2-naphthol; X = O or N.

Table 1:								
Compound	Colour	Melting	Molar	Magnetic	Analysis % found (calcd.)			
		Decomp.	cond.	moment	с	н	N	м
Quinaldinic acid (HQuinA)	White	157m		-	.	-	-	<u>.</u>
Cu(8HQ) ₂ . HQuinA	Dark green	275d	5.2	1.96	64.00 (64.10)	3.00 (3.60)	7.90 (8.00)	12.00 (12. 9 0)
Cu(1N2N) ₂ . HQuinA	Greenish grey	260md	5.5	2.00	61.90 (62.06)	3.10 (3.27)	7.24 (7.30)	12.00 (12.18)
Pd(1N2N) ₂ . HQuinA	Greenish brown	280d	6.2	Diamag	55.78 (57.08)	2.90 (3.04)	6.34 (6.74)	16.90 (17.01)

Table 2:							
Compound	Selected IR	Selected IR absorption bands (in cm ⁻¹)					
Quinaldinic acid (HQuinA)	3400br 2700-1800v	1680m	1620s	1580sh 1500s			
Cu(8HQ) ₂ .HQuinA	3500-3200∨ 2910w	1675sh	1640s	1565m 1535sh	1510s	640, 575	
Cu(1N2N) ₂ .HQuinA	3400-3200v 2950m,2910r	1680sh n	1630s	1565m 1540sh	1515w	800, 570	
Pd(1N2N) ₂ .HQuinA	3560-3450v	1750sh	1660m	1590sh 1570m	1500s	950,655,480	

Table O.

Br = broad, m = medium, s = strong, sh = shoulder, v = very strong, w = week.

Table 3 :

Compound	Minimum inhibitory concentraction (in µgmol ⁻¹)					
	E. Coli	S. aureus	C. albicans			
Cu (8HQ) ₂ .HQuinA	40	20	40			
Cu(1N2N) ₂ .HQuinA	20	20	40			
Pd(1N2N) ₂ .HQuinA	20	20	20			

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