

ACOUSTIC PARAMETERS OF SOME HALOSUBSTITUTED CHALCONEIMINES IN BINARY MIXTURE OF DIOXANE+WATER SOLVENT

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ABSTRACT

Chalconeimines are the phenolic schiff base an important class of compound. In present investigation complex formation between chalconeimine and solvant are studied by evaluating certain acoustic parameter. The **Complex formation in binary liquid mixtures** of 1-4 dioxane and water with benzaldehyde chalconeimine -K1 and 2chloro 3- fluro benzaldehyde chalconeimine -K2 has been studied by measuring ultrasonic velocity at 2 MHz, by using in concentration range of 0.01 M at 32 °C the under investigation parameters are ultrasonic velocity (u), adiabatic compressibility (\(\beta\)s), intermolecular length,(Lf) Apparent molal volume(Φv), Apparent molal Compressibility(Φk), specific acoustic impedance(Z) and Relative association(RA) These parameters have been utilized to study the solute-solute interactions in these systems. The ultrasonic velocity shows a decrease in value with respect to concentraction and adiabatic compressibility show corresponding increase in value with respect to concentraction for these ligands. The results indicate the occurrence of complex formation between unlike molecules through intermolecular hydrogen bonding between oxygen atom of dioxane molecule and hydrogen atom of ligands. The variation of these parameters with percentage indicates the possibility of the complex formation in these systems.

Keywords: Acoustical impedance, Dioxane+water mixtures, halo substituted chalconeimine

1 Introduction

Ultrasonic is the branch of acoustic, which consists of waves of high frequencies. Ultrasonic study has wide range of application in material science, agriculture, medicine, biology, industry, oceanography, sonochemistry research due to its non-destructive nature[1]. It is the technique used for the study of molecular interaction in liquids. Ultrasonic technique is used to study the nature of molecular interaction in liquids [2-5.] also to study the ion-solvent interactions in aqueous and non aqueous solutions[6]. Recent developments have made it possible to use ultrasonic energy in medicine, engineering, agriculture and other industrial applications [7-8]. In solution of ionic solute the attraction between the solute and solvent is essentially of ion-dipole interaction depends mainly on ion size and polarity of the solvent [9]. The measurements of ultrasonic waves are useful in study of molecular interactions in liquids which provides valuable information regarding internal structure, complex formation, internal pressure and molecular association [10]. The dissolution of electrolyte in a solvent causes a volume contraction due to interaction between ions and solvent molecules and this may influence other acoustical properties of solution. In recent years, the studies acoustical properties of aqueous electrolytic solutions [11] have been found to be useful in understanding the specific ion-ion and ion-solvent interaction in solutions. strength of ion-dipole attraction is directly proportional to the size of the ion, charge and the magnitude of the dipole, but inversely proportional to the distance between the ion and the dipolar molecule. The accurate measurement of density, viscosity, ultrasonic velocity and hence the derived parameters such as adiabatic compressibility, apparent molal compressibility and apparent molal volume will give significant information regarding the state of affairs in a solution [12].

2 .Experimental

2.1 Material and Methods

halosubstituted chalconeimine are synthesized by amination of chalcone with substituted amines. this chalcone are synthesized by using general claisen Schmidt method. For evaluating the acoustic properties the very pure and analytical grade solvent and extra pure double distilled water is used. The dencities of pure solvent and solutions are determined by using specific gravity bottle. The ultrasonic velocity measurements were made using a crystal controlled variable path ultrasonic interferometer (Mittal Enterprise, Model F-05) of 2 MHz with accuracy of (±) 0.03 %

The ligands used for viscometric study are 4-fluro benzaldehyde chalconeimine -K1 2 -

chloro 3- fluro benzaldehyde chalconeimine - K2

Theory and calculation

Adiabatic compressibility (β), Apparent molal volume (Φ v), Apparent molal compressibility (Φ ks), Intermolecular free length (Lf), Specific acoustic impedance (Z) and Relative association (R) were calculated by using following equations

$$\begin{split} \beta s &= 100 \, / \, Us^2 ds & (1) \\ \beta o &= 100 \, / \, Uo^{2*} \, do & (2) \\ \varphi v &= (M/ds) + [(do - ds)10^3] \, / \, m \, ds do & (3) \\ \varphi k(s) &= [1000 \, (\beta s do - \beta o \, ds) \, / \, m \, ds do] + (\beta s \, M) \\ /ds) & (4) \\ Lf &= K \times \sqrt{\beta} s & (5) \\ Z &= Us ds & (6) \end{split}$$

where,

do = Density of pure solvent,

m = Molality, M = Molecular weight of solute,

 β o = Adiabatic compressibility of pure solvent

 β s =Adiabatic compressibility of solution

K = Jackbson's constant

Us = Ultrasonic velocity in the solution in m/s. β s is in bar⁻¹ and φ k(s) is in cm³ mol⁻¹ bar⁻¹

Table - 1 Acoustic parameters of fluro benzaldehyde chalconeimine -K1

% Dioxa ne	Mole fracti on of Diox ane	Ultraso nic Velocity u×10 ³	Adiabatic compressi bility βs ×10 ⁻⁷	Intermole cular free length Lf×10 ⁻⁹	Apparent molal volume Φv ×10 ⁻³	Apparent molal Compressi bilityΦk×1	Relativ e associa tion R _A	Specific acoustic impeda nce Zs×10 ⁶
100	1	3332	0.853244	0.601	7075.684	1.05	0.5240	3517.39 9
90	0.656	2684.4	1.4538	0.7.84	14046.09	2.03	0.8369	2562.39 9
80	0.458 4	2074	2.57487	1.04	14889.72	3.59	1.1451	1872.56 1
70	0.330	1906	3.20516	1.16	18202.23	4.50	s1.2524	1636.92 2
60	0.240 9	1712.8	4.150907	1.32	23345.91	5.94	1.3832	1406.50 9

Table - 2 Acoustic parameters of 2 -chloro 3- fluro benzaldehyde chalconeimine -K2

% Diox ane	Mole fracti on of Diox ane	Ultras onic Velocit y (u)× 10 ³	Adiabatic compressi bility βs (bar- 1)×10 ⁻⁷	Intermole cular free length Lf×10 ⁻⁹	Appar ent molal volum e Φv ×10 ⁻³	Apparent molal Compressibility Φ k×10 ⁴	Relativ e associa tion (R _A)	Specifi c acousti c imped ance Zs×10 ⁶
100	1	1698.8	3.18288	1.16	4.2232	4.57	1.0599 9	1849.4 2
90	0.656	1660.8	3.58725	1.23	8.2431	5.24	1.4322	2611.1
80	0.458 4	1584	4.12	1.32	8.7182 0	5.88	1.4976	1642.9 9
70	0.330	1465.2	5.07479	1.46	10.720 4	7.24	1.7412 6	1344.8
60	0.240	1424.8	5.48278	1.5	12.872 5	7.88	1.8193 6	1280.0 9

Results and Discussion

In the present investigation different acoustic parameters such as adiabatic compressibility (B), apparent molal volume (Φv), apparent molal compressibility (Φk), acoustic impedance (Zs), relative associastion (RA) and intermoleculer free length (Lf) of the solutions in 0.01M concentrations of solute are determined at 303k and presented in table 1 and 2. The ultrasonic velocity of both the ligand is going to decrease with decrease in percentage of dioxane this indicates decrese the cohesion which is caused due to hydrogen bonding. The value of adeabetic compressibility increase (Bs) with decrease in percentage of solution may be due to departure of solvent molecules around ions supporting weak ion-solvent interactions. increases linearly in both the compound which represents the interaction of solute and solvent is weak. Value of adeabetic compressibility bs is more for FB ligand than CFB show that FB have strong solute solvent interaction. The value of apparent molal volume (Φv) is high in case of more polar substituent than less polar substituents. The apparent molal volume fv values increases with decrease in concentration of chalconesimines in binary mixtures this represents strong solute -solute and solute solvent interactions. Relative association RA increases linearly with decrease in percentage. It is influenced by breaking of bond in ions Decrease in RA with decrease in concentration suggests that breaking up of aggregation predominates the solvation of ions. . The value of relative association increases with decrease in concentration in all systems. It is found that there is weak interaction between solute and solvent. Relative association is more case of bulky and more polar substituents. The value of apparent molal compressibility(Φ k) increases with decrease in percentage of all systems, Showing weak electrostatic attractive force in the vicinity of ions causing electrostatic salvation of ions. Compressibility is more in case of bulky substituents. The values of intermolecular free length (Lf) indicates that with increase in concentration there is increase in ion-solvent interactions. The intermolecular free length increases due to greater force of interaction between solute and solvent by forming hydrogen bonding and less interaction between two solute molecules.

Conclusion

From the experimental data, acoustic parameter have been determined for halosubstituted chalconeimine in binary liquid(dioxane+water) at 32^{0} C the results have been used to study the ion-solvent interaction exists in the solvent mixture. From the magnitude of ϕk and β -coefficient values it can be concluded that the existence of ion-solvent interaction.

Acknowledgement

The author is thankful to dept of chemistry, Bar.R.D.I.K & Ny K.D. College Badnera for providing all the necessary facility.

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