

Date of Publication
June 2022

vidyavartaTM

International Multilingual Research Journal



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Art, Science & Commerce
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An Ultrasonic Study of 3-(2-Methoxyphenoxy)-1, 2-Propanediol in Different Binary Mixtures of Solvent At 308 K.

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ABSTRACT :

The ultrasonic velocity, density, viscosity and absorption have been measured for solutions of 3-(2-methoxyphenoxy)-1, 2-propanediol are presented at 308K. By taking measurements of 3-(2-methoxyphenoxy)-1, 2-propanediol at the concentration range is 0.002 to 0.01 mol dm⁻³ of each solution. The ultrasonic velocity, density and viscosity have been measured at 2MHz for the aqueous solutions of (i) 3-(2-methoxyphenoxy)-1,2-propanediol+ Acetone solvent (ii) 3-(2-methoxyphenoxy)-1,2-propanediol+ Ethanol solvents, at different concentrations at 308K. The acoustical parameters such as adiabatic compressibility (β), intermolecular free length (L_f), and acoustical impedance have been computed. The aim of the study is to increase the solubility, stability of the drugs by the formation of complexation. These properties are attributed to solute-solvent interactions through hydrogen bonding, molecular association, Solvent-solvent interaction etc. The results are interpreted in terms of molecular interactions occurring in these solutions.

Introduction :

In recent years acoustic and

thermodynamic behavior in many binary liquid mixtures has been analyzed by several researchers more than 5 decades. The analysis has been adequately employed in understanding the nature of molecular interactions in pure liquids and liquid mixtures.

Ultrasonic propagation parameters yield valuable information regarding the behaviors of liquid systems, because intramolecular and intermolecular association, dipolar interactions, complex formation and related structural changes affect the compressibility of the system which in turn produces corresponding variations in the ultrasonic velocity. The acoustical and thermodynamical parameters obtained in ultrasonic study show that the ion solvation is accompanied by the destruction or enhancement of the solvent structure¹⁻⁴. Excess thermodynamic properties of liquid mixtures are of great interest to conveniently design industrial processes and also to provide useful information on the molecular interactions required for optimizing thermodynamic models⁵. When two or more liquids are mixed there occur some changes in physical and thermodynamic properties because of free volume change, change in energy and change in molecular orientations. Derived thermodynamic and acoustical parameters like internal pressure, free volume and acoustic impedance are of considerable interest in understanding the intermolecular orientations in binary liquid mixtures⁶⁻⁸. Excess thermodynamic properties of mixtures are useful in the study of molecular orientations and arrangements⁹⁻¹⁰.

The scientific investigation in the field of ultrasound velocity measurement in 3-(2-methoxyphenoxy)-1,2-propanediol drug solution in various solvent system presented. This drug is used as treat coughs and congestion caused by the common cold, bronchitis, and other breathing illnesses infections. The acoustic properties of 3-(2-methoxyphenoxy)-1,2-propanediol have been studied in 20% Acetone-water and 20% Ethanol-water solutions at 308K.

Experimental :

All the studied chemicals which were purchased from S.D. fine chemicals Mumbai, India, with on purity of 99.6% were used as such without further purification solvents acetone and ethanol used in the present work were of AR grade. Densities, viscosities and ultrasonic velocities were measured at 30S K over a wide range of composition. The densities were determined by using specific gravity bottle by relative measurement method with accuracy $\pm 1 \times 10^{-5}$ gm./cm³. The viscosities were measured by precalibrated Ostwald type viscometer. Ultrasonic velocity measurements were made by using an ultrasonic interferometer (Mittal Enterprises, New Delhi) at a frequency of 2MHz with a tolerance of $\pm 0.005\%$. All the measurements were carried out at 30S K.

Acoustic parameters-

Acoustic parameters such as apparent molar compressibility (ϕ_c), apparent molar volume (ϕ_v), adiabatic compressibility (β_s), specific acoustic impedance (Z), intermolecular free length (L), Limiting apparent molar volume (ϕ_v^0) Limiting apparent molar compressibility (ϕ_c^0) were determined using following relations.

- Ultrasonic velocity $u = \lambda \nu$ ----- 1
- Adiabatic compressibility $\beta_s = 1/\rho^2 \nu^2$ ----- 2
- Apparent molar volume $\phi_v = 10^3(\rho_0 - \rho) / \rho_0 \rho$ ----- 3
- Apparent molar compressibility $\phi_c = 10^3(\beta_s - \beta_0) / \rho_0 \rho$ ----- 4
- Intermolecular free length $L_c = K(\beta_s)^{1/2}$ ----- 5
- Specific acoustic impedance $Z = \rho \nu$ ----- 6
- Limiting apparent molar volume $\phi_v^0 = \phi_v + K_1 C^{-1}$ ----- 7
- Limiting apparent molar compressibility $\phi_c^0 = \phi_c + K_2 C^{-1}$ ----- 8

Table no.1

Experimental Data of Density, Ultrasonic Velocity and Viscosity of 3-(2-methoxyphenoxy)-1,2-propanediol in different solvent at 30SK

Solvent	Conc. mol dm ⁻³	Density ρ kg m ⁻³	Ultrasonic Velocity (m/s)	Viscosity η (mPa.s)
20% Acetone-Water Medium	0.01	1002.51	1472.8	1.000
	0.05	1002.27	1470.8	1.000
	0.10	1002.03	1468.8	1.000
	0.15	1001.79	1466.8	1.000
	0.20	1001.55	1464.8	1.000
	0.25	1001.31	1462.8	1.000
20% Ethanol-Water Medium	0.01	982.1	1451.4	0.950
	0.05	981.6	1449.4	0.950
	0.10	981.1	1447.4	0.950
	0.15	980.6	1445.4	0.950
	0.20	980.1	1443.4	0.950
	0.25	979.6	1441.4	0.950

Table no.2

Variation of some acoustical parameters with concentration of 3-(2-methoxyphenoxy)-1,2-propanediol in different solvents at 30S K

Solvents	Conc. mol dm ⁻³	$\rho \times 10^3$ m ³ mol ⁻¹	$\phi_v \times 10^3$ m ³ mol ⁻¹	$\phi_c \times 10^3$ m ³ mol ⁻¹ Pa ⁻¹	$L_c \times 10^{11}$ (m)	$Z \times 10^3$ kg m ⁻² s ⁻¹
20% Acetone-Water Medium	0.01	4.605	0.99308	80.6208	1.2939	14.74542
	0.05	4.619	1.24135	83.0911	1.2960	14.72175
	0.10	4.632	1.65514	72.9857	1.2985	14.69945
	0.15	4.640	2.48272	65.7012	1.3003	14.68754
	0.20	4.642	4.06545	55.4575	1.3020	14.68017
	0.25	4.654	6.09308	420.8666	1.2994	14.67504
20% Ethanol-Water Medium	0.01	4.585	1.24135	124.7384	1.2912	14.70086
	0.05	4.599	1.65514	129.0791	1.2952	14.71051
	0.10	4.614	2.48272	137.5279	1.2953	14.74185
	0.15	4.624	4.06545	129.0747	1.2967	14.73547
	0.20	4.624	4.06545	129.0747	1.2967	14.73547
	0.25	4.624	4.06545	129.0747	1.2967	14.73547

Table-3

Limiting values of ϕ_v and ϕ_c along with slope (S_v & S_c) for 3-(2-methoxyphenoxy)-1,2-propanediol in different medium at 30SK temperature

Temp T (K)	Medium	Parameters			
		$\phi_v^0 \times 10^3$ m ³ mol ⁻¹	$\phi_c^0 \times 10^3$ m ³ mol ⁻¹ Pa ⁻¹	$S_v \times 10^3$ m ³ mol ⁻¹ Pa ⁻¹	$S_c \times 10^3$ m ³ mol ⁻¹ Pa ⁻¹
30K	20% Acetone-W	12.758	-23.33	628.1	185.7
	20% Ethanol-W	11.698	52.12	66.22	-1258.1

Table-4

A and β_s coefficient values at 30SK in different medium for 3-(2-methoxyphenoxy)-1,2-propanediol

Medium	Coefficient	Temp 30K
20% Acetone-Water medium	A	0.553
	β_s	-0.112
20% Ethanol-Water medium	A	0.428
	β_s	-0.119

Results and discussion:

Table 1 shows that ultrasonic velocity (u), density (ρ), and viscosity (η) increases with increase in concentration for all systems. The increase in ultrasonic velocity is due to decrease in intermolecular free length (L_c) as shown in table 2. This suggests that there is a strong interaction between 3-(2-methoxyphenoxy)-1,2-propanediol and solvent molecule. Adiabatic compressibility (β_s) is a measure of intermolecular association or repulsion calculated from the measured ultrasonic velocity (u) and density (ρ). Adiabatic compressibility is found to decrease with increase in concentration. Since adiabatic

compressibility is inversely related to the product of density and ultrasonic velocity based on this the compressibility is expected to decrease which has observed in the present case. When the sound waves travels through the solution, certain part of it travels through the medium and rest gets reflected by the ion¹¹ i.e. restriction for flow of sound velocity by the ions. The character that determines the restriction movement of sound waves is known as acoustic impedance (Z). It has been found that acoustic impedance increases with increase in concentration. The apparent molar compressibility (κ_a) explains the solute-solvent and solute-solute interactions in solution and was calculated by using the equation no. 4. The apparent molar volume (ϕ_v) is defined as the change in volume of solution for the added one mole of a particular component at constant temperature and pressure. It is thermodynamic property which helps in elucidating solvation behavior of electrolyte in solution. Apparent molar volume was evaluated from the density of solution and solvent.

It is evident from the table 3 that ϕ_v^0 values are negative for 20% Acetone-water and for 20% Ethanol-water ϕ_v^0 values are positive. The negative ϕ_v^0 values are suggest solute-solvent interaction whereas positive values are due to solute-solute interaction, is further confirmed by ϕ_v^0 values which are positive for both solution 20% Ethanol-water and for 20% Acetone-water of the drug. S_v is a measure of solute-solvent interaction. It is observed from the table 3 that S_v values are higher in 20% Acetone-water and low in 20% Ethanol-water solution. This confirms that in 20% Ethanol-water solution solute-solute interactions and in 20% Acetone-water solute-solvent interaction predominate.

The viscosity B-Co-efficient has been derived from Jones-Dole equation

$$(c > 0.1m) \eta_r - 1 / C^{1/2} = A + B C^{1/2}$$

Where η_r is the relative viscosity. A and B are the characteristics of the solute and solvent.

A is Falkenhagen coefficient represent the contributor from solute-solute interaction and B is Jones Dole coefficient known to depend on the size of the solute particle and on the interaction between solute and solvent.

They were obtained by a least-squares treatment as intercept and slope of the linear plot of $\eta_r - 1 / C^{1/2}$ Vs $C^{1/2}$. The graph for each system given linear straight line showing validity of Jones-Dole equation. The slope of straight line gives value of B co-efficient.

The viscosity A coefficient represent the ion-ion interactions and negative values have shown some physical significance. However negative A values have also been reported to be in other solvents in some studies.²²⁻²³

The large and small value of 'A' shows the stronger and weaker solute-solute interactions respectively. When solute is introduced into solvent of organic-water mixture it will interfere with the ordered structure of water in the solutes co-sphere. As only one solute is present so such variation in the values of A can be explained.

In the present study viscosity of liquid solutions increases with increase in concentration of drugs solution in 20% Acetone-water 20% ethanol-water mixture. The increase in viscosity with increase in concentration may be attributed to the increase in solute solvent interactions.

Viscosity B coefficients have been established to arise from ion-solvent interactions and are responsible for introducing order or disorder in the structure of the solvent. Solute with negative B Coefficient is characterized as structure breakers indicating weak solute-solvent interactions. Such type of results is also shown by Reddy et al.²⁴⁻²⁵

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