

STUDIES ON VISCOSITY, DENSITY AND REFRACTIVE INDEX OF 3-Ethyl-4-methyl N-[2-(4-{[(trans-4-methylcyclohexyl) carbamoyl] sulfamoyl}phenyl)ethyl]-2-oxo-2,5-dihydro-1H-pyrrole-1-carboxamide IN MIXED SOVENT at 300.15K

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Abstract: Refractive index, molar refractivities and molar polarizability constant of 3-ethyl-4-methyl-N-[2-[4-[(4-methylcyclohexyl)carbamoylsulfamoyl]ethyl]-5-ethyl-3-methyl-2-oxo-1H-pyrrol-1-yl]benzamide in 20% Methanol media at $303.15 \text{ K} \pm 0.10^\circ\text{C}$ temperature and different concentrations (0.625×10^{-3} to $10.0 \times 10^{-3} \text{ M}$). The values of molar refraction (R_m) and molar polarizability (α) constant are found to be decreased with decreasing concentration of solute in solvent. Viscosity coefficient (A, B) evaluate by using John-Dole equation. These parameters throw the light on the solute-solvent interaction and solute-solute interaction.

Key words: molar polarizability constant, Molar refractivities and Viscosity coefficient.

INTRODUCTION

The substituted heterocyclic compounds 3-ethyl-4-methyl-N-[2-[4-[(4-methylcyclohexyl)carbamoylsulfamoyl]ethyl]-5-ethyl-3-methyl-2-oxo-1H-pyrrol-1-yl]benzamide has an antidiabetic medication within the sulfonylurea class, primarily prescribed for the management of type 2 diabetes [1][2]. The refractive index is an important additive property of molecular structure of liquid. For pure hydrocarbon, one can get an idea of aromatic content of liquid using refractive index. The refractive index is the ratio of angle of incident to the angle of refraction and it depends on the temperature and wave length of light. The extent of refraction depends on -i) the relative concentration of atom or molecule ii) The structure of atom or molecule. So refractive index gives idea about geometry and structure of molecule. Refraction of light is additive property, but also depends on the structural arrangement of atom in molecule. This can some time be used to determine the structure of an unknown compound whose molecular formula is known.

Has been studied density and refractive index of binary liquid mixture Eucalyptol with Hydrocarbon at different temperature [3]. Have been studied refractivity properties of some homologous series such as n-ethanoate, methyl alkanoates, ethyl alkanoates etc. were measured in the temperature range from 298.15 to 333.150K [4]. The properties of liquid such as refractive index, ultrasonic velocity and viscosity of binary mixture are studied by many workers [5]. Has studied relationship of refractive index to mass density and consistency of the mixing rule use to calculate these two quantities of multicomponent mixture like ambient aerosols with the index-density relationship [6]. Has studied refractive indices of binary mixture of bromoalkane and non polar hydrocarbons, also studied molecular interaction between the components of binary mixtures [7]. An additive properties such as molar refractivity and molar polarizability constant of allopurinol, acenocoumarol, warfarin and amoxicillin in different media [8]. Ultrasonic velocity and viscosity of PEG-8000, PEG- study of acoustical properties, viscosity coefficient of substituted heterocyclic compounds under suitable condition. Number of researchers studied the molar refraction and viscosity coefficient for binary and tertiary system. [9-14]

The present work is undertaken to make the systematic study of above Glimepiride refractometrically and viscometry at 300.15 K.

EXPERIMENTAL

Above substituted heterocyclic compounds have most important. The solution of above compound is prepared in solvent like 20% Methanol by dissolving an appropriate amount by weight. For density measurement all the weight took on contech balance (0.001gm). The refractive index of solvent and solutions are measured at different (0.625×10^{-3} to 10×10^{-2}) by Abbe's refractometer having accuracy with ± 0.0001 unit. The constant temperature of the prism box is

maintained by circulating water from thermostat at 300.15 K. Refractometer was calibrated using glass test piece of known refractive index supplied with the instrument.

The molar refraction of solvent and solution are determined by using Lorentz-Lorentz equation.

$$R_M = \frac{(n^2-1)}{(n^2+2)} \times \frac{M}{d} = \frac{4}{3} \pi N \alpha \quad (1)$$

The entire viscosity data have been analyzed by using Jones-dole equation.

$$\eta_{sp} / \sqrt{c} = A + B \sqrt{c} \quad (2)$$

n- Refractive index M- Molecular weight d- Density of solution R_m - Molar refraction N- Avogadro's number α - Molar polarizability constant X_1 and X_2 Mole fraction of solvent and solute in solution.

The refractive index of solvent and solution at different concentrations are measured by Abbe refractometer. The calculated values of molar refractivities and polarizability constant shown in table-1. For viscosity measurement ostwald viscometer (10ml) was used. The flow time measured by using digital clock (0.01Sec).

However study of molar refractivity , molar polarizability constant and viscosity coefficients of substituted heterocyclic compounds such Glimepiride in 20% Methanol media at 303.15 K $\pm$ 0.10C temperature and different concentrations (0.625x10⁻³ to 10.0x 10⁻²M) under identical set of experimental condition. This could cover mini fold aspect of solute-solvent interactions scanty.

RESULTS AND DISCUSSION

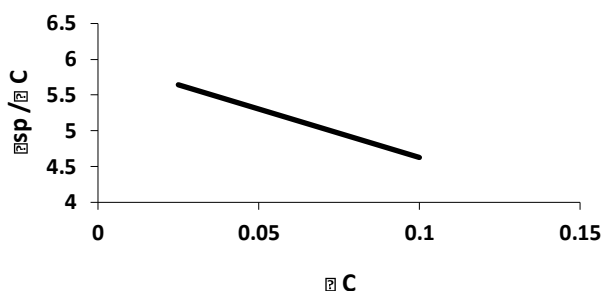
Table-1: The values of Molar polarization and polarizability constant at 300.15K.

Conc ⁿ in Moles/lit.	Density Kg/Cm ³	R.I.(n)	$R_m \times 10^{-5}$ cm ³ / mole	$\alpha \times 10^{-27}$ cm ³
10 x10 ⁻³	0.9794	1.3613	1.110008	4.4023
5 x10 ⁻³	0.9774	1.3576	0.551020	2.1853
2.5 x10 ⁻³	0.9744	1.3544	0.274133	1.0872
1.25 x10 ⁻³	0.9714	1.3510	0.136301	0.5406
0.625 x10 ⁻³	0.9684	1.3484	0.067905	0.2693

Table-2 - The values of η_r , Density, $\eta_{sp} / \sqrt{C}$, A and B-coefficient of substituted drug in 20% Ethanol.

Conc. Mol./lit	Density gm/Cm ³	Flow time (Sec)	η_r	$\eta_{sp} / \sqrt{C}$	A	B
10 x10 ⁻²	0.9794	219	1.47973	4.797297	5.9819	-13.543
5 x10 ⁻³	0.9774	199	1.344595	4.873303		
2.5 x10 ⁻³	0.9744	186	1.256757	5.135135		
1.25 x10 ⁻³	0.9714	176	1.189189	5.351078		
0.625 x10 ⁻³	0.9684	170	1.148649	5.945946		

Fig.1 Plots of $\frac{n_{sp}}{C}$ Vs C



From above tables (2), it could be seen that molar refractivity and molar polarizability constants decreases with decreasing the concentration of solution.

From Table-2 it is observed that the value of 'A' (Falkenhagen coefficient) are positive in all system studied. 'A' is measure of ionic interaction. The value of A is positive; it indicates that there is a strong solute-solute interaction in solute molecules. 'B' is Jones-Dole coefficient measures solute-solvent interaction. The value of "B" coefficient is negative in drug. Solute with negative 'B' coefficient is characterized as "Surface former" indicating weak solute-solvent interactions. This may be attributed to strong ion-solvent interaction in system.

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