

Study of Sonochemical in Mixture of Acetone with O – Benzo Quinone



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Abstract

Ultra sound velocity, density and viscosity are Experimently Determind in the mixture of acetone with o – Benzo quinone over entire range of composton at 303. ook and 308k These measerements have been used to evaluate some important parameter such as isentropic compressibility, specific acoustic impetonce, intermolecular free length, molar sound velocity, apparent molal compressibility, specific viscosity, reduced and relative viscosity and some other parameters for the same concentrations and temperature range. These parameters have been discussed in the terms of solute solvents interaction ons. The Ultrasound velacity increases on increasing concentration of solute.

Keywords: Somochemical Study, Acetone, O-Benzo Quinone

Introduction

Investigations of the physicochemical properties of the denatured state of o – benzo quinone are required to identify and characterize the contribution of various non-bonding interactions¹⁻² between different functional groups of o – benzoquinone and acetone – solvent interactions which are the driving force of covalent process the study of ultra sonic leads to a better understanding of the nature of interactions between the solute – solovent³ in recent years these have been as increasing interest in the study of non – electrolity components in equeous and non aqueous solutions⁴ The physical properties of dilute solutions of non – electrolyles depends that the solute in a Acetone structure breaker or maker the influence of small quantity of o – benzoquinone on hydrogen bonded structure of acetone in the solution in order to study the nature of molecular interactions in the above solutions the results have been discussed in tems of various interactons operating in these system.

Experimental

The Ultrasound velocities of the solutions were measured with a multi frequency ultrasonic interferometer of model f – 81.Quartz crystal have different frequency and have the accuracy of about $\neq 0.05\%$ the average of ten reading's treated as a final value of ultrasound velocity the temperature around the cell was controlled with in ± 0.01 by circulating the water from thermostat density measeerement were reformed with a pre collaborated bicapillary pycknometer. The accuracy of density measurements was up to ± 0.0001 gm/cm³ visosities were measured using an ubbelohde type capillary viscometer. Which was callibrated with doubly distilled water at three temperatures. The efflux time was measured using an electronic watch with a resolution of 0.01 see. An overaze of four the meassed viscoties was 0.005 cp The ultrasonic parameter such as molar sound relacity (R), Apparent molal compressibility (ϕ_k), insetropic pressibility, (β_s), specific aeoustic impedance (z), Specific viscosity (η_{sp}) and reduced viscosity, (η_{red}) were computed in there solutions using the deseribed standend formulae⁵.

Molar Sound Velocity (R)

$$R = M/\rho \quad v^{1/3} \dots\dots\dots(i).$$

Apparent molal compressibility

$$\phi_k = (\rho_0 \beta_s - \beta_{s0} \rho) + \beta_{s0} M/\rho_0 \dots\dots\dots(2)$$

$$\beta_s = \frac{1000}{C \times \rho_0} \dots\dots\dots(3)$$

$$Z = V \cdot \rho \dots\dots\dots v^2 \cdot \rho \dots\dots\dots(4).$$

$$\text{Relative association } R = M/\rho \dots \times 1/3 \dots\dots\dots(5).$$

Where $M = n_1 m_1 + n_2 m_2$ is ettctive molecular, weight. Where n_1 and n_2 are the number of moles of solvent and solute and m_1 and m_2 are their molecular weight respectivity. P and P₀ are there donsities of solution and solvent and V and V₀ are their U.V respectivity. β_s and β_{s0}

are the isentropic compressibility of solution and solvent and c is the concentration of solute in mole/litre.

The present work the chemical *o* – benzo quinone and Acetone used of AR (BHD) quality.

Aim of the Study

The Ultra Sound Study of Somochemical in mixture Acetone with *O*-Benzo Quinone. The Spectral study also have been carried out by I R Spectroscopy.

Result and Discussion

It can be seen from there data that the ultrasound velocity of solutions increases with increasing the concentration of *o* – benzo quinone in the solution of acetone at different - temperature. The ultrasound velocities of some sample decreases on increming temperature this trend of variation of ultrasound velocity also obseved for this solute.

The viscosity and reduced viscosity are given for the sample of *o* – benzo quinone with acetone solvent is tabulated in table and the value of molar sound velocity and isentropic compressibility is given in the table and the value of specific acoustic impedance and appaent molar compressibility (ϕ_k) are given in the table.

The viscosity of solutions for the given sample inceases with increase in the concentration of *o*-benzo quinone. The variation of density, viscosities of the present work also agreed with singh at.al.6 The

linear results of desity and ultrasound velocity have also been reported for glucose in binary aqueous solution of mannose maltose and raffinose by pal A and chauhan N 7 Linear result for molar volume for the electrolyte in aqueous solution of alcohols also shown by parmer M Lard guleria M.K and others8-9

Molar Cone. <i>O</i> - benzo. Quinone (mole/Lit)	Viscosity		Specific viscosity	
	303.K	308K	303K	308K
0.0216	0.3175	0.2804	0.0037	0.0042
0.0432	0.3196	0.2825	0.0105	0.0119
0.0649	0.3218	0.2847	0.0174	0.0197
0.0865	0.3240	0.2869	0.0243	0.0275
0.1081	0.3261	0.2890	0.0311	0.0353
0.1297	0.3283	0.2912	0.0380	0.0430
0.1513	0.3304	0.2933	0.0448	0.0508
0.1730	0.3326	0.2955	0.0517	0.0586
0.1946	0.3348	0.2977	0.0585	0.0664
0.2162	0.3369	0.2998	0.0654	0.0741

the results are also agreed with kumar P. et. Al10 the physicochemical behaviour of liquid mixture is also studied by R. Mehra and gaur A.K11 and the results agreed with the present work.

Molar Cone. <i>O</i> - benzo. Quinone (mole/Lit)	Ultrasound velocity		Density (gm/ml))	
	303.K	308K	303K	308K
0.0216	1162	1132	0.7796	0.7608
0.0432	1164	1134	0.7817	0.7629
0.0649	1166	1136	0.7839	0.7651
0.0865	1168	1138	0.7861	0.7673
0.1081	1170	1140	0.7882	0.7694
0.1297	1172	1142	0.7904	0.7716
0.1513	1174	1144	0.7925	0.7737
0.1730	1176	1146	0.7947	0.7759
0.1946	1178	1148	0.7969	0.7781
0.2162	1180	1150	0.7990	0.7802

Molar Cone. <i>O</i> - benzo. Quinone (mole/Lit)	Isentropic compressibility		Lowering compressibility	
	303.K	308K	303K	308K
0.0216	95.00	102.58	0.59	0.65
0.0432	94.42	101.93	1.17	1.30
0.0649	93.83	101.28	1.76	1.95
0.0865	93.25	100.64	2.34	2.59
0.1081	92.68	100.01	2.41	3.22
0.1297	92.11	99.38	3.48	3.85
0.1513	91.55	98.75	4.04	4.48
0.1730	90.99	98.14	4.60	5.09
0.1946	90.43	97.52	5.16	5.71
0.2162	89.88	96.91	5.71	6.32

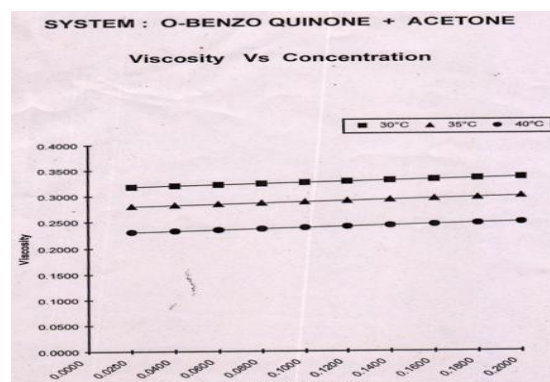
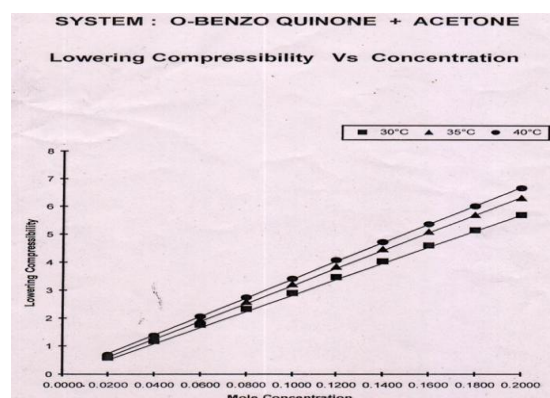
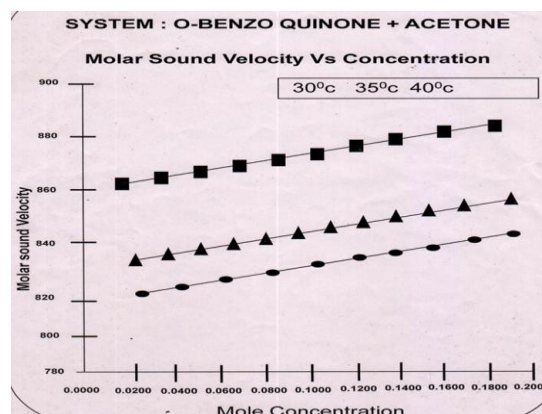
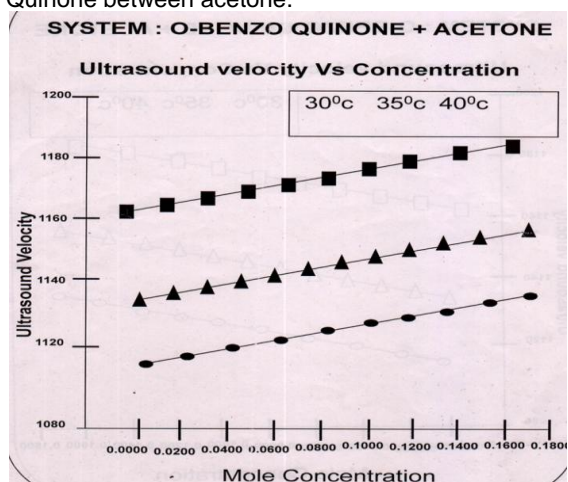
Depending on the polarity of side chain *o* – benzo Quinone vary in their hydrophilic or hydrophobic character. These properties are impotent

in *o* – benzo quinone stru stone and *o* – benzo quinone – Acetone interoetions. The importance of the

Molar Cone. O- benzo. Quinone (mole/Lit)	Reduce viscosity (η)		Molar sound velocity	
	303.K	308K	303K	308K
0.0216	0.1703	0.1930	862.77	1.0034
0.0432	0.2437	0.2762	865.66	1.0069
0.0649	0.2683	0.3041	868.56	1.0103
0.0865	0.2805	0.3179	871.45	1.0138
0.1081	0.2878	0.3262	874.34	1.0172
0.1297	0.2927	0.3317	877.24	1.0207
0.1513	0.2962	0.3357	880.13	1.0241
0.1730	0.2988	0.3387	883.05	1.0276
0.1946	0.3009	0.3410	885.95	1.0311
0.2162	0.3025	0.3428	888.85	1.0345

Molar Cone. O- benzo. Quinone (mole/Lit)	Solvation Number (Sn)		Apparent Molar volume (k)	
	303.K	308K	303K	308K
0.0216	0.1023	0.1051	10.7758	11.0518
0.0432	0.2048	0.2102	11.1302	11.4240
0.0649	0.3066	0.3146	11.4843	11.7958
0.0865	0.4074	0.4179	11.8349	12.1637
0.1081	0.5074	0.5203	12.1835	12.5296
0.1297	0.6065	0.6218	12.5302	12.8935
0.1513	0.7048	0.7225	12.8751	13.2553
0.1730	0.8025	0.8226	13.2196	13.6168
0.1946	0.8991	0.9215	13.5607	13.9745
0.2162	0.9950	0.0197	13.8999	14.3304

Side chain come from the influence, this has on the O – benzo quinone residues interaction with other structures, both with in a single o – benzo Quinone between acetone.



Conclusion

These measurements have been used to evaluate some important parameter such as isentropic compressibility, specific acoustic impetnce, etc.

Investigations of the physicochemical properties of the denatured state of O-benzoquinone are required to identify and characteristics the contribution of various non bonding interaction between different functional groups of O-benzoquinone and acetone solvent interaction which are the driving force of co-valent process. The study nature of molecular interactions in the above solutions the results.

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