

STUDIES IN ADDITIVE PROPERTIES SUCH AS MOLAR REFRACTIVITIES AND MOLAR POLARIZABILITY CONSTANT OF SOME SUBSTITUTED DRUGS WITH DIFFERENT CONCENTRATION IN VARIOUS MEDIA

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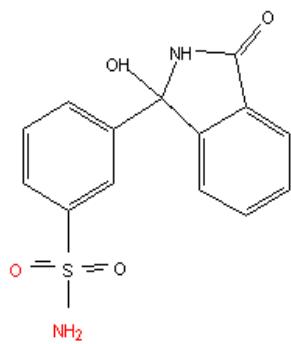
ABSTRACT: The additive properties such as Refractive index, Molar polarizability and molar refractivities of some heterocyclic drugs such as **Chlorothalidon** -2-chloro-5-(1-hydroxy-3-oxo-1,2-dihydroisoindol-1-yl)-benzenesulfonamide **Idoquinol** -5.7 di-iodo-8 – quinolinol have been studied in DMSO, DMF and THF media at $27 \pm 0.1^\circ\text{C}$ temperature and concentration.

The values of molar refractivity(R) and molar polarizability (α) are found to be decreased with decreasing the concentration of solute.

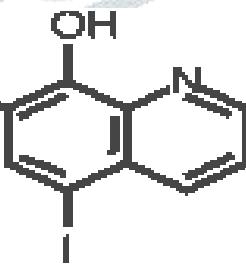
Key words: Additive property, **Chlorothalidon**, Refractive index, Molar polarizability and molar refractivities

INTRODUCTION- Refractive index is one of the important additives properties of liquid. When a ray of light passes from one medium to another, it suffers to refraction, that is a change of direction. If it passes from less dense to more dense medium, it is reflected towards the normal to form angle of refraction (r) which is less than angle incidence (i). The refractive index is the ratio of the velocity of light in vacuum to that in the medium and it depends upon the temperature and wave length of light. The properties of liquid such as refractive index, viscosity and ultrasonic velocity of binary mixtures are studied by many workers¹⁻³. Mahajan⁴ has studied molar refraction and polarizability constant of 2-amino-5-chloro benzene sulphonic acid in different percentage of dioxane-water mixture. Burghate⁵ and Agrawal⁶ studied the refractive indices in mixed solvents. Oswal et al⁷ have studied dielectric constants and refractive indices of binary mixtures. Ikhe⁸ has studies molar refraction and polarisability constant of pyrazoles and isoxazoles in different percentage of dioxane-water mixtures. Meshram et al³⁶ studied the molar refraction and polarizability constant of Al(III), Ce(III) and Fe(III) complexes with some substituted isoxazolines, pyrazole and pyrazoline refractometrically.

However study of molar refractivity and molar polarizability constant of novel compounds such as **Chlorothalidon**-2-chloro-5-(1-hydroxy-3-oxo-1,2-dihydroisoindol-1-yl)-benzenesulfonamide and **Idoquinol** -5.7 di-iodo-8 – quinolinol in non aqueous solvent such as THF,DMF and DMSO under identical set of experimental conditions which could cover manifold aspect of solute-solvent interaction is scanty.



Chlorothalidon



Idoquinol

Therefore the present work is undertaken to make the systematic study of above novel compounds refractometrically at 27°C temperature.

EXPERIMENTAL: Above novel compound are extensively used as drugs in pharmaceutical. These compounds provide the photographic material with good storage stability even at high temperature and high humidity. The compounds are synthesized by standard method and purity is checked by M.P, TLC, IR, and NMR.etc.

The solution of the compounds are prepared in different solvents (THF,DMF,DMSO etc.) by dissolving an appropriate amount by weight. All the weighing are made on Mechaniki Zactady Precyzyng Gdansk balance made in Poland ($\pm 0.001\text{gm}$).The accuracy of density measurements is within 0.1Kg/m^3 .

The refractive index of solvent and solutions are measure at different concentrations by Abbe's refractometer having accuracy with ($\pm 0.01\text{unit}$) .The temperature of prism box maintained constant by circulating water form thermostat at $27 \pm 0.1^\circ\text{C}$.Refractometer is initially calibrated with glass piece ($n=1.5220$) provided with the instrument.

The molar refraction of solvent and solution mixtures are determined from

$$\begin{aligned} R_m &= \left(\frac{n^2+1}{n^2-1} \right) \frac{M}{d} \\ R_m &_{(\text{solute})} = X_1 R_m_1 + X_2 R_m_2 \end{aligned}$$

R_m → molar refraction n → refractive index, d → density of solution,

No → Avogadro's number, α → polarizability constant,

$R_m 1$ & $R_m 2$ → molar refractivity of solvent and solute and

X_1 & X_2 → mole fraction of solvent and solute in solution.

The molar refraction represents actual or true volume of the substances molecules in mole. The molar refraction of solute can be calculated as:

$$R_m (\text{solute}) = R (\text{mixture}) - R (\text{solvent})$$

The refractive index of solvent and solution at different concentration are measured from Abbe's refractometer and the values of molar refraction and polarizability constants are evaluated and presented in tables 1 to 4 for different systems.

Table

Molar refraction and polarizability constant for Ligand 1 (Chlorothalidon) in DMF

Molarity (M) in moles/liter	R.I.	$R_m (\text{cm}^3 \text{ mole}^{-1})$	$\square \square \times 10^{-23} \text{ cm}^3$
0.01	1.321	0.1499	0.005975
0.005	1.320	0.07537	0.002903
0.0025	1.318	0.01637	0.0006309
0.00125	1.316	0.009546	0.0003677
0.000625	1.315	0.004794	0.0001847

Molar refraction and polarizability constant for Ligand 1 (Chlorothalidon) in THF

Molarity (M)	R.I.	$R_m (\text{cm}^3 \text{ mole}^{-1})$	$\square \square \times 10^{-23} \text{ cm}^3$
0.01	1.467	0.1207	0.004654
0.005	1.466	0.06144	0.002367
0.0025	1.464	0.03114	0.001199
0.00125	1.463	0.01573	0.000606
0.000625	1.462	0.007430	0.0003055

Molar refraction and polarizability constant for Ligand 1 (Chlorothalidon) in DMSO

Molarity (M)	R.I.	$R_m (\text{cm}^3 \text{ mole}^{-1})$	$\square \square \times 10^{-23} \text{ cm}^3$
0.01	1.360	0.1611	0.006207
0.005	1.359	0.08118	0.003127
0.0025	1.357	0.04095	0.001577
0.00125	1.355	0.02073	0.0007986
0.000625	1.354	0.01043	0.00040199

Molar refraction and polarizability constant for Ligand2(Iodoquinol) in DMF

Molarity (M)	R.I.	$R_m (\text{cm}^3 \text{ mole}^{-1})$	$\square \square \times 10^{-23} \text{ cm}^3$
0.01	1.421	0.1243	0.004788
0.005	1.420	0.06262	0.002412

0.0025	1.419	0.03142	0.001210
0.00125	1.418	0.01584	0.0006102
0.000625	1.417	0.002670	0.000102

Molar refraction and polarizability constant for Ligand 2(Iodoquinol) in THF

Molarity (M)	R.I.	Rm ($\text{cm}^3 \text{mole}^{-1}$)	$\alpha \times 10^{-2} \text{ cm}^3$
0.01	1.468	0.08094	0.003118
0.005	1.467	0.04058	0.001563
0.0025	1.467	0.02003	0.0007827
0.00125	1.466	0.01018	0.0003940
0.000625	1.464	0.005201	0.00006570

Molar refraction and polarizability constant for Ligand 2(Iodoquinol) in DMSO

Molarity (M)	R.I.	Rm ($\text{cm}^3 \text{mole}^{-1}$)	$\alpha \times 10^{-23} \text{ cm}^3$
0.01	1.363	0.08584	0.003306
0.005	1.362	0.04517	0.001740
0.0025	1.360	0.02286	0.0008807
0.00125	1.358	0.01153	0.0004442
0.000625	1.356	0.005805	0.0002236

Result and discussion: it could be seen from above tables (1 to 4) that molar refractivity and polarizability constants decrease with decreasing the concentration of solution. It is also observed that the values of Rm and α are found to be grater in polar solvents, .THF and DMF. Polar solvents involve H-bonding, may form complex with solute and non polar solvent does not involve H-bonding and does not associate with solute. This may also be attribute to the fact that the dipole in the compound lies perpendicular to the longer axis of the molecules considerable dipole association (inter molecular attraction) takes place which would be accompanied by increase in polarizability constants (α) as well as molar refractions (Rm) with increasing the concentration because of mutual compensation of the dipoles.

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