# APPENDIX -A

# Page No.40 in Chapter Two :

Reijen - gave a quantiative treatment on the role of defects on the sintering rate. He has studied the micreostructure of the sintering of Spinel Ferrites with an excess of  $Fe_20_3$  or deficiency of  $Fe_20_3$ . As in the spinel oxygen vacancies in an optimum concentration have to be present to give higher sintering rate. This condition is full filled in spenels that are deficit in  $Fe_20_3$ . Spinels with excess  $Fe_20_3$  which are swintered in an attemosphere sufficiently exidising to have cation vacancies will have a very low concentration of oxygen vacancies. The sintering rate is then reduced strongly and competing Mechanism such as pore growth can be predominent.

- 1. P.J.L. Reijen Sci. coram 4 (1968) 169
- P.J.L. Reijen VI Sym. Reactivity of solids schenactady (Aug. 1968)

### Page- No. 68 in chapter Two :

Fig 2.4 mark X' indicates observed points on the figure where the spinel structure has been confirmed for particular composition and temperature. '0' are the points on the same figure, that indicate the composition and temperature of final product containing (spinel + Fe<sub>2</sub>0<sub>3</sub> + CuFe<sub>2</sub>0<sub>2</sub>). The amount of phases will vary with composition and reaction temperature.

2) 
$$\cos(\sqrt{x}) = \frac{n_B + 5(1 - x)}{6 + x}$$

where  $x = \text{composition of } \mathbf{C}u$  ion  $n_B = \text{Magnetic moment}$ 

-2-

$$\propto$$
 is Y, K angles.

Refer Page No. 137 in Chapter -IV : Details of calculation of  $n_B$  : Emu/gm of Ni = 53.34 on Oscilloscope Division of Ni = 14

$$N = \frac{Emu/gm \text{ of Ni}}{Division \text{ of Ni}} = \frac{53.34}{14} = 3.81$$

$$O_{S} = \frac{N \times No. \text{ of divisions}}{Mass \text{ of samples}} \quad (Oscilloscope.)$$

$$MS = \frac{(1-P)O_{S} \times d_{S}}{S}$$

$$P = \frac{5.1 - d_{S}}{5.1} \quad \text{where p is porosity.}$$

$$(1-P) = 1 - \frac{5.1 - d_{S}}{5.1}$$

$$= \frac{d_{S}}{5.1}$$

$$n_{B} = \frac{MS \times Molecular \text{ weight.}}{5585 \times d_{S}}$$

$$n_{B} = \frac{(1-P)O_{S} \times Molecular \text{ weight}}{5585}$$

### Page 95, 96, 97 in Chapter - Three :

Ghani explained these breaked on the basis of phase transition. CuFe<sub>2</sub>0<sub>4</sub> undergoes from tetragonal to cubic phase at this temperature. This temperature will vary respect to the tetragonality ratio.

The condifuction mechanism in Cu-Ni ferrite was Comparison of two breaks in conductivity temperature curves. He exaplinate ed the conduction mechanism in these three regions is as follows :-

 Region-I. Due to impurities or impurity phases.
 Region-II. Due to phase transition (For CuFe<sub>2</sub>0<sub>4</sub> tetragonal to cubic

3. Region-III: Thermally .activated hopping conduction. In Region-II and III the conduction mechanism is mainly due to hopping . (i.e. hopping of polaron.) It is also ution observed that the activation energy is more for polarons than electrons. The observation of high activation energy is attributed to polaron conduction mechanism. Sawant et al have suggested the activation energy  $\mathbf{x}$ .  $\hat{\mathbf{E}}_{i}$  0.2 is due to electron conduction.

> (G.K. Joshi , A.Y. Khot, S.R. Sawant) J.Mat. Sci. (G.B) 22 (1987) 1694)

The conductivity in  $CoFe_2^{04}$  is mainly due to the formation  $Co^+3$ . The activation energy for first region is minimum and for III Region it is maximum. The changes in activation energies in second region have been found to be

sensitive to addition of CO. Applying concept of Elwell and Disxon the results, i.e. activation energy in II to III region is mainly due to thermally activated hopping processes.

(D.Elevell A.Dixon solid State Comm. 6 (1968) 585)

In this case temperature dependence of conductivity is mainly determined by temperature variation in the mobility of charge carriers. In such case the charge carriers are strongly localised on  $Fe^{+2}$  cations. The localisation may be attributed to electron-phonon inteaction (formation of Polarons) or to the strong exchange interaction between carriers and the magnetic sub-lattice (Magenetic polarons). An additional localisation of electrons at  $Fe^{+2}$  may arise from the inhomogenity, distributions of ions over octahedral sites (Ref. N. Rezlescs, D. Condurache P.Pertrawiu, E. Luca, J.Am. Ceran Soc. 57 (1974) 40)

The Curie temperatures measured and observed from resistivity plots generally differ. Parker et al have observed similar behaviou. and the difference is about 70°C at higher concentrations of Fe in Ni e ferrite system (Ref. S.A. Patil, Ph.D.Thesis, Shivaji University Kolhapur) (P. 164 (1980).

# Page No. 130 in Chapter -IV :

Effect of quenching temperature magnetisation a re studied on these samples. Table 1,2,3, the values of  $n_B$ , Tc are given. It is observed  $n_B$  and  $T_c$  values increase with increase of quenching temperature.Patil <sup>(1)</sup> have observed similar behaviour in  $CuFe_2^{0}a_4$  samples. He has attributed this on the basis of transfer of Cu ions from B site to A site and amount of transfer

-3-

ion increases w-ith increase of quenching temperature.

If Cu Co Ferrite the Similar behaviour may occur, where Cu ions transfer from B site to A site and Fe and Co ion from A to B site. The cation distribution shown in these tables clearly indicates the more and more cu ion is on A site at higher quenching temperature.

The phases like  $CO_3O_4$  and  $CuFe_2O_4$  reduce their amount. (1. S.A. Patil Ph.D.Thesis:

study of physical properties of Cu Fe  $3-x^0$  ferrites, Shivaji university Kolhapur (1980) )