1% with a magnetic balance. Electrical resistance was measured by a standard fourprobe method in the range from 77 K to 700 K with current reversal. The current leads and potential probes were spot-welded on the specimens with manganin wires of 0.2 mm in diameter.

## §3. Results

Figure 1 shows, as an example, the magnetization curves along the direction parallel and perpendiculer to the c axis at 4.2 K for Fe<sub>2</sub>P and (Fe<sub>0.8</sub>Ni<sub>0.2</sub>)<sub>2</sub>P compounds. The easy direction of magnetization was the c axis and no magnetocrystalline anisotropy in the basal plane was observed for all the compounds. The spontaneous magnetization  $\sigma_s$ was determined by extrapolating the linear part of the c axis magnetization back to zero field. The temperature denendence of  $\sigma_{\rm e}$ is shown in Fig. 2 for  $(Fe_{1-x}Ni_x)_2P$  compounds. The transition from ferromagnetic to paramagnetic at  $T_c$  for all the compounds shows a second order transition except for Fe<sub>2</sub>P, where the transition at  $T_{\rm c}$  is of the first order accompanying by the deformation of the lattice.

Figure 3 shows the inverse susceptibility  $\chi_g^{-1}$  vs temperature curves for the compounds. Here, no correction for the temperature independent susceptibility was made except for the compound with x=0.7, for which the  $\chi_g$  in Fig. 3 is the estimated value by sustraction of the  $\chi_g$  for Ni<sub>2</sub>P from the measured one. The  $\chi_g$  for the compounds in the range of  $0 \le x \le 0.7$  well obeys the Curie-Weiss law



Fig. 1. Magnetization curves along the direction parallel and perpendicular to the c axis at 4.2 K for x=0 and x=0.2 in  $(Fe_{1-x}Ni_x)_2P$ .



Fig. 2. Spontaneous magnetization per gram  $\sigma_s$  plotted against temperature for  $(Fe_{1-s}Ni_s)_2P$ .



Fig. 3. Inverse susceptibility per gram  $\chi_g^{=1}$  plotted against temperature for  $(Fe_{1-x}Ni_x)_2P$ .

except in the vicinity of  $T_c$ . The compounds with  $x \ge 0.8$  show the Pauli paramagnetism and the value of  $\chi_g$  was  $3.2 \times 10^{-6}$  (emu/g) independent of x.

Figure 4 shows the concentration dependence of the magnetic moment  $\bar{\mu}$  at 4.2 K and the effective numbers of Bohr magneton  $\bar{n}_{eff}$ per metal atom. Both  $\bar{\mu}$  and  $\bar{n}_{eff}$  decrease monotonically with increasing x and become zero at about x=0.7. The ferromagnetic Curie temperature  $T_c$  and the paramagnetic Curie temperature  $\theta_p$  are plotted against the concentration of Ni in Fig. 5. The value of  $\theta_p$  decreases monotonically with increasing x. On the other hand, the curve of  $T_c$  takes a maximum at about x=0.1 in agreement with the result reported by Fruchart *et al.*<sup>3)</sup>

Figure 6 shows the temperature dependence of the electrical resistivity along the c axis for  $(Fe_{1-x}Ni_x)_2P$  compounds. In the figure, the data of Ni<sub>2</sub>P are those obtained by employ-



Fig. 4. Concentration dependence of the mean magnetic moment at 4.2 K p and the mean effective numbers of Bohr magneton  $\bar{n}_{eff}$  per metal atom for  $(Fe_{1-x}Ni_x)_2P$ .

0.5

X

1.0



Fig. 5. The ferromagnetic and paramagnetic Curie temperatures,  $T_c$  and  $\theta_P$ , respectively, plotted against composition for  $(Fe_{1-x}Ni_x)_2P$ .

ing the polycrystal specimen since no single crystal was available. The resistivities of Fe rich compounds increase with increasing temperature at low temperatures which might be



Fig. 6. Temperature dependence of the electrical resistivity along the c axis for  $(Fe_{1-x}Ni_x)_2P$ .

associated with spin disorder, but decrease with increasing temperature at higher temperatures in paramagnetic state. On the other hand, the resistivities of the compounds with  $x \ge 0.8$  increase with increasing temperature at all temperature ranges. These results might be attributable to the hole in the narrow d electron band. It may be noted that the jump of the resistivity due to the first order transition observed in Fe<sub>2</sub>P disappears already by substituting only 8% of Fe by Ni.

## §4. Discussion

For the hexagonal crystal in ferromagnetic state, the magnetocrystalline anisotropy is conventionally expressed by

$$E = K_1 \sin^2 \theta + K_2 \sin^4 \theta + \cdots, \qquad (1)$$

where  $K_1$  and  $K_2$  are the first and second order anisotropy constants, respectively, and  $\theta$  is the angle between the magnetization and the *c* axis. When a magnetic field is applied in the basal plane for the sample with the easy direction of magnetization in the *c* axis, the following equation can be obtained from the minimum condition of the free energy,<sup>4)</sup>

$$\frac{H}{M} = \frac{2K_1}{M_s^2} + \left(\frac{4K_2}{M_s^4}\right) M^2 + \cdots$$
 (2)

Here,  $M_s$  is the saturation magnetization and M is the magnetization along the direction perpendicular to the c axis at the effective field H. Since M is proportional to H as

98

Metal atom ( $\mu_{B}$ )

per

Magnetic Moment

Mean

3.0

2.0

1.0

0L 0 (Vol. 44,