Volume 29A, number 8

30 June 1969

FEB 23 1970

EFFECT OF PRESSURE ON T_c AND BAND STRUCTURE IN TRANSITION METAL ALLOYS

W. GEY and D. KÖHNLEIN Physikalisches Institut der Universität Karlsruhe, 75 Karlsruhe, Germany

Received 9 May 1969

Measurements on ten alloys of the series Zr-Nb-Mo show a close correlation of dT_c/dp with the d-band structure of these alloys and suggest that the d-band is virtually rigid with respect to pressure.

In an attempt to detect some unifying principle for the complex reaction of superconductivity of transition metals and their alloys to application of hydrostatic pressure, the alloy series Zr-Nb-Mo was chosen on the following grounds : 1) The series is homologous in structure [1]. 2) Data on the electronic specific heat coefficient γ and θ_D exist [2]. 3) Presumably no effects of electronmagnon interactions complicate the situation.

The main result is that for all 10 alloys the variation of $T_{\rm C}$ with pressure is strongly influenced by the shape of the d-band.

Pressure up to 75 kbar was generated in a piston-cylinder type cell [3]. At first sight, the results are rather complex. For Nb and Nb-rich alloys a pronounced kink near 20 kbar occurs in the otherwise $T_{\rm C}(p)$ behaviour. For Nb-Zr alloys with 20, 40 and 60 at. % Zr and for Nb73Mo27 no kink is observed; $T_{\rm C}(p)$ is nearly linear for all pressures.

The results on alloys are collected in fig.1, to-

gether with other relevant parameters. The occurence of the kink in $T_c(p)$ for the Nb-rich alloys is represented by splitting of $d \ln T_c/dp$ into two branches for p < 20 kbar and p > 25 kbar, respectively. We discuss here the results for p > 25 kbar where the slope d ln T_c/dp decreases monotonically with increasing number n of valence electrons per atom at zero pressure*. One notes that $d \ln T_c/dp$ changes sign at a concentration near that at which $T_{c}(n)$ has a maximum. Since in these alloys $T_{c}(n)$ is closely related to the d-band structure, this suggests that $T_{c}(p)$ is likewise connected with the d-band. Analyzing the data in this respect it is found that a best data fit is obtained by the simple relation

* *n* is thus defined by the chemical composition. It is not to be confused with the actual electron concentration which may be pressure dependent.

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 $\partial \ln T_c / \partial p = a \left(\partial \ln T_c / \partial n - \partial \ln \theta / \partial n \right) + b$ (1)

with $a = 0.40 \times 10^{-6} \text{ bar}^{-1}$ and $b = 0.62 \times 10^{-6} \text{ bar}^{-1}$, θ = Debye temperature. Substitution of the BSC expression for T_c in eq. (1) yields

$$\partial \ln \theta / \partial p + (NV)^{-1} \cdot \partial \ln NV / \partial p =$$
 (2)

 $= a (NV)^{-1} \cdot \partial \ln NV / \partial n + b$ V = V_{BCS}. (We note that comparison of the detailed shape of $\theta(n)$ and $T_c(n)$ near n = 5 makes

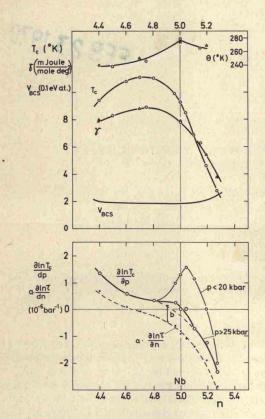


Fig. 1. Lower part: measured slopes $\partial \ln T_C / \partial p$. For p > 25 kbar comparison is made with $a \cdot \partial \ln / \partial n \equiv \equiv a (\partial \ln T_C / \partial n - \partial \ln \theta / \partial n)$ (eq. (1) in the text). Relevant parameters are plotted in the upper part in the usual nomenclature. Data for θ and $\gamma \propto N$ were taken from ref. [2] and own measurements.

interpretation in terms of the more recent analysis of McMillan [4] less favourable). For niobium $\partial \ln \theta / \partial p = \gamma_{\rm G} \cdot \kappa$ can be calculated from the experimental data on $\gamma_{\rm G}$ (Grüneisen parameter) [5] and κ (compressibility) [6]. $\partial \ln \theta / \partial p$ becomes 0.65×10^{-6} bar⁻¹ which equals our empirical term b within experimental error. Under the reasonable assumption that $\partial \ln \theta / \partial p$ does not change much with composition, eq. (2) thus reduces to

$$\partial NV/\partial p = a \cdot \partial NV/\partial n \tag{3}$$

with the solution

$$NV = f(ap+n) \tag{4}$$

Thus we arrive at the result that for p > 25 kbar the shape of the function NV(n) remains unaffected by pressure, i.e. is rigid with respect to pressure.

Since in this alloy system V is a slowly varying function of n, the shape of NV(n) is mainly determined by N. One is led to assume that this is also true for NV(p), and to conclude from eq. (4) that the effect of the application of pressure is primarily a filling of the d-band. A simple mechanism which would provide this can be seen in a pressure-induced shifting of the broad s-band relative to the narrow d-band.

We wish to thank Professor W. Buckel for stimulating our interest in this problem and discussing it in all stages, Dr. G. Dummer for specific heat measurements on some of the samples, and Dr. H. Wühl for preparing most of the samples

References

- 1. C.W. Berghout, Phys. Letters 1 (1962) 292.
- F. Heininger, E. Bucher and J. Muller, Phys. kondens. Materie 5 (1966) 243.
- 3. W. Gey, Phys. Rev. 153 (1967) 422.
- 4. W. L. McMillan, Phys. Rev. 167 (1968) 331.
- 5. J.G. Collins and G.K. White, Progr. Low Temp. Phys. 4 (1964) 450.
- 6. P.W.Bridgman, The physics of high pressure, (London, G. Bell and Sons, Ltd., 1958) p.160.