

where n_v is the degeneracy factor defined in Ref. 4, and $a^* = \kappa_b \hbar^2 / m^* e^2$ is the effective Bohr radius.

The long-wavelength approximation (10a) to the static dielectric constant leads to the same result as does the Thomas-Fermi approach to the screening,⁷ with a potential whose asymptotic form at large r is $\varphi \sim Zes(1+sd)/\kappa_b(sr)^3$.⁷ Because of the discontinuity in $d\kappa(q, 0)/dq$ at $q = 2k_F$, we must add to this an oscillatory term which can be evaluated from (9) and (10) using a theorem of Lighthill.⁸ The oscillatory term dominates at large r , and has the asymptotic form

$$\varphi(r) \sim -\frac{Zes}{\kappa_b} \frac{4k_F^2 \exp(-2k_F r)}{(2k_F + s)^2} \times \left[\frac{\sin \xi}{\xi^2} + \frac{8^{1/2} s \cos(\xi - \pi/4)}{\pi^{1/2} (2k_F + s) \xi^{5/2}} + \dots \right], \quad (11)$$

where $\xi = 2k_F r$. The leading term in (11) has the same dependence on r as does the result of Roth, Zeiger, and Kaplan⁹ for a three-dimensional semiconductor with cylindrical energy bands.

The screened Coulomb potential due to external charges is discussed more fully in a forthcoming paper,⁷ where it is used to calculate bound states and ionized impurity scattering of the electrons confined to the plane.

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⁴We assume that the electrons occupy a single band, and ignore transitions to higher bands. If the occupied band has N electrons/cm², a spin degeneracy of 2, and an orbital degeneracy of n_v ($n_v = 1$ for a free-electron gas, and $n_v = 2$ for the results of Ref. 1), then the Fermi wave vector is $k_F = (2\pi N/n_v)^{1/2}$.

⁵The sum rule (4) applies to the model that led to (3), but does not have the generality of the corresponding three-dimensional result [see, for example, F. Stern, in *Solid State Physics*, edited by F. Seitz and D. Turnbull (Academic Press, Inc., New York, 1963), Vol. 15, p. 299, esp. Sec. 23]. In a physical system such as a thin metallic film or the inversion layer of a semiconductor, our results are valid only for excitation energies $\hbar\omega$ smaller than the separation between the two lowest electric sub-bands.

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⁷F. Stern and W. E. Howard, to be published.

⁸M. J. Lighthill, *Introduction to Fourier Analysis and Generalised Functions* (Cambridge University Press, Cambridge, England, 1958), p. 52.

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PHONON-DISPERSION MEASUREMENTS ON A KRYPTON SINGLE CRYSTAL*

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Phonon-dispersion relations for the symmetric [100], [110], and [111] branches in fcc krypton have been measured by triple-axis neutron spectrometry. Measurements were carried out at 79°K on a single-crystal sample grown from the melt at a pressure of 2.31 kbar.

The phonon-dispersion relations in fcc krypton have been measured on the triple-axis spectrometer at the Brookhaven high-flux-beam reactor. The single crystal used for the experiment was grown from the melt at a pressure of 2.31 kbar in an aluminum-alloy pres-

sure cell incorporating a nucleation tip at the bottom. The cell was cylindrically shaped with an inside diameter of 12 mm and an outside diameter of 47 mm. The growth process was carried out in a temperature-controlled Dewar with heaters appropriately placed so as to pre-

vent blockage of the high-pressure tubing and to maintain a suitable temperature gradient along the sample holder. After complete solidification the sample was annealed for three weeks at 166°K, i.e., about 10° below the 2.31-kbar melting temperature, and then was cooled to 79°K, the temperature at which the dispersion-curve data were taken. This process yielded a crystal with a measured lattice parameter of 5.725 ± 0.010 Å. On the basis of this lattice parameter, compressibility data, and previously reported values of the lattice parameter at 79°K and 1 atm,¹ it is estimated that the sample pressure decreased to about 0.3 kbar in cooling to 79°K.

The size of the krypton crystal, estimated by beam-masking experiments, was approximately $8 \times 8 \times 11$ mm³. The small size of the crystal, for an inelastic neutron-scattering experiment, has somewhat limited the accuracy obtainable in the present investigation. The quality of the crystal was very good, however, as indicated by the measurement of a Bragg rocking curve with a full width at half-maximum of only 7'. This high degree of perfection, which is of some interest in itself, is probably due not only to the annealing treatment, but also to the high-pressure growth technique. Note that in this method of crystal preparation, in which the solidified rare gas fills an entire cell at a high pressure, the high vapor pressure exhibited by such crystals near their melting temperatures generates no additional problem in containment of the crystal during a long-term, high-temperature annealing process. Moreover, it is felt that the reason the crystal was not damaged on cooling slowly from 166 down to 79°K is that the pressure on the sample never becomes negative in that interval.

This apparent absence of crystal damage on cooling a solidified rare gas is consistent with the observations of Peterson, Batchelder, and Simmons² on crystals of argon grown and handled in an essentially stress-free manner within a thin-walled Mylar tube, but is in contrast with the behavior of crystals of solidified rare gases grown in Inconel tubes by White and Woods³ at pressures near 1 atm. In the latter case, thermal conductivity data indicated the crystals to be quite imperfect. This could well have been due to crystal damage in pulling away from the tube, because of the large thermal contraction of the solidified gas relative to Inconel.

A positive external pressure at the solidus eliminates this effect, however, unless the sample is cooled down to a temperature at which the sample pressure passes through 0. It should be possible to grow krypton crystals at some higher pressure and cool them at constant density to 0°K with no danger of pressure reversal.

The measurements reported here were carried out by constant- \bar{Q} scan⁴ with several fixed incident neutron energies between 18 and 35 meV. Typical neutron groups corresponding to the momentum transfer at the zone boundaries gave neutron counts of approximately 3/min above a background of 2/min. Better signals were obtained for the neutron groups at smaller q values, however. The points in Figs. 1 and 2 show the positions of neutron groups observed on the [100], [110], and [111] branches at a temperature of 79°K. The error bars were estimated from the shapes of the neutron groups. The [110] axis of the crystal lay about 15° away from the Dewar axis, and this somewhat inconvenient orientation rendered it impossible to observe the transverse branch propagating along [110] polarized in the [110] direction. Since no data exist at present for the elastic stiffnesses of krypton, it is not possible to compare initial slopes of the dispersion curves with those associated with velocity of long-wavelength acoustic disturbances.

Calculations of the phonon-dispersion curves

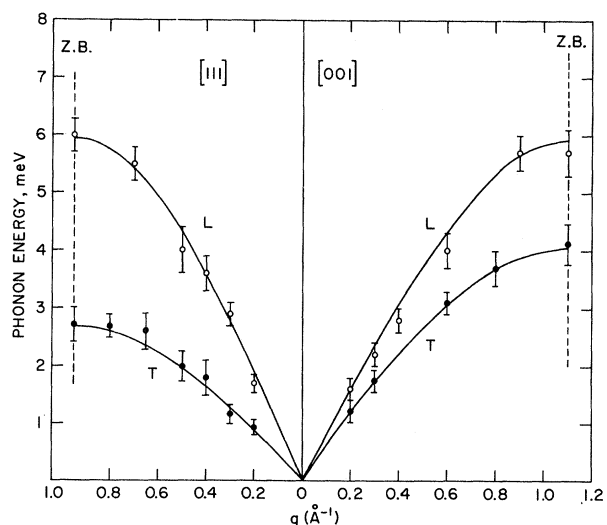


FIG. 1. Phonon-dispersion relations of Kr at 79°K along [001] and [111] directions. The conversion factor from meV to the frequency (10^{12} cps) is 0.242.

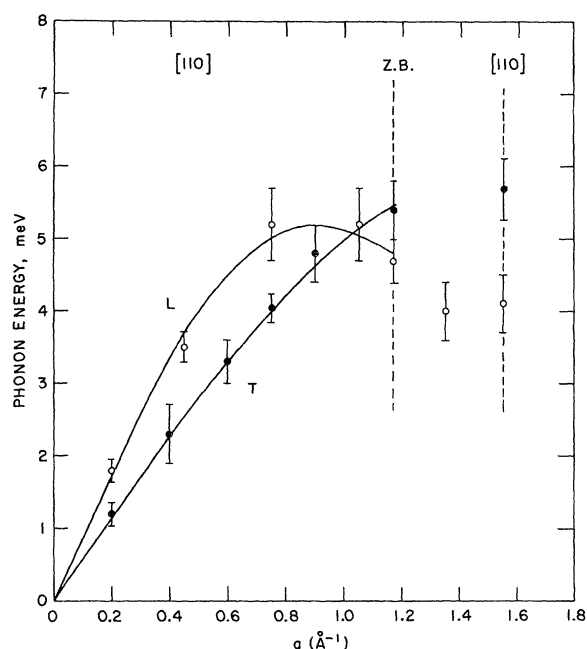


FIG. 2. Phonon-dispersion relations of Kr at 79°K along [110] direction. The transverse mode shown has its polarization vector in the [100] direction.

of krypton have been made by Grindlay and Howard,⁵ Horton and Leech,⁶ and Wallace.⁷ A meaningful comparison of calculated curves with the present experimental results will require a thorough analysis, including consideration of anharmonic effects as well as the detailed form of the potential function. However, a tentative comparison with the Grindlay and Howard results is shown by the solid curves of Figs. 1 and 2. Those curves were obtained by reducing the Grindlay and Howard calculated phonon energies by a factor of 0.92. Their dispersion curves were calculated in the harmonic approximation using a 6-12 Lennard-Jones potential with the potential constants evaluated for a lattice parameter of 5.674 Å. The atomic volume in the crystal used in the present experiment was 2.7% greater than that assumed in their calculations. Therefore, taking a value of $\gamma = 3$ for the Grüneisen parameter,⁶ a decrease of 8% in the phonon energies could be explained

by a quasiharmonic shift. However, agreement of the experimental results with a substantially harmonic theory based on the 6-12 Lennard-Jones potential may well be fortuitous.

The experimental work reported on krypton in this Letter is being continued with the following objectives: (1) to obtain improved values of the phonon energies proper, (2) to measure at constant volume the anharmonic shift of phonon energies with temperature, and (3) to measure the quasiharmonic volume dependencies of phonon energies. In carrying out (3), inelastic neutron-scattering studies are planned on crystals grown at a series of pressure points, exploiting thereby the change of density of the solid along the melting line. To achieve an acceptable level of accuracy in a reasonable length of experimental time, it will first be necessary to prepare larger single crystals than the one used in the present work.

A detailed analysis of this experiment and its results will be published later.

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