

vided by Barker's study of the TO phonon mode in GaP [A. S. Barker, Phys. Rev. 165, 917 (1968)]. The shape of the large-angle Raman line and IR absorption peak deviate strongly from the Lorentzian shape. Barker has carried out an analysis of his data by introducing a frequency-dependent damping factor. By studying the shape of the polariton spectrum of this mode, in principle one may extend Barker's analysis to a much wider range of frequencies. In essence, in the polariton regime, one can sweep the line center over a wide frequency region by varying the scattering angle.

⁵L. Van Hove, Phys. Rev. 95, 249 (1954); 95, 1374 (1954).

⁶J. J. Hopfield, Phys. Rev. 112, 1555 (1958).

⁷For example, see E. Burstein, S. Ushioda, A. Pinczuk, and J. F. Scott, 1969, paper A-1, in *Proceedings of the International Conference on Light Scattering Spectra of Solids, New York, 1968*, edited by G. B. Wright (Springer, New York, 1969).

⁸See the thesis of S. Ushioda cited in Ref. 1. Our expression for the scattering efficiency is one-half that given by Ushioda. We believe Ushioda's result is in error, because he presumed the cross section associated with both the Stokes and anti-Stokes line (in the classical theory) is proportional to the mean-square value $\langle u^2 \rangle_T$ of the *total* relative displacement of the ions in the unit cell. In fact, $u(t)$ contains a positive frequency component u_+ and a negative frequency part u_- , with (in the classical theory) $\langle u_+^2 \rangle_T = \langle u_-^2 \rangle_T = \frac{1}{2} \langle u^2 \rangle_T$. The cross section for anti-Stokes scattering is proportional to $\langle u_+^2 \rangle_T$, while that for Stokes scattering is proportional to $\langle u_-^2 \rangle_T$ rather than $\langle u^2 \rangle_T$. Thus, Ushioda's cross section should be reduced by a factor of 2. This will bring it into agreement with our result.

⁹D. T. F. Marple, J. Appl. Phys. 35, 539 (1964).

¹⁰B. Tell, S. P. S. Porto, and T. C. Damen, Phys. Rev. Letters 16, 450 (1966).

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COMMENTS AND ADDENDA

The Comments and Addenda section is for short communications which are not of such urgency as to justify publication in *Physical Review Letters* and are not appropriate for regular Articles. It includes only the following types of communications: (1) comments on papers previously published in *The Physical Review* or *Physical Review Letters*; (2) addenda to papers previously published in *The Physical Review* or *Physical Review Letters*, in which the additional information can be presented without the need for writing a complete article. Manuscripts intended for this section may be accompanied by a brief abstract for information-retrieval purposes. Accepted manuscripts will follow the same publication schedule as articles in this journal, and galley proofs will be sent to authors.

Cooling Curve in a One-Dimensional Crystal

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A region of $2N+1$ particles in a harmonic one-dimensional crystal is given temperature T_0 at time $t=0$. The central particle in this region cools down as follows¹:

$$T(t) = T_0 \sum_{n=-N}^N J_n^2(Nt), \quad (1)$$

where t is a suitable dimensionless time variable. By very complicated asymptotics Rubin¹ was able to determine the following limiting behavior:

$$\lim_{N \rightarrow \infty} T(t)/T_0 = 1, \quad \text{for } t < 1$$

$$= 2\pi^{-1} \arcsin(1/t), \quad \text{for } t > 1. \quad (2)$$

The present note demonstrates how this limiting cooling curve may be found in a simple manner. The standard representation²

$$J_n^2(z) = 2\pi^{-1} \int_0^{\pi/2} J_0(2z \sin \phi) \cos 2n\phi \, d\phi$$

allows the summation over n in Eq. (1) to be performed. Introducing a new variable $x = 2N \sin \phi$, we obtain

$$T(t)/T_0 = 2\pi^{-1} \int_0^{2N} dx \, x^{-1} J_0(xt) \times \sin[2N \arcsin(x/2N)] + J_N^2(Nt). \quad (3)$$

Taking formally the limit $N \rightarrow \infty$ in (3), we have

$$\lim_{N \rightarrow \infty} T(t)/T_0 = 2\pi^{-1} \int_0^\infty dx x^{-1} J_0(tx) \sin x,$$

equal to² the right-hand side of Eq. (2). The formal step involved is easily justified but details are omitted here.

¹R. J. Rubin, Phys. Rev. **131**, 964 (1963). See in particular his Appendix C.

²W. Magnus and F. Oberhettinger, *Formulas and*

Theorems for the Special Functions of Mathematical Physics (Chelsea Publishing Co., New York, 1949), pp. 32-34.

Specular Reflection in Al Films

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Experimental results on the size-affected resistivity of aluminum films are discussed.

We wish to comment on two conclusions reached by von Bassewitz and Mitchell¹ (BM) from their experimental results:

(a) The value of the mean free path l , for which they give a value of 17.5μ for polycrystal films as well as for single-crystal films.

(b) The reflection coefficient p , BM claim that the reflection at a smooth surface of a single-crystal film of Al is nearly specular.

1. DETERMINATION OF l

To obtain the mean free path l from the measurements of the size-affected resistivity ρ , BM plot $1/\rho t$ (t is the thickness of the film) as a function of $\ln t$ and draw a straight line through the experimental points. Extrapolation of the line of the axis where $1/\rho t = 0$ gives an intersection point $t_1 = 17.5 \mu$. BM conclude that the intersection point should be the point where $t_1/l = 1$. This leads to $l = 17.5 \mu$ and $\rho_\infty l = 11.5 \times 10^{-12} \Omega \text{ cm}^2$, where ρ_∞ is the bulk resistivity. To obtain this, BM use an approximation of the Fuchs-Sondheimer² (FS) theory for the case $t/l \ll 1$ which reads

$$\rho_\infty/\rho = (3t/4l) (1+p)/(1-p) \ln(l/t). \quad (1)$$

If $l = 17.5 \mu$, then the experiments of BM were done in a regime $0.01 < t/l < 0.2$. However, in this regime the extrapolation procedure described above is misleading. This is most easily seen by plotting $1/\rho_\infty t$ versus $\ln t$, where ρ_∞ is the (numerical) value of the full FS theory. When a straight line is drawn through the calculated points in the regime $0.01 \leq t/l \leq 0.2$, one finds that extrapolation to $1/\rho_\infty t = 0$ gives $t_1/l = 1.9$ if $p = 0$, $t_1/l = 1.3$ if $p = 0.1$, $t_1/l = 0.9$ if $p = 0.3$, and $t_1/l = 0.75$ if $p = 0.5$,

instead of $t_1/l = 1$ as follows from Eq. (1). These differences clearly show that if $l = 17.5 \mu$, Eq. (1) may not be used. The procedure used by BM to obtain $l = 17.5 \mu$ and thus $\rho_\infty l = 11.5 \times 10^{-12} \Omega \text{ cm}^2$ is, in fact, inconsistent with the FS theory.

2. DISCUSSION OF REFLECTION COEFFICIENT p

When concluding something about the reflection coefficient p from the FS theory, the dependence of l on t must be known. As BM have done no other experiments from which l can be determined, they had to make an assumption about the dependence of l on t . In fact, they reason in terms of a constant mean free path and then conclude from the fact that the slope of the straight line in the $1/\rho t$ -versus- $\ln t$ graph is different for single-crystal films and polycrystal films that $p \neq 0$. Using Eq. (1) they found that for smooth surfaces of single-crystal films $0.42 < p < 0.84$. From the discussion given above about the validity of Eq. (1) for the experiments of BM, it is clear that the obtained value of p seems not to be correct. The conclusion that $p \neq 0$ can, however, be right if indeed l is independent of t . It is clear, however, that one could also interpret the experiments by assuming $p = 0$ and letting l be a function of t , a possibility which is not at all unlikely in view of the results obtained by Cotti *et al.*³ Let us consider the consequences of such an interpretation; BM give that ρ varies from 8.5×10^{-8} to $1.1 \times 10^{-8} \Omega \text{ cm}$ for single-crystal films, and from 18×10^{-8} to $2.0 \times 10^{-8} \Omega \text{ cm}$ for polycrystal films for thicknesses between 0.2 and 3.6μ . Using the results of a two-band model,⁴ one finds that for single-crystal films, l varies from 6.1 to 18μ and from