HELICON PROPAGATION IN SODIUM AND COPPER

A Thesis Presented to The Faculty of the Department of Physics University of Houston

In Partial Fulfillment of the Requirements for the Degree Master of Science

ΒY

Stephanie Dietz Yesavage December, 1971

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ABSTRACT

Helicons and Gantmakher-Kaner oscillations were observed for the sodium metal. For a cylindrically symmetric Fermi surface approximation for copper, we observed the beating of two Gantmakher-Kaner waves for the [111] direction of copper and helicon propagation for the [001] direction of copper. Using the true Fermi surface of copper, we observed helicon propagation in the [001] direction. In all instances of helicon propagation, we found that a relaxation time of 10^{-10} sec. produced smooth sinusoidal shaped helicon curves, whereas a relaxation time of 10^{-9} sec. produced a delta-function shaped curve. TABLE OF CONTENTS

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INTRODUCTION

An electromagnetic wave striking a metal surface is generally damped to 1/e of its initial amplitude within a distance of δ , the skin depth. However, if a static magnetic field is present and the metal is pure, it is possible under certain conditions for the wave to propagate through the metal. This propagated wave is called a helicon.

Given a slab of metal of thickness Q upon which an electromagnetic wave of frequency ω is incident in the presence of a static magnetic field \vec{B} , we calculate in this paper the ratio of the transmitted to incident electric fields, E_t/E_i . The expression E_t/E_i is a complex quantity, thus supplying an amplitude and a phase factor.

The metals chosen for investigation were sodium and copper. The calculation of E_t/E_i requires the knowledge of the Fermi surface of the metal. Whereas, such calculations have been done for a spherical Fermi surface which sodium possesses and for approximate Fermi surfaces of copper, no calculation has yet been done using the exact Fermi surface of copper. It is the main purpose of this thesis to calculate E_t/E_i using the correct_Fermi surface of copper and to-thus make data available for comparison with experimental data recorded previously by others. A second purpose of this paper was an investigation of the effect of the relaxation time τ on the shape of the curve produced by a plot of E_t/E_i versus B.

The sodium metal was investigated because of the simplification of the calculations brought about by its spherical Fermi surface. After finding that the sodium data was in good agreement with experimental data, we advanced to a cylindrically symmetric Fermi surface for copper and, at last, to the true non-cylindrically symmetric Fermi surface. The former Fermi surfaces provide an accuracy check for the final copper calculations.

THEORY

Let us first describe the condition for helicon propagation. Suppose an electromagnetic wave with wave number \vec{q} and frequency ω is propagating in a metal in the presence of a static magnetic field \vec{B} . Suppose, also, that \vec{q} and \vec{B} are both parallel to the Z-axis. Now the force \vec{F}_{e} on an electron in the metal with an average velocity in the Z-direction, \vec{V}_{e} , due to the B field is found from

$$\vec{F}_e = e(\vec{v} \times \vec{B}) \cdot (1)$$

The electron will move in a helix about an axis parallel to the Z-axis due to the initial velocity of the electron. The cyclotron frequency of the electron is given by

$$\omega_c = (eB)/(mc)$$
 (2)

Now, the electron sees the electromagnetic wave as having the Doppler shifted frequency ω_e given by

$$\omega_e = \omega + q \, \overline{V_z} \, . \tag{3}$$

Let γ equal the relaxation time of the electron. If $\omega_c \gamma \gg 1$, then the electron makes many spirals of its orbit before it has a collision. In this time, it would be able to absorb energy from the electric field, if the condition $\omega_e = \omega_c$ is satisfied. This is called Doppler shifted cyclotron resonance. Thus, we have two required conditions for energy absorption,

$$\omega_e = \omega_c \quad , \tag{4}$$

and

$$\omega_c \tau \gg 1 \quad . \tag{5}$$

Let the maximum value of \overline{V}_{Z} of an electron in the metal be denoted by \overline{V}_{ZMAX} . Thus, the maximum Doppler shifted frequency seen by an electron is given by

$$\omega_{\text{EMAX}} = \omega + q \overline{V}_{\text{EMAX}} . \tag{6}$$

Now, for a given B field, ω_c is constant. Thus, if

 $\omega_c \langle \omega_{eMAX} \rangle$, then some electrons with a $\overline{V_z} \langle \overline{V_{zMAX}} will$ be able to satisfy the condition $\omega_c \omega_e$, and energy will be absorbed from the electromagnetic wave or helicon. If $\omega_c = \omega_{eMAX}$, only the electrons with $\overline{V_{z}} = \overline{V_{zMAX}}$ would be capable of absorbing energy. We call the condition $\omega_c = \omega_{eMAX}$ the absorption edge. Now, if we increase the value of B such that $\omega_c > \omega_{eMAX}$, then no electron could absorb energy, and the helicon would propagate with no -cyclotron damping effects.

For three-fold symmetry about the magnetic field direction, we have the following expression for E_t/E_c from Antoniewicz, et al.¹:

$$\frac{E_{\pm}}{E_{i}} = \frac{-i4\omega}{Q_{c}} \sum_{m=1}^{\infty} \frac{(-i)^{m}}{\left(\frac{m}{Q}\right)^{2} - \left(\frac{\omega}{c}\right)^{2}} + \frac{i4c}{Q\omega} + \frac{i4c}{Q\omega} + \frac{i4c}{Q\omega} , \quad (7)$$

where

$$\vec{\sigma} = \text{conductivity tensor} = \begin{pmatrix} \sigma_{xx} & \sigma_{xy} & \sigma_{xz} \\ \sigma_{yx} & \sigma_{yy} & \sigma_{yz} \\ \sigma_{zx} & \sigma_{zy} & \sigma_{zz} \end{pmatrix}, \quad (8)$$

$$\sigma_{+} = \sigma_{xx} - \iota \sigma_{xy} , \qquad (9)$$

and

$$\epsilon_{+}\left(\frac{m\pi}{Q}\right) = \frac{4\pi}{\omega} \tau_{+}\left(\frac{m\pi}{Q}\right) + K \qquad (10)$$

Now, $K \leftrightarrow \frac{4\pi}{\omega} \sigma_{+}$, and we ignore the term $\frac{14c}{Q \omega \varepsilon_{+}(o)}$ because of its negligible contribution. Thus, we have

$$\frac{E_{\pm}}{E_{i}} = -\frac{i4\omega}{QC} \sum_{m=1}^{d} \frac{(-1)^{m}}{(\frac{m}{Q})^{2} + \frac{i\omega 4\pi}{C^{2}} \Gamma_{\pm}(\frac{m\pi}{Q})}, \qquad (11)$$

and

$$\frac{E_{\pm}}{E_{c}} = -\frac{i4\omega}{Q_{c}} \sum_{m=1}^{\infty} \frac{(-i)^{mu}}{\left[\left(\frac{m\pi}{Q}\right)^{2} + \frac{\omega4\pi}{c^{2}} \tau_{xy}\right] + \frac{i\omega4\pi}{c^{2}} \tau_{xy}} \quad (12)$$

Now, let us denote the real and imaginary parts of $\sigma_{\bar{x}x}$ and $\sigma_{\bar{x}y}$ as follows:

Real
$$T_{XX}$$
 = ANSRX, (13)

Imaginary
$$\nabla_{XK} = ANSIX$$
, (14)

Real
$$\sigma_{xy} = ANSRY$$
, (15)

and Imaginary
$$\tau_{xy} = ANSIY$$
 (16)

Then we have,

$$\frac{E_{t}}{E_{c}} = -\frac{i4\omega}{QC} \sum_{m=1}^{\infty} \frac{(-1)^{mu}}{\left\{ \left(\frac{m\pi}{Q}\right)^{2} - \frac{\omega 4\pi}{C^{2}} \left[ANSIX - ANSRY \right] \right\} + i \left\{ \frac{\omega 4\pi}{C^{2}} \left[ANSRX + ANSIY \right] \right\}}$$
(17)

Let
$$S = \left\{ \left(\frac{m \pi}{Q} \right)^2 - \frac{\omega 4 \pi}{c^2} \left[ANSIX - ANSRY \right] \right\},$$
 (18)

and
$$V = \left\{ \frac{\omega 4T}{c^2} \left[ANSRX + ANSIY \right] \right\}$$
 (19)

Thus, we have for the real and imaginary parts of E_{t}/E_{i} ,

$$\operatorname{REAL}\left(\frac{E_{t}}{E_{t}}\right) = \frac{4\omega}{QC} \sum_{m=1}^{2} \frac{(-1)^{m+1}V}{S^{2}+V^{2}}, \qquad (20)$$

and

IMAGINARY
$$\left(\frac{E_{t}}{E_{c}}\right) = \frac{4W}{QC} \sum_{m=1}^{\infty} \frac{(-1)^{m+1}S}{S^{2}+V^{2}}$$
 (21)

Now, we have stated that E_{t}/E_{i} gives the amplitude, $|E_{t}/E_{i}|$, and the phase Θ of the E_{t}/E_{i} ratio. Thus, we have

$$\left|E_{L}/E_{i}\right| = \left(\left[\operatorname{ReA}\right]\left(\frac{E_{t}}{E_{i}}\right)\right]^{2} + \left[\operatorname{Imaginary}\left(\frac{E_{t}}{E_{i}}\right)\right]^{2}\right)^{\frac{1}{2}}, \qquad (22)$$

and

$$\operatorname{Arctan} \Theta = \frac{\operatorname{Imaginary}\left(\frac{E_{\pm}}{E_{\pm}}\right)}{\operatorname{ReAI}\left(\frac{E_{\pm}}{E_{\pm}}\right)} \quad (23)$$

However, we are concerned only with Real E_{\pm}/E_{c} , since this is the only part seen experimentally.

From the development of Mertsching², we have the following expressions:

$$\underbrace{\overrightarrow{U}}_{a,\lambda_r}(\overrightarrow{q}) = \underbrace{\underbrace{e^2}_{a\Pi f \lambda^2}}_{-KMAX} \underbrace{\underset{m=-\infty}{\overset{oo}{\text{m}_c}} \underbrace{\overrightarrow{V}_{ma} \, \overrightarrow{V}_{m\lambda_r}}_{m=-\infty} \underbrace{\underbrace{V}_{ma} \, \overrightarrow{V}_{m\lambda_r}}_{+i(\overrightarrow{q}, \overrightarrow{V} - \omega - m\omega_c)}, \quad (24)$$

and

$$\vec{V}_{m\alpha} = \frac{1}{2\pi} \int_{0}^{2\pi} \vec{V}_{\alpha}(\theta) e^{-\frac{1}{2\omega_{c}}} \int_{0}^{0} \left[\vec{q} \cdot \{ \vec{V}(\theta'') - \vec{V} \} + m\omega_{c} \right] d\theta'' d\theta , \qquad (25)$$

where

KMAX = maximum value of $\&_z$ on the Fermi surface,

 ω = frequency of electromagnetic wave,

1

 Υ = relaxation time,

 \hbar = Planck's constant,

e = charge of electron,

 $m_c = cyclotron mass,$

 ω_c = cyclotron frequency

Q =thickness of metal slab,

 $q = wave vector = \frac{m\pi}{Q}$ m = 0, 1, 2...,

m = an integer,

 α = indice for X, Y, or Z component,

k = indice for X, Y, or Z component,

 $\vec{V}_{\alpha}(e) =$ a-component of electron velocity at Fermi surface, and $\overline{\vec{v}}$ = average electron velocity at Fermi surface.

In our calculation of E_t/E_i , we need expressions for τ_{xx} and τ_{xy} . Thus, we have, assuming that $\frac{1}{2}$ is in the Z-direction,

$$\sigma_{XX} = \frac{e^2}{2\pi \hbar^2} \int_{-\kappa_{MAX}}^{+\kappa_{MAX}} dz \sum_{m=-\infty}^{\infty} \frac{\vec{v}_{mX} \cdot \vec{v}_{mX}}{\vec{\tau} + i(m\pi \nabla_z - \omega - m\omega_c)}, \quad (26)$$

$$\vec{\nabla}_{mx} = \frac{1}{2\pi} \int_{0}^{2\pi} V_{x}(\theta) e^{-\frac{1}{W_{c}} \int_{0}^{\infty} \left[\frac{m\pi}{Q} \left\{ V_{z}(\theta') - \overline{V_{z}} \right\} + mW_{c} \right] d\theta''} d\theta, \quad (27)$$

$$\overline{J}_{XY} = \frac{e^2}{2\pi \hbar^2} \int_{-KMAX} m_c d d z \sum_{M=-\infty}^{\infty} \frac{\overline{V}_{MX} \overline{V}_{MX}}{\frac{1}{T} + i(\frac{M}{Q} \overline{V}_z - \omega - M \omega_c)}, (28)$$

and
$$\nabla_{my}^{*} = \frac{1}{2\pi} \int_{0}^{2\pi} \nabla_{y}(\theta) e^{\frac{i}{\omega_{z}} \int_{0}^{\theta} \left[\frac{m\pi}{Q} \sum_{z} V_{z}(\theta'') - \overline{V_{z}} \right] + m\omega_{z} d\theta''} d\theta$$
. (29)

We have the following dispersion relation for a circularly polarized electromagnetic wave in a metal, such as a helicon:

$$q^{2} = \frac{4\pi\omega}{c^{2}} \left(IMAGINARY \, \sigma_{+} - iREAL \, \sigma_{+} \right) \quad . \tag{30}$$

Since

$$T_{+}$$
 = (ANSRX + ANSIY) + i (ANSIX - ANSRY), (31)

we have

$$q^{2} = \frac{4\pi\omega}{c^{2}} \left\{ (ANSIX - ANSRY) - i (ANSRX + ANSIY) \right\}.$$
 (32)

We may express $|q|^2$ in polar form as:

$$q^{2} = \left\{ \left(\frac{4\pi\omega}{c^{2}} \operatorname{Img}_{U} \overline{U}_{+} \right)^{2} + \left(\frac{4\pi\omega}{c^{2}} \operatorname{REAL} \overline{U}_{+} \right)^{2} \right\}^{2} \stackrel{i}{e} , \qquad (33)$$

where

$$\Theta = \arctan \frac{-\text{ReAl } T_+}{\text{Img. } T_+} = \arctan \frac{(-\text{ANSIX} + \text{ANSRY})}{(\text{ANSRX} + \text{ANSIY})}$$
 (34)

Let us call the roots of q^2 , q_1 and q_2 . Then we have that

$$q_{1} = \left\{ \left(\frac{4\pi\omega}{c^{2}} I_{mg} \nabla_{+} \right)^{2} + \left(\frac{4\pi\omega}{c^{2}} R_{eA} | \nabla_{+} \right)^{2} \right\}^{2} \left(\cos \frac{\Theta}{2} + i \sin \frac{\Theta}{2} \right), \quad (35)$$

and

$$g_{2} = \left\{ \left(\frac{4\pi\omega}{c^{2}} \operatorname{Img} \nabla_{+} \right)^{2} + \left(\frac{4\pi\omega}{c^{2}} \operatorname{R}_{eA} | \nabla_{+} \right)^{2} \right\}^{4} \left(\cos \left[\frac{\Theta}{2} + \pi \right] + i \sin \left[\frac{\Theta}{2} + \pi \right] \right). (36)$$

An electromagnetic wave propagates as $e^{i\left(\frac{\Theta}{2} + -\omega t\right)}$. If
 $\operatorname{Img} \nabla_{+} \gg \operatorname{R}_{eA} | \nabla_{+}, \text{ then } \Theta \text{ is close to zero. This makes}$
 g_{i} and g_{2} mostly real, and there is a negligible damping

factor. As the Real ∇_{+} increases, Θ approaches $-\frac{T}{2}$ which means that the imaginary part of $\frac{1}{2}$ is increasing which causes the damping factor to increase. At $\Theta: -\frac{T}{2}$, we have

$$q_{1} = \left\{ \left(\frac{4\pi\omega}{c^{2}} \operatorname{Img.} \nabla_{+} \right)^{2} + \left(\frac{4\pi\omega}{c^{2}} \operatorname{Real} \nabla_{+} \right)^{2} \right\}^{4} \left(\frac{1}{\sqrt{2}} - i \frac{1}{\sqrt{2}} \right), \quad (37)$$

and

$$q_{2} = \left\{ \left(\frac{4 \Pi \omega}{c^{2}} I_{mg} \sigma_{+} \right)^{2} + \left(\frac{4 \pi \omega}{c^{2}} R_{eA} | \sigma_{+} \right)^{2} \right\}^{4} \left(-\frac{1}{\sqrt{2}} + \frac{1}{\sqrt{2}} \right) .$$
(38)

The root q_1 is discarded because the imaginary part of q_1 carries a minus sign giving rise to an exponentially increasing wave. Thus, q_2 would be our solution.

Therefore, when helicon propagation exists, we will expect the condition $\operatorname{Tm}_{\mathfrak{g}} \mathcal{T}_{+} \gg \operatorname{Real} \mathcal{T}_{+}$, greater than implying at least one order of magnitude.

It will be noticed that the expression for E_{\star}/E_{i} contains a sum going from one to infinity over the variable m. The following explains the approach taken to approximate this sum.

When cylindrical symmetry about the magnetic field direction exists, we have from Wood³ the following expression for τ_+ , for the local regime (q(<< 1) :

where L = the mean free path of the electron, and

m = density of electrons per unit volume of real space.When: ω_{τ} , this reduces to

$$\sigma_{+} = \frac{Mec}{6} \left[\frac{1}{\omega_{c} \tau} + i \right]$$
 (40)

Now we have the equations

$$\operatorname{Real} \ \nabla_{+} = \frac{\operatorname{Mec}}{\operatorname{B}} \frac{1}{\omega_{c} \Gamma} , \qquad (41)$$

and

$$\operatorname{Img.} \, \overline{\mathbf{U}_{+}} = \frac{\operatorname{mec}}{B} \quad . \tag{42}$$

The real part of E_{t}/E_{i} is largest when

$$\left\{ \left(\frac{m\pi}{Q}\right)^2 - \frac{\omega 4\pi}{c^2} \operatorname{Img} \left(\tau_+ \right)^2 = 0 \quad . \tag{43} \right\}$$

Thus, solving for that M* which makes this expression equal to zero, we obtain

$$M^* = \frac{Q}{C} \left(\frac{\omega 4 \text{ mec}}{\pi B} \right)^{k_2} \qquad (44)$$

With the assumption that our finite relaxation time still gives a sufficiently large $\omega_c \gamma$, we sum over a finite number of muscies centered about M*. The contribution to E_t/E_c made by an mudecreases as its distance from M* increases.

From McGroddy, et. al.⁴, we have the following expressions for the case of spherical symmetry:

REAL
$$T_{+} = \frac{(3\omega_{p}^{2}/16V_{F}q)(1-1/X^{2})}{X \times 1}$$
, (45)

Img.
$$T_{+} = (\omega_{p}^{2}/4TT(\omega_{c}-\omega))F(x)$$
, (46)

where

$$X = \frac{q V_F}{\omega_c (1 - \frac{\omega}{\omega_c})}, \qquad (47)$$

$$F(x) = \frac{3}{x^2} \left[\frac{1}{2} - \frac{1 - x^2}{4x} l_{1} l_{1-x} \right], \qquad (48)$$

$$\omega p^2 = 4\pi m e^2 / m^{**}$$
, (49)

and $w_c = eB/m^{**}c$. (50) The effective mass of the electron at the Fermi surface is denoted by m_{*}^{**} and the velocity by V_F .

These expressions hold for an infinite value of $\omega_c \gamma$. Thus, we have that when $\times < 1$, helicons propagate without attenuation, and when $\times \geq 1$, the helicons are damped so severely they disappear.

It should be noted that McGroddy's expressions for ∇_+ can not be substituted into our expression for E_{\pm}/E_{\pm} for any meaningful results. This is brought about because McGroddy lets the real part of ∇_+ equal zero during propagation. For a finite relaxation time, for which E_{\pm}/E_{\pm} is derived, we never have a zero real part of ∇_+ . Thus, we will always get a real part of E_{\pm}/E_{\pm} . If we do assume the real part of ∇_+ is zero during helicon propagation, then we have

$$\frac{E_{t}}{E_{L}} = -\frac{i}{Q} \frac{4\omega}{QC} \sum_{m=1}^{\infty} \frac{(-i)^{m}}{\sum_{m=1}^{\infty} \frac{($$

which gives us a zero real part of E_{\pm}/E_{i} . Thus, we would conclude that there is no helicon propagation, when we know that according to McGroddy, there is helicon propagation. In our calculations for copper and sodium, it will be found that the real part of ∇_{\pm} is smaller than the imaginary part by at least one order of magnitude during propagation. This nonzero real part is necessary to exhibit $R_{eAl} = E_{\pm}/E_{i}$.

We now have a means of predicting at what approximate value of magnetic field we should expect to see helicon propagation start. Setting X=1 and using $m = M^*$, we find that

$$B = \left(\frac{m_{e}^{++2} V_{p}^{2} \pi \omega 4 m C}{e}\right)^{3}.$$
 (52)

It was stated that the knowledge of the Fermi surface is necessary for the calculation of E_t/E_t . This is because we must know the velocity of an electron on the Fermi surface. If $E(\bar{\lambda})$ is the equation of the Fermi surface, then we can relate the Fermi velocity to this surface by

$$(V_{x}, V_{y}, V_{z}) = \frac{1}{\pi} \left(\frac{\partial \mathcal{E}(\overline{h})}{\partial h_{x}}, \frac{\partial \mathcal{E}(\overline{h})}{\partial h_{y}}, \frac{\partial \mathcal{E}(\overline{h})}{\partial h_{z}} \right) .$$
 (53)

In addition to the helicons, we will encounter Gantmakher-Kaner oscillations which are caused by electrons with an extremum in $m_c \overline{V_2}$ along the magnetic field. The period of the oscillations is a constant and is defined by the equation

$$\left[m_{c}\overline{V_{z}}\right]_{ex.} = \frac{eQ}{a\pi c} \left(\frac{B}{m}\right) \qquad m = integer, (54)$$

where the period (β/m) is measured in units of the magnetic field.

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REAL Et/E: FOR SODIUM

Sodium has a spherical Fermi surface. Now we have

$$\frac{\omega}{\omega_{m}} = \frac{e^{2}}{2\pi^{2} \hbar^{2}} \int_{-\kappa MAX} m_{c} dk_{z} \sum_{m=-\infty}^{\infty} \frac{\overline{v}_{m} \overline{v}_{m}}{\frac{1}{\tau} + i(\overline{q}, \overline{v} - \omega - m\omega_{c})}, \quad (55)$$

and

$$\vec{V}_{m} = \frac{1}{2\pi} \int_{0}^{2\pi} \vec{V}(\theta) e^{-\frac{1}{2\pi}} \int_{0}^{0} \left[\frac{1}{2} \cdot \left\{ \vec{V}_{2}(\theta') - \vec{V}_{2} \right\} + m \omega_{2} \right] d\theta'' d\theta .$$
(56)

Now $\overleftarrow{\nabla}_{m}$ is an integral over the variable $\&_{2}$ and will be calculated by numerical integration methods. We must evaluate the integrand for a finite (preferably odd) number of values in order to this integral numerically. This means we must evaluate $\overrightarrow{\nabla}_{m}$ at these particular $\&_{2}$ values also.

We will use a cylindrical co-ordinate system where z, θ , and ρ correspond to A_2, θ , and b_1 where $b_2 = (b_{Fx}^2 + b_{Fy}^2)^{\frac{1}{2}}$. A spherical Fermi surface is drawn in such a co-ordinate system below.



A j-orbit will be designated as a path of intersection of the Fermi surface with a plane $\&_2 = \text{constant.}$ If we evaluate the integrand of $\overleftarrow{\tau}_m$ at j* number of values of $\&_2$, then we will have to deal with j* number of j-orbits. The expression \overrightarrow{V}_m involves an integration over Θ around a j-orbit. To evaluate \overrightarrow{V}_m , we would again use numerical methods. Each j-orbit would be divided into N intervals. The integrand of \overrightarrow{V}_m would be evaluated at the N+1 points in the interval [0,2T].

For a spherical Fermi surface, however, it is possible to simplify matters so that the \bar{v}_m integrations will not have to be done.

Knowing that for a Fermi sphere we have

$$\mathcal{E}(\mathbf{b}_{z}) = \frac{\hbar^{2}}{2m} \left(\mathbf{b}_{x}^{2} + \mathbf{b}_{y}^{2} + \mathbf{b}_{z}^{2} \right), \qquad (57)$$

and recalling Equation (53), we find that \vec{v}_F , the velocity of an electron on the Fermi sphere, and the wave vector, \vec{k}_F , are related by

$$m \vec{v}_F = \hat{\pi} \hat{b}_F \quad (58)$$

Now for any j-orbit $k_{FZ_j}(\Theta) = k_{FZJ} = \text{constant.}$ Thus, we have

$$\overline{V}_{FZj} = \frac{1}{2\pi} \int_{0}^{\infty} \frac{\pi}{m} h_{FZj}(\theta) d\theta = \frac{1}{2\pi} \frac{\pi}{m} \int_{0}^{\infty} h_{FZj} d\theta , (59)$$

and

$$\overline{V}_{FZj} = \frac{A}{2\pi m} \frac{m}{A} V_{FZj} \int_{a}^{2\pi r} d\Theta = V_{FZj}$$
(60)

Therefore, we have that

$$\left\{ V_{z}(\Theta'') - \overline{V}_{z} \right\} = 0 , \qquad (61)$$

which simplifies $\vec{\nabla}_{n}$ to

$$\vec{V}_m = \frac{1}{2\pi} \int_0^{2\pi} \vec{V}(\theta) e^{-im\theta} d\theta \qquad (62)$$

Also, we have that

$$m_{c} = \frac{f_{c}}{2\pi} \int \frac{dbs}{V_{\perp}}, \qquad (63)$$

where

$$V_{\perp} = (V_{xF}^{2} + V_{yF}^{2})^{y_{2}}, \qquad (64)$$

and Thus,

$$d k_{s} = \frac{m}{2} V_{1} d \theta \qquad (65)$$

$$m_{c} = \frac{\pi}{2\pi} \int_{0}^{2\pi} \frac{1}{\pi} V_{1}(\frac{1}{\sqrt{2}}) d\theta \qquad (66)$$

and

$$m_c = m_{c}$$
 (67)

Hence, we also have

$$\omega_c = cyclotron frequency = \frac{eB}{m_c c} = \frac{eB}{m c}$$
. (68)

In the evaluation of $\vec{\nu}_{\varkappa}$, we note that

$$V_{FXj}(\Theta) = V_F \sin \Theta_j \cos \Theta$$
, (69)

$$V_{Fyj}(\Theta) = V_F \operatorname{pin}\Theta_j \operatorname{pin}\Theta$$
, (70)

and

$$V_{FZj}(\theta) = V_F \cos \theta_j \qquad (71)$$

Recalling that

$$aTS_{m,m} = \int_{0}^{2T} i(m-m)\Theta di\Theta, \qquad (72)$$

we obtain

$$V_{mxj} = (V_F \sin Q_j/2) [S_{m,1} + S_{m,-1}],$$
 (73)

$$V_{myj} = (V_F \sin \Theta_j / 2i) [S_{m,1} - S_{m,-1}],$$
 (74)

and
$$V_{mzj} = V_F \cos \varphi_j \delta_{m,o}$$
. (75)

From this, we see that the only contribution to the sum in $\overleftarrow{\tau}_{m}$ will be from the integers n=1 and n= -1. Computing

$$\tau_{xx}$$
 and τ_{xy} , we get

$$T_{XX} = \frac{-c^2 V_F b_F m}{4\pi^2 \pi^2} \int_0^{\pi} \frac{a \sin^3 \varphi d\varphi}{a^2 + b^2} , \quad (76)$$

and

$$T_{xy} = \frac{e^2 V_F k_F m}{4\pi^2 f_{x}^2} \int_0^{\pi} \frac{lr \sin^3 \varphi d\varphi}{e^2 + h^2}, \quad (77)$$

where

$$\alpha = + i q \nabla_2 - i \omega , \qquad (78)$$

and
$$f_r = \omega_c$$
.

$$\sigma_{\pm} = \sigma_{xx} \mp i \sigma_{xy} \qquad (79)$$

so we obtain

$$\overline{\sigma_{\pm}} = -\frac{3}{4} \overline{\sigma_0} \int_0^{\pi} \frac{\sin^3 q \, d\varphi}{1 + iT(\pm \omega_e - \omega + q V_F \cos \varphi)}, \quad (80)$$

where
$$r_0 = m e^{T/m}$$
. (81)

This integral can be evaluated as follows:

$$\chi = \cos \phi$$
, (82)
 $\sin \varphi = (1 - \chi^2)^{\frac{y_2}{2}}$.

and

Such a substitution will give us

$$\nabla_{+} = \frac{3i\nabla_{0}}{4qv_{F}T} \left\{ 2 \Omega - (-\Omega^{2} - 1) \log\left(\frac{-\Omega + 1}{-\Omega - 1}\right) \right\}, \quad (83)$$

where

$$-\Lambda = ((\dot{V}_{\tau}) - \omega + \omega_{E})/q V_{F} . \qquad (84)$$

We now put σ_{+} into E_t/E_i and obtain

$$\frac{E_{\pm}}{E_{i}} = \frac{-i4\omega}{Q_{c}} \sum_{m=1}^{\sigma} \frac{(-1)^{m}}{\left(\frac{m\pi}{Q}\right)^{2} - \left(\frac{\omega}{c}\right)^{2} \frac{4\pi}{i\omega} T_{+}, \quad (85)$$

and

$$\frac{E_t}{E_t} = -\frac{i4\omega}{Q_c} \sum_{m=1}^{7} \frac{(-1)^m}{\left[\left(\frac{m\pi}{Q}\right)^2 - \frac{\omega 4\pi}{c^2} \right] \operatorname{Imag} \left[\Gamma_+\right] + i\frac{\omega 4\pi}{c^2} \operatorname{ReA} \left[\nabla_+\right] \cdot (86)$$

Let

$$S = \left[\left(\frac{m\pi}{Q} \right)^2 - \frac{\omega 4\pi}{c^2} I_{mag} \cdot \tau_+ \right] , \qquad (87)$$

and

$$V = \frac{\omega 4\pi}{c^2} \operatorname{ReA} (T_+ .$$
 (88)

Then we have that

$$\operatorname{ReAl}\left(\frac{E_{t}}{E_{t}}\right) = \frac{4\omega}{Q_{c}} \sum_{m=1}^{c_{1}} \frac{(-1)^{m+1} V}{S^{2} + V^{2}} \qquad (89)$$

A computer program was run using the τ_+ we have just derived for two different cases. In one case, we chose a B-field range of 1,000 gauss to 10,000 gauss, $\omega = 10^6 \text{ sec}$, and $\tau = 10^9 \text{ sec}$. In the second case, we chose the same B-field range and frequency but made $\tau = 10^{10} \text{ sec}$.

The sum in the expression $R_{e_A} (E_t/E_i)$ which is over m going from one to infinity was approximated by a sum going from one to seventy-five. An attempt was made to check the accuracy of this approximation in the following way. We ran a computer program to check the rate of convergence of

$$m = M^{\frac{4}{2} + \chi} \frac{(-1)^{m+1}}{5^{2} + \sqrt{2}}$$

where X equals an arbitrary integer. The results are given in Chart I. A sum was used over a finite number of m's centered about M*. The B-field range was from 7,000 gauss to 8,200 gauss. The number of terms included in each sum increased as B increased. However, the sum over 9 terms was also noted for each B-field so that a comparison could be made to a sum taken over more terms ranging from 19 to 37. The difference in these two sums was calculated, and for $\Upsilon = 10^{9}$ sec., there was a 1% difference or less while for $\Upsilon = 10^{9}$ sec., there was a difference as high as 115%. Thus, we could assume that for $\Upsilon = 10^{9}$ sec., it was slow so that more terms than 9 would have to be included.

We must remember that the M* which makes S zero was defined only when $\omega_c \tau$ was infinite. We see that for $\tau = 10^9 \sec$, $\omega_c \tau \approx 17$ for B = 1,000 gauss and 170 for B = 10,000 gauss. For $\tau = 10^{10} \sec$, $\omega_c \tau \approx 1.7$ for B = 1,000 gauss and 17 for B = 10,000 gauss. Since our values of $\omega_c \tau$ are not particularly large, we may doubt

DISPLAY OF RAPIDITY OF CONVERGENCE OF REAL E_{t}/E_{c}

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CHART I

FOR RELAXATION TIME $\tau = 10^{9}$ sec.

B Field	Sum of 9 Terms	Final Sum	Number of Terms in Final Sum	Difference in Final Sum and <u>Sum of Nine Terms</u>
7000	326-02	326-02	19	0.0%
7100	223-02	223-02	23	0.0%
7200	178-03	180-03	23	1.1%
7300	+.545-04	+.529-04	25	3.0%
7400	+.369-03	+.367-03	27	0.5%
7500	+.184-02	+.184-02	27	0.0%
7600	+.258-02	+.259-02	31	0.4%
7700	+.484-03	+.486-03	31	0.4%
7800	+.103-03	+.104-03	31	1.0%
7900	792-04	776-04	33	2.1%
8000	365-03	363-03	35	0.6%
8100	160-02	160-02	35	0.0%
8200	263-02	263-02	37	0.0%
	FOR RELAXA	TION TIME	1=10 SEC.	
7000	943-04	705-04	19	33.7%
7100	514-04	747-04	23	30.6%
7200	206-04	431-04	23	51.1%
7300	+.289-04	+.133-04	25	115.4%
7400	+.819-04	+.612-04	27	34.4%
7500	+.121-03	+.101-03	27	20.8%
7600	+. 961-04	+.117-03	31	12.0%
7700	+.746-04	+.950-04	31	21.0%
7800	+.235-04	+.394-04	33	41.0%

the validity of using M* in a meaningful way. For B = 1,000 gauss, we have M* = 71, and for B = 10,000 gauss, we have M* = 17. Thus, it might seem that $\omega_c \tau'$ for the case of $\tau = 10^{-10}$ sec. is not sufficiently large to use M* as that m which gives the largest contribution, thus leaving open the question as to what should be chosen as a centering point for the summation. Since our graphs of $R_{e,R} | (E_t/E_t)$ versus B for $\tau = 10^{-3}$ sec. produced smooth regularly occuring waves, and a sum over an insufficient number of terms gave jagged irregular graphs, it was assumed that for the case of $\tau = 10^{-10}$ sec., we were summing over a sufficient number of terms if we got a smooth curve for our graphs. Our choice of summing from one to seventy-five gave us the desired results.

Graphs I through IV give the output for $\Upsilon = 10^{-9}$ sec. Graphs V through IX give the output for $\Upsilon = 10^{-9}$ sec.

Let us recall our predicted value of the magnetic * field at which we expect to see helicon propagation start,

$$B = \frac{m^2 V_F^2 T W + MC}{e}$$
 (90)

From this definition we obtain B = 5,800 gauss. Thus, if we plot R_{eA} (E_t/E_t) versus B, we would expect to see helicon propagation around 5,800 gauss. From our graphs for $\gamma = 10^9$ sec., we see that helicon propagation is

definitely starting at about this point. In the case of $\Upsilon = 10^{10}$ sec., our graph shows helicon propagation starting at about 6,200 gauss. This is not as close in agreement with 5,800 gauss as the previous case. However, as mentioned before, our $\omega_c \Upsilon$ is also one order of magnitude smaller than before, and thus we would not expect a close agreement.

Let us describe the graphic results in more detail. First we consider the graphs for $\Upsilon = 10^{10}$ sec. On Graph I, we see the initiation of helicon propagation at about 6,300 gauss. Graphs II, III, and IV show the helicons increasing in amplitude and in distance between peaks. The curve has a smooth sinusoidal nature.

Now let us look at the graphs for $\Upsilon = 10^{\circ}$ sec. On Graph V we observe Gantmakher-Kaner oscillations. They have a constant period of 375 gauss. Recalling Equation (54), we find that a period of 375 gauss corresponds to $[m \bar{v}_z]_{ext} = 9.3 \times 10^{-20} \text{ gm-cm/sec}$. This, in turn, corresponds to a $b_{z ExT}$ of .88 $\times 10^{+3} \text{ cm}^{-1}$. This is extremely close in value to KMAX, which was taken as .9 $\times 10^8 \text{ cm}^{-1}$. We also note that the amplitude of the Gantmakher waves is smaller than the amplitude of the helicons by a factor of 10^4 .

Graph VI is the same as Graph V with a slight

increase in amplitude. Graph VII shows the start of helicon propagation. Graphs VIII and IX show the helicons gaining in amplitude and increasing in period. We notice that while the Gantmakher-Kaner oscillations were sinusoidal in shape, the helicons are not. They have more of a delta-function appearance about them. Recalling that for the same magnetic field range, the helicons for $T = 10^{-10}$ sec. were sinusoidal shaped, we conclude that the larger relaxation time has definitely caused a distinct change in curve shape.

Graphs X and XI are reproductions of experimental data taken by Wood³ exhibiting these two distinct wave shapes. Here also, we notice the increasing amplitude and period of the helicons.

The relaxation time also affects the magnitude of the amplitudes. We notice that the amplitudes for $T = 10^{10}$ sec. are two orders of magnitude smaller than the amplitudes for $T = 10^{-9}$ sec. in the area of helicon propagation.



GRAPH I CALCULATED - SPHERICALLY SYMMETRIC















GRAPH JUI


<u>GRAPH IX</u> CALCULATED SPHERICALLY SYMMETRIC



SIGNAL

TRANSMITTED

EXPERIMENTAL SPHERICALLY SYMMETRIC

 $\frac{3}{2}$



GRAPH XI

Now we will advance to a cylindrically symmetric Fermi surface, which will include the spherical case. For cylindrically symmetric surfaces we have

$$V_{x} = V_{\perp} \cos \Theta$$
, (91)

$$V_y = V_\perp \sin \Theta$$
, (92)

$$V_{z}(\theta) = V_{z}$$
, (93)

where
$$V_{\perp} = (V_{\chi}^{2} + V_{y}^{2})^{\gamma_{2}}$$
, (94)

and
$$\pi_{\lambda} = m_{\lambda}$$
.

Now again our equation for $\frac{-r}{v_{mv}}$ simplifies to

$$V_{m,\chi} = \frac{1}{2\pi} \int_{0}^{2\pi} V_{\perp} \cos \Theta e^{-im\Theta} d\Theta , \qquad (96)$$

$$V_{m,X} = (V_{\perp}/2) (\delta_{m,1} + \delta_{m,-1}) , \qquad (97)$$

$$V_{my} = (V_{\perp}/2i)(\delta_{m,1} - \delta_{m,-1})$$
 (98)

Now we have

Again using the equations

$$\alpha = + i q V_2 - i \omega , \qquad (100)$$

and
$$\lambda = \omega_c$$
, (101)

we have that

$$\sigma_{xx} = \frac{e^2}{4\pi^2 m_c} \int_{-KMAX}^{+KMAX} \frac{a b_{\perp}^2 db_{\pi}}{a^2 + b^2}, \qquad (102)$$

.

(95)

and

$$\nabla_{xy} = \frac{-e^2}{4\pi^2 m_c} \int \frac{4\pi h_x}{a^2 + h^2} .$$
 (103)

Now we have that

$$\overline{\tau_{+}} = \overline{\tau_{XX}} - i \, \overline{\tau_{XY}} \qquad (104)$$

Thus, we have

Now, as before, we have gotten around evaluating the integral expression of \vec{v}_m . However, while previously we had derived an exact expression for \vec{v}_+ , now it must be evaluated numerically.

We mentioned earlier that sodium was investigated mainly for the purpose of an accuracy check on our computer programs for copper. The integral expression for σ_{\mp} was used to evaluate Real Et/E: over a magnetic field range of 8,800 gauss to 9,500 gauss. We then checked our results with the data gathered for the closed expression of σ_{\mp} . This integral expression of σ_{\mp} will later be used for copper, and from this check, we will know how accurate our results will be.

Now we rewrite our expression for σ_+ as:

$$\overline{v_{+}} = \frac{e^{2} \Upsilon}{4\pi^{3} m_{c}} \int_{-\kappa_{MAX}}^{+\kappa_{MAX}} \frac{(\pi b_{+})^{2} db_{z}}{1 + i \Upsilon (-\omega_{c} - \omega + q \overline{v_{z}})} \qquad (106)$$

We are dealing with a spherical Fermi surface. Let us

look at the drawing below.



For a sphere, we have that KMAX equals the radius $\mathbf{k}_{\rm F}$. We divide the $\mathbf{k}_{\rm z}$ -axis into 52 equally spaced intervals, so we will be evaluating the integrand of $\nabla_{\rm H}$ at 53 points in the interval [-KMAX, +KMAX]. Thus, we will have 53 j-orbits. A typical j-orbit is shown. Now $(\pi \mathbf{k}_{\perp}^2) = A_3^2$ is the cross-sectional area of a j-orbit. We know that for sodium KMAX = 0.9 X 10⁸ cm⁻¹. Thus, from the drawing we can see that

$$A_{j} = \pi (k_{F}^{2} - k_{z}^{2}) \quad . \tag{107}$$

Also recalling that for a Fermi sphere, we have for any j-orbit for which $k_z = \text{constant}$,

$$V_z = \frac{\hbar k_z}{m}$$
 (108)

Thus, we can calculate all the data we need to evaluate v_{\downarrow} numerically. We used this expression of v_{\downarrow} in our

CHART II

Comparison of Real V, and Imaginary V, for Nine mo 's Centered About M* NOTE: In number notation, +18 means 10

B = 8,900 gauss M* = 23

:

Value of	m Real V7	Imaginary V+
19	.281+18	.412+20
20	. 285+18	•414+20
21	. 288+18	.416+20
22	. 292 + 18	.417+20
23	. 296+18	. 419+20
24	•301 + 18	.421+20
25	•306+18	.424+20
26	•311+18	.426+20
27	• 317+18	.428+20
	B = 9,375 gauss	M* = 23
19	•251+18	.390+20
20	•253 + 18	•391+20
21	•256 + 18	• 393+20
22	. 259+18	• 394+20
23	. 263 + 18	• 396+20
24	.266+18	• 398+20
25	.270+18	•399+20
26	. 274 + 18	.401+20
27	•279 + 18	•403+20



expression for Real E_t/E_i using a sum over m going from 1 to 75. Chart II shows a comparison of the real and imaginary parts of σ_+ . We know that for helicon propagation Img. $\sigma_+ \gg \operatorname{Real} \sigma_+$. Chart II verifies that this is so.

Graph XII shows a comparison of our data, written with a solid line, with the data obtained by our previous closed form expression of v_{+} , drawn in a dotted line.

We see from the graphs that the numerically integrated ∇_+ gives a curve shape identical with that which resulted from the closed form ∇_+ , except that the curves are out of phase with one another by about 85 gauss. This phase difference is probably caused by the fact that there is some error inherent in numerical integration techniques. In conclusion, we see that our numerical integrations are accurate enough to give us the correct shape of the curve which is what we are most concerned with.

We now advance to our most complicated fashion of calculating R_{eA} E_{\pm}/E_{\pm}^{\prime} . This method will be the one used to calculate E_{\pm}/E_{\pm}^{\prime} for copper using the true Fermi surface of that metal, but we now wish to check the accuracy of the computer program for the simpler case of sodium. The complication lies in that we must now

evaluate \vec{v}_m by numerical integration methods.

As before, we divide the $\&_2$ axis into fifty-two equally spaced intervals so that we will be evaluating ∇_+ at 53 points in the interval [-KMAX, KMAX]. We then calculate the values of the x, y, and z components of the Fermi velocity at 45 different points equally spaced around each j-orbit. Thus, we will have the v_x , V_y , and V_z values for 45 points in the interval [$o_{,2\pi}$]. These Fermi velocity values were calculated as follows and as demonstrated in the diagram below:





Thus, we have

$$V_{ZP} = \frac{f_{U}k_{Z}}{m_{U}}, \qquad (109)$$

$$V_{XP} = \frac{\hbar}{m} (K M A X^2 - b_z^2)^{V_2} \cos \Theta$$
, (110)

and

$$V_{yp} = \frac{\hbar}{m} \left(K M A X^2 - b_z^2 \right)^{1/2} \sin \Theta \qquad (111)$$

Now, let us recall that

$$\vec{\nabla}_{m} = \frac{1}{2\pi} \int_{0}^{2\pi} \vec{\nabla}(\theta) e^{-\frac{1}{2\pi} \int_{0}^{\infty} \frac{m}{Q} \left\{ V_{2}(\theta') - \overline{V_{2}} \right\} + m\omega_{c} d\theta''} \qquad (112)$$

We have that

$$\bar{V}_{z} = \frac{1}{2\pi} \int_{0}^{2\pi} V_{z}(\theta) d\theta \qquad (113)$$

For each j-orbit, we must do a numerical integration of \overline{v}_z . Let us store this value in an array called AVGVZ(J). Since we are going to do a numerical integration of $v_{m,x}$ and $v_{m,y}$, we must evaluate the integrand at the 45 different values of Θ . Let us denote these angles by Θ_m , where n goes from 1 to 45, and $\Theta_1 = 0$ and $\Theta_{15} = 2\Pi$. Also, let

$$\frac{m\pi}{Q} \left\{ V_{2}(\theta'') - \overline{V}_{2} \right\} = X(\theta'') \qquad (114)$$

We evaluate the expression

$$e^{-\frac{1}{2}} \left\{ \int_{0}^{0} \frac{mT}{Q} \left\{ V_{z}(\theta'') - \overline{V}_{z} \right\} + m w_{c} d\theta''$$

in the following way. For $\Theta = \Theta_1$, we have $e^{\frac{i}{2}\omega_0} \int_0^{\Theta_1} X(\Theta'') + m\omega_c d\Theta'' = e^{-im\Theta_1} e^{-\frac{i}{2}\omega_0} \int_0^{\Theta_1} X(\Theta'') d\Theta''$ (115)

For
$$\Theta = \Theta_2$$
, we have

$$e^{-\frac{i}{\omega_c}\int_0^{\Theta_2} X(\Theta'') + m\omega_c d\Theta''} = e^{-im\Theta_2} e^{-\frac{i}{\omega_c}\int_0^{\Theta_2} X(\Theta'') d\Theta''}, \quad (116)$$

or

$$e^{-i\omega_{2}\int_{0}^{0}X(\theta'') + m\omega_{2}d\theta''} = e^{-im\theta_{2}-i\omega_{2}\int_{0}^{0}X(\theta'')d\theta''} e^{-i\omega_{2}\int_{0}^{0}X(\theta'')d\theta''}, \quad (117)$$

and so forth. We evaluate the integrals using the Trapezoidal Rule. Now, let

$$A_{p,r} = \int_{\theta_p}^{\theta_r} X(\theta'') d\theta'' = E_X P(N(r)) \qquad (118)$$

Thus, we have

$$A_{4,5} = E \times P(N(5)) = \int_{\Theta_4}^{\Theta_5} \overline{X}(\Theta'') d\Theta'', \qquad (119)$$

and also

$$\int_{0}^{0} X(0'') d\theta = A_{1,2} + A_{2,3} + \cdots + A_{n-1,m} = EXPIN(2) + EXPIN(3) + \cdots + EXPIN(m), \quad (119)$$

and

$$\int_{0}^{m-1} X(\theta'') d\theta = A_{1,2} + A_{2,3} + \dots + A_{m-2,m-1} = EXPIN(2) + EXPIN(3) + \dots + EXPIN(m-1), \quad (120)$$

and so forth. Now let

$$PARTP(m) = e^{-\frac{1}{2}\omega_c} SUMm \qquad (121)$$

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and

where

$$SUM_{m} = \sum_{\alpha=1}^{M} EXPIN(\alpha) \qquad (123)$$

If we now let

$$CART(m) = e^{-iN\Theta_m}$$
, (124)

then at each of the 45 points, the value of the integrand will be given by the expressions

$\begin{pmatrix} Value \text{ of Integrand} \\ \text{for Points on the} \\ Positive \not{}_{z} Axis \end{pmatrix}_{mx} = V_{x}(\Theta_{m}) PARTF(m) CART(m), \quad (125) \\ \text{and} \\ \begin{pmatrix} Value \text{ of Integrand} \\ \text{for Points on the} \\ \text{Negative} \not{}_{z} Axis \end{pmatrix}_{mx} = V_{x}(\Theta_{m}) PARTN(m) CART(m). \quad (126)$

Now, let us recall Equations (73) and (74) which state

$$V_{MX} = V_F \operatorname{Din} Q_j / 2 \left[S_{m,1} + S_{m,-1} \right],$$

and

$$V_{my} = V_F \sin Q_j / 2i \left[\delta_{m,1} - S_{m,-1} \right]$$

Thus, we have that

REAL
$$V_{1,X} = V_F \sin Q_{j}/2$$
 Imag. $V_{1,X} = 0$, (127)

REAL
$$V_{-1,X} = V_F Ain O_j / 2$$
 Imag. $V_{-1,X} = O_j$ (128)

Real
$$V_{1y} = 0$$
 Imag. $V_{1,y} = -V_F \sin \varphi_j/2$, (129)

and

REAL
$$V_{-1y} = 0$$
 I mag. $V_{-1,y} = V_F \sin q_j / 2$. (130)

Now for sodium $V_F = 1.03 \times 10^8$ cm/sec, and for the j=1-orbit, $\theta_j = \frac{1}{2}$. Thus, we should expect the following values:

Real $V_{1,X} = .515 \times 10^8$ Imaginary $V_{1,X} = 0.0$ Real $V_{-1,X} = .515 \times 10^8$ Imaginary $V_{-1,X} = 0.0$ Real $V_{1,y} = 0.0$ Imaginary $V_{1,y} = -.515 \times 10^8$ Real $V_{1,y} = 0.0$ Imaginary $V_{-1,y} = .515 \times 10^8$

Using data from a computer run for B=7,000 gauss for the j=1 orbit or $\lambda_z = 0.0$, we obtain Chart III. Thus, we have checked the accuracy of calculating the N_m 's with our program and found an error of 1.1%.

In calculating E_{t}/E_{t} , we approximate the sum over m going to infinity with a sum of nine terms centered about M*, since it was shown earlier in Chart I that for $T = \sqrt{0}^{9}$ sec., we had very rapid convergence of Real E_{t}/E_{t} . Recalling that during helicon propagation, $\operatorname{Imag} T_{t} > \operatorname{Real} T_{t}$, we note the sample data given in Chart IV. The data given by our previous program, where the T_{t} 's but not the V_{m} 's were calculated numerically, is on the left. The data given by this last program, where both T_{t} and V_{t} are calculated numerically, is given on the right. The value of the magnetic field is 8,850 gauss. Helicons are propagating in this region, and we notice that $\operatorname{Imag} T_{t} > \operatorname{Rea} T_{t}$ by a factor of 100. The close values of the T_{t} serve as another accuracy check of our last program.

NUMERICALLY INTEGRATED V_{NX} AND V_{NY} VALUES FOR SODIUM AT B=7,000 G.

CHART III

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<u>N</u>	Real VNK	Imaginary V _{NX}	Real VNY	Imaginary VNY
-5	+.2440+03	2083-02	3605+00	+.1656+01
-4	+.2428+03	6373-01	+.1024+01	+.6730+00
-3	+.2423+03	3705+00	5982-01	+.3839-01
-2	+.2431+03	4984+00	1394+00	3860+00
-1	+.5209+08	+.2924+00	+.4569-02	+.5209+08
0	±.2427+03	+.0000	+.2924+00	.0000
+1	+.5209+08	2924+00	+.4569-02	+.3860+00
+2	+.2431+03	+.4984+00	1394+00	3839-01
+3	+.2423+03	+.3705+00	5982-01	6730+00
+4	+.2428+03	6373-01	+.1024+01	1656+01
+5	+.2440+03	+.2083-02	3605+00	+.1656+01

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VALUES OF REAL J. AND IMAGINARY J. FOR SODIUM FOR B=8,850 GAUSS AND M-RANGE OF 18 TO 26

CHART IV

М			J Calculated With Numerically Integrated Vals		
	Real T+	Imaginary V.	Real J+	Imaginary Tr	
18	.2870+18	•4145 + 20	.292+18	.418+20	
19	.2907+18	.4162+20	. 296 + 18	.420+20	
20	.2947+18	.4180+20	•300+18	.422+20	
21	•2991+18	.4200+20	•310+18	.426+20	
22	•3038+18	.4221+20	• 316+18	.429+20	
23	•3089+18	•4243+20	• 322+18	•431+20	
24	•3144 + 18	.4267+20	• 328+18	•434+20	
25	• 3205+18	.4292+20	• 335+18	•437+20	
26	.3270+18	.4319+20	•352+18	•443+20	

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The program used is included in the appendix. We used a B-field range from 8,800 gauss to 9,450 gauss in order to check our graphic results with Graph XII. Our data is plotted on Graph XIII. Graph XII is superimposed on our data for comparison.



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Et/E: For Copper

Summarizing our approach to sodium, we found E_t/E_i by calculating σ_+ in three ways: (1) A closed form σ , (2) An integrated σ with closed expressions for the \vec{v}_m 's, and (3) An integrated σ with numerically integrated expressions of \vec{v}_m .

In the case of the true Fermi surface for copper, we have as a good approximation

$$V_{z}(\theta) = V_{z} + \alpha(b_{z}) \cos 4\theta,$$
 (131)

$$V_{X} = b \sin \theta, \qquad b \neq b(b_{2}) \neq b(\theta) \qquad (132)$$
$$V_{y} = -b \cos \theta, \qquad (133)$$

for b_z parallel to the [001] direction (see Figure I). Thus, our expression for \vec{v}_m would be

$$V_{mx} = \frac{l_{w}}{2\pi} \int_{0}^{2\pi} \sin \Theta e = i \frac{d}{d\Theta} \frac{1}{d\Theta} \frac{d\Theta}{d\Theta},$$
 (134)

and

$$V_{my} = \frac{1}{2\pi} \int_{0}^{2\pi} \cos \theta e^{-i\frac{\alpha}{4w_{c}}} \sin 4\theta_{-im\theta} = d\theta.$$
 (135)

This is not a simple expression to evaluate, as in the case of sodium where the V_{mx} and V_{my} were evaluated as delta functions. Therefore, methods (1) and (2) are eliminated for use with the true Fermi surface of copper. This leaves method (3) for use. However, as an accuracy check we can use a cylindrically symmetric approximation

of the copper Fermi surface with method (2). We have already described method (2) in the previous section on sodium. Our equation for ∇_+ was given as

$$T_{+} = \frac{e^{2}T}{4\pi^{3}m_{c}} \int \frac{(\pi b_{1})^{2} db_{2}}{1 + iT(-w_{c} - w + qV_{2})}, \quad (136)$$

which can be rewritten as

$$T_{+} = \frac{e^{2}T}{4\pi^{3}m_{c}} \left(\frac{A(b_{z})db_{z}}{1+iT(-w_{c}-w_{-}\frac{\hbar}{2\pi}\frac{\partial A}{\partial b_{z}})}, (137) \right)$$

where

$$TT b_{\perp}^{2} = A(b_{\perp}), \qquad (138)$$

and

$$mV_2 = -\frac{\pi}{2\pi} \frac{2A}{2b_2} , \qquad (139)$$

where $A(b_2)$ is the cross-sectional area made by a plane perpendicular to the b_2 axis intersecting the Fermi surface.

We did computations for the directions [001] and [111]. From Wood, data for $A(b_1)$ and $\frac{\Im A}{\Im b_2}$ for these two directions was available. The drawing below shows the orientation of the b_2 axis with the Fermi surface for these two directions.



Let us first examine the [III] direction. A Υ of 10⁻⁹ sec. was used, and two different frequencies $f = 395.2 \times 10^3$ cycles/sec. and $f = 6 \times 10^5$ cycles/sec. were investigated.

Graph XIV is experimental data obtained by Wood for a frequency of 400.0 X 10³ cycles/sec. and a B-field range of 10 to 30 thousand gauss. We see Gantmakher-Kaner oscillations with a B-field period of 600 gauss. However, the graph displays beats, and thus there must be present two Gantmakher-Kaner waves of slightly different frequencies. Our frequencies are expressed in the units cycles per unit gauss magnetic field.

If we have two waves expressed as

$$y_{i} = A \cos\left(2\pi f_{i} t + \phi_{i}\right) , \qquad (140)$$

and

$$y_2 = A \cos(2\pi f_2 t + \phi_2)$$
, (141)

then, adding, we get

$$y = y_1 + y_2 = A \left[\cos \left(2\pi f_1 t + \varphi_1 \right) + \cos \left(2\pi f_2 t + \varphi_2 \right) \right] .$$
 (142)
Making use of the trigonometric identity

$$\cos \alpha + \cos b = 2 \left[\cos \frac{1}{2} (\alpha + b) \right] \left[\cos \frac{1}{2} (\alpha - b) \right], \qquad (143)$$

we obtain

$$y = 2A \cos \left[2\pi \left(\frac{f_1 - f_2}{2} \right) t + \left(\frac{\phi_1 - \phi_2}{2} \right) \right] \cos \left[2\pi \left(\frac{f_1 + f_2}{2} \right) t + \left(\frac{\phi_1 + \phi_2}{2} \right) \right]$$
(144)

The composite wave may be regarded as having a frequency equal to $\frac{1}{2}(f_i + f_2)$, which is the average of the frequencies of the original waves. The amplitude of the composite

wave is given by the quantity in the first set of brackets in Equation (144). The amplitude varies with a frequency of $\frac{1}{2}(f_1 - f_2)$, and the number of beats per second is given as $(f_1 - f_2)$. The drawing below is a typical representation of the presence of beats.

Graphs XV and XVI show the results for a frequency of 395.2 X 10^3 cycles/sec. for a B-field range of 2,000 to 14,000 gauss. Graph XVII displays the beats which begin at 5,700 gauss and end at about 13,700 gauss. The results for a frequency of 6 X 10^5 cycles/sec. are shown in Graphs XVIII through XX for a B-field range of 100 to 20,000 gauss. Graph XXI displays the beats which begin at 6,400 gauss and end at about 18,000 gauss. In both cases, the envelope of the wave is irregular in shape, but still displays the periodic recurrence of increasing and decreasing wave amplitudes. Since our Gantmakher-Kaner oscillations do not have the simple forms of Equations (140) and (141) and may not have the same amplitude, we understand why our graphic results



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GRAPH XI



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GRAPH XVII









are not as symmetric as the above drawing. The irregular envelope of the wave which we see would probably result mathematically by summing the two Gantmakher-Kaner waves expressed as Fourier series. Because of the irregularity of the envelope, we are unable to determine the beat frequency accurately from the graph.

Charts V and VI show the calculation of the period of the Gantmakher-Kaner waves, both cases of frequency giving a period of about 600 gauss. This agrees with the experimental data mentioned before.

Recall Equation (54) which states that

 $(AB)\frac{CQ}{\partial \Pi C} = [m_e V_z]_{EXT.} = -\frac{\hbar}{2\pi}\frac{2A}{2h_z},$

where $\Delta B = 6/m_{\nu}$ = the period of the G-K wave. Now, Graph A shows a plot of $\partial A/\partial K_{z}$ versus $a|\Delta_{z}|$ for the [111] direction of copper. We see two extremal values of the function $\partial A/\partial K_{z}$ which have been labeled P₁ and P₂. P₁ corresponds to $\partial A/\partial K_{z}$ equaling 6.57 X 10⁸ cm⁻¹, and P₂ corresponds to $\partial A/\partial K_{z}$ equaling 8.1 X 10⁸ cm⁻¹. Thus, there are two values of $[m V_{z}]_{z \times T}$. which should give Gantmakher-Kaner waves, each having a different period. For P₁, we obtain $\Delta B = 538$ gauss, and for P₂, we obtain $\Delta B = 664$ gauss.

Calculation of Frequency of Waves for the L(1) Direction of Copper for $f = 395.2 \times 10^3$ cyc/sec

CHART V	
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Peak No. on Graph	Value of Et/E:	B-Field	Period in Gauss
1	64×10^{-10}	5,700	
2	$78 \times 10^{-1.0}$	6,400	700
. 3	98×10^{-10}	7,000	600
4	49×10^{-10}	7.700	700
5	89×10^{-10}	8,300	600
6	132×10^{-10}	8,900	600
7	122×10^{-10}	9.500	600
8	58×10^{-10}	10.100	600
q	91×10^{-10}	10 700	600
10	117×10^{-10}	11 300	600
11	117×10^{-10}	11 000	600
10	10^{-10}	11,900	600
12	40×10^{-10}	12,500	600
13	60×10^{-10}	13,100	600
14	63 x 10 - °	13,700	

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Calculation of Frequency of Waves for the (111) Direction of Copper for $f = 6 \times 10^5 \text{ cyc/sec}$

CHART VI

<u>Peak No. on Graph</u>	Value of Et/E:	B-Field	Period <u>in Gauss</u>
1	123×10^{-10}	6,400	
2	65×10^{-10}	7.000	600
3	119×10^{-10}	7.700	700
4	121×10^{-10}	8,300	600
5	107×10^{-10}	9,000	700
6	135×10^{-10}	9,600	600
7	167×10^{-10}	10,200	600
8	107×10^{-10}	10,200	600
0	107×10^{-10}	10,000	600
7	109×10^{-10}	11,400	700
10	108 x 10	12,100	600
11	69×10	12,700	600
12	75 x 10 -10	13,300	500
13	101 x 10	13,800	700
14	70×10^{-10}	14,500	600
_	48×10^{-10}	15,100	600
-	61×10^{-10}	15,700	600
-	85×10^{-10}	16,300	600
-	55×10^{-10}	16,900	500
~	37×10^{-10}	17,400	500
-	55×10^{-10}	18,000	600

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Now the frequency is related to the period by

$$f = \frac{1}{(\Delta B)} \qquad (145)$$

Thus, we have that

or

 $f_1 = \frac{1}{538} \text{ gauss}^{-1} = .0018 \text{ gauss}^{-1}$, and $f_2 = \frac{1}{664} \text{ gauss}^{-1} = .0015 \text{ gauss}^{-1}$. Therefore, the frequency of the composite wave is given by

> $f^* = \frac{y_2}{2} (.0018 + .0015) \text{ gauss}^{-1}$, $f^* = .00165 \text{ gauss}^{-1}$

Now the ΔG from the graph was 600 gauss, corresponding to a frequency of .0016 gauss⁻¹. Thus, we have a perfect agreement with theory. That is, the two frequencies of the Gantmakher-Kaner waves making up the beats are .0015 gauss⁻¹ and .0018 gauss⁻¹. We also note that the experimental value of the frequency of the composite Gantmakher-Kaner wave corresponds to a value of $\Im A/\Im b_Z$ of 7.6 X 10⁸ cm⁻¹ on Graph A. This corresponds to a frequency of .0016 gauss⁻¹. Thus, our data agrees with experimental data as well. Let us now look at the results for the [001] direction of copper. Graph XXXII is an experimental curve obtained by Wood showing helicon propagation starting at about 16,000 gauss.

We have computed data for a frequency of 395.2 X13^{hz}. for $\Upsilon = 10^{-9}$ sec. and $\Upsilon = 10^{-10}$ sec. Graphs XXII through XXVII show the results for $\Upsilon = 10^{-9}$ sec. for a B-field range of 8,000 to 19,000 gauss. Helicons begin propagating at about 15,275 gauss as seen from Graph XXV. The helicons have a delta function shape, just as the helicons for sodium had at $\Upsilon = 10^{-9}$ sec. Graphs XXII through XXV show waves which are not constant in period nor increase in amplitude. These can be called neither helicons nor Gantmakher-Kaner oscillations. Their source is unknown to us.

Graph B is a graph of $\Im A/\Im b_z$ versus b_z for the [001] direction of copper. We see an extremal value of $\Im A/\Im b_z$ equal to 2.2 X 10⁸ cm⁻¹, which means that we would expect a Gantmakher-Kaner wave contribution from those electrons having a N_z velocity corresponding to this value of $\Im A/\Im b_z$. However, experimentally, no Gantmakher-Kaner oscillations are seen. Realizing that experimental results would come from use of the true


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GRAPH XXV CALCULATED CYLINDRICALLY SYMMETRIC



GRAPH XXVI



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CALCULATED CYLINDRICALLY SYMMETRIC



noncylindrically symmetric Fermi surface, one might suspect that our assumption of a cylindrically symmetric Fermi surface, which Graph B describes, causes the Gantmakher-Kaner oscillations. This will be discussed further in the data analyzation of the copper calculations using the true Fermi surface values.

On Graph B we also note the value of $|\lambda/\partial b_z|_{MAX}$, which is responsible for the absorption edge. The value of $[\lambda/\partial b_2]_{MAX}$ is 9.4 X 10⁸ cm⁻¹. Using the equation

$$\omega_c = eB/mc = \omega + gV_z , \qquad (146)$$

we saw that the last group of electrons which can experience Doppler - shifted cyclotron resonance are those having the velocity V_{ZMAX} . Now since $\omega << q V_{Z}$, we have

$$eB^{*/mc} = qV_{ZMAX}$$
, (147)

or

$$B^{+} = q V_{ZMAX} mc/e = q c \hbar (2A/2bz)/2\pi e . (148)$$

Thus, B* is the value of B at which we would expect helicons to start propagating. Using $[2A/2]_{MAX} = 9.4 \times 10^8 \text{ cm}^{-1}$, we obtain that

where m = 41 for $q = m \pi / Q$. This value of m corresponds to M*, defined earlier as that value of m in \vec{q} which



gives the largest contribution to REAL E_t/E_i . Helicons are seen to start at about 15,275 gauss; however, as usual, the exact point is impossible to locate.

Data was calculated for $\Upsilon = 10^{-10}$ sec. for a B-field range of 13,250 to 17,500 gauss. Graphs XXVIII through XXX show the data. Helicons are seen to begin propagating at about 15,750 gauss, which is in good agreement with the experimental data. We also note the smooth sinusoidallike shape of the helicons, differing from the delta shaped peaks for the previous Υ . This result from the change of Υ is identical with that seen in sodium.





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GRAPH XXX ED CYLINDRICALLY SYMMETRIC

We now conclude this thesis with the results for the final case of copper in which we use the true Fermi surface. Again we use the computer program included in the appendix in which the \vec{V}_{mx} and \vec{V}_{my} are evaluated by numerical integrations. The values of the V< , Vy and V_z components of the velocity for an electron on the Fermi surface were made available by Wood. Recalling that

$$\vec{V} = \frac{1}{2\pi} \left(\frac{\partial \mathcal{E}(\mathbf{h})}{\partial \mathbf{h}} \right)$$

and that

$$\vec{v} = \lim_{\Delta \vec{b} \to 0} \frac{1}{\vec{b}} \frac{\mathcal{E}(\vec{b} + \Delta \vec{b}) - \mathcal{E}(\vec{b})}{\Delta \vec{b}}$$
, (149)

we see that it is necessary to know the constant energy surface \overrightarrow{bk} away from the Fermi surface. In the calculation of the velocity by Wood, it was assumed that the constant energy surface \overrightarrow{bk} away from the Fermi surface was concentric in shape with the Fermi surface. If this is not a valid assumption, then our velocity on the Fermi surface is not accurate at all points and a difference in the data results and the experimental curve might be expected. The results are shown on Graphs XXXI and XXXIII. A frequency of 395.2 X 10³ hertz and a T of 10⁻⁹ sec. was used. Graph XXXI displays a B-field range of 16,500 gauss to 17,500 gauss, and Graph XXXIII shows a B-field range of 11,100 to 11,900 gauss. The [001] direction of copper was investigated; thus, we were able to compare our data with Graphs XXIII, XXIV, XXVI, and XXVII which show the data for the cylindrically symmetric Fermi surface approximation. The amplitudes of the peaks of Graphs XXXI and XXXIII are smaller by two orders of magnitude than the amplitudes of their respective peaks on Graphs XXIII, XXIV, XXVI, and XXVII. The numerical integration techniques used in the program cannot be responsible for the decrease since their accuracy was checked in the case of sodium. Recalling Graph XII, we see an amplitude for a helicon calculated by our numerical integrations which is identical with that calculated by the earlier two methods in which numerical integrations were not used.

Graph XXXI does, however, show the familiar delta shape due to the relaxation time of 10^{-9} sec. and agrees with the experimental curve on Graph XXXII in that helicons are propagating in this B-field region. Graph XXXIII does not agree with the experimental curve which shows nothing propagating in this region. Our only explanation of this is that the velocity on the Fermi surface is not accurate at all points due to a non-concentric constant energy surface Δb away from the Fermi surface. Also, the loss of cylindrical symmetry of the Fermi surface is believed to be the cause of the reduction in amplitude of the peaks by a factor of 10^{-2} which was mentioned earlier. As for the case of sodium, limited computer time forced us to investigate only small B-field ranges in this final case.

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GRAPH XXXI



GRAPH XXXIII







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COMPUTER PROGRAM FOR THE CALCULATION OF THE RATIO OF THE MAGNITUDE OF THE TRANSMITTED ELECTRIC FIELD TO THE INCIDENT ELECTRIC FIELD , .* The following are the variables contained in the program: Q = thickness of metal slab (cm.) AMXK= maximum value of the wave vector in the z-direction W= frequency of incident electromagnetic wave T= relaxation time PE= electron density KPS= the number of magnetic field values for which we are evaluating E_{+}/E_{i} BO= the lowest magnetic field value that we are using SU= the step interval between consecutive magnetic field values IO= the number of integers which when added to IPED is considered as a sufficient approximation to the infinite sum in the expression for E_t/E_i JS= number of orbits into which the Fermi Surface has been divided between $k_z=0.0$ and $k_{z=AMXK}$ J= DO-LOOP index for orbits N= number of Fermi surface data points for an orbit. N may be even or odd, but the points are assumed as evenly distributed in the interval K= DO-LOOP index for wave number. F = wave number = K*PI/QKS= integer which gives maximum wave number NN= DO-LOOP index for the sum contained in the expression NNO= the sum in the expression which goes from plus to minus infinity is replaced by a sufficiently approximate sum going from plus to minus (NNO-1)/2 NT = IO + 1

The values of the variables in the program for sodium that were used are as follows:

Q = 0.1 cm
AMXK = 0.9 X
$$10^8$$
 cm⁻¹
W = 10^6 hz.
PE = 2.5 X 10^{22} electrons/cm³
IO = 5
JS = 27
N = 45
NNO = 11
NNT = 6

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The values of the variables in the program for copper that were used are as follows:

$$Q = 0.084 \text{ cm}$$

$$AMXK = 1.44 \times 10^8 \text{ cm}^{-1}$$

$$W = 10^6 \text{ hz.}$$

$$PE = 8.5 \times 10^{22} \text{ electrons/cm}^3$$

$$IO = 4$$

$$JS = 27$$

$$N = \text{variable}$$

$$NNO = 9$$

$$NT = 5$$

DOUBLE PRECISION A, ANSRX, ANSRY, ANSIX, ANSIY, BETA, GAMMA, S. V. DET, DARTR, DARTI, ÉTEIR, ÉTEII COMPLEX CART. CSUMA, CSUMB, CSUMC, CSUMD, PARTP, PARTN, SUMAXN, SUMAÝN, SUMBXN, SUMBYN, SÚMCXN, SUMAXP, SUMAYP, SUMBXP, SUMBYP, SUMCXP, SUMCYP, SUMCYN, DENOMP, DENOMN, DENM, DENP, VNXPP, VNYPP, VNXNN, VNYNN, VNXN, VNYN, VNYP, VNXP, SAXX, SAXY, SBXXA, SBXXB, SBXYA, SBXYB, SIGXX, SIGXY, CIA, CVYXX, CVYXY DIMENSION DENP(NNO), DENM(NNO), VNXPP(NNO), VNYPP(NNO), VNXNN(NNO), VNYNN(NNO), VX(NMAX), VY(NMAX), VZ(NMAX), EXPIN(NMAX), PARTP(NMAX), PARTN(NMAX), CART(NMAX), NAN(JS), WC(JS), ARRAY(JS), AVGVZ(JS), ARMC(JS), CVYXX(JT,KS), CVYXY(JT,KS) Q =AMXK= CIA = (0.0.1.0)H= 1.0545E-27 TM = 9.11E - 28R = 9.11E - 28W =T= D = 2.99E + 10E = 4.8E - 10PE= PI = 3.1416ST=(W*PE*E)/(PI*D) $A = ((E^{*}2)^{AMXK})/((H^{*}2)^{(PI^{*}2)}((2^{JS})-2))$ REWIND 8 DO 1000 KP=1.KPS B = (BO + SU*(KP-1))PEB=(2.0*Q*SQRT(ST))/SQRT(B)IPEC=INT(PEB) IPED=IPEC + IO IPEG=IPEC - IO DO 265 J=1.JS READ(8) ARRAY(J), NAN(J) N=NAN(J)READ(8) (VX(I), I=1, N)READ(8) (VY(I), I=1, N)READ(8) (VZ(I), I=1, N) CALCULATION OF THE ARRAY AVGVZ IF(N-2*(N/2)) 110,115,110 SUMA = VZ(1) + VZ(N-1)SUMB=0.0 KN = N - 2DO 111 L=2,KN,2 SUMB=SUMB + VZ(L)SUMC=0.0KN = N - 3DO 112 L=3,KN,2

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SUMC = SUMC + VZ(L)112 SUMD = (VZ(N-1) + VZ(N))/(2.0*(N-1))AVGVZ(J) = (SUMA + 4.0*SUMD + 2.0*SUMD)/(3.0*(N-1)) + SUMDGO TO 120 SUMA=VZ(1) + VZ(N)110 SUMB=0.0 KN = N - 1DO 116 L=2,KN,2 SUMB=SUMB + VZ(L)116 SUMC=0.0 KN=N-2DO 117 L=3,KN,2 SUMC = SUMC + VZ(L)117 AVGVZ(J) = (SUMA + 4.0*SUMB + 2.0*SUMC)/(3.0*(N-1))GO TO 120 120 CONTINUE CALCULATION OF THE ARRAYS ARMC(J) AND WC(J) ARMC(J) = 1.37 * R*(NAN(J) * 1.0) / (NAN(1) * 1.0)WC(J) = (E*B) / (D*ARMC(J))DO 600 K=IPEG. IPED F = (K*PI)/QEXPIN(1)=0.0DO 125 M=2.N EXPIN(M) = (VZ(M) + VZ(M-1) - 2.0 * AVGVZ(J)) * F*PI/((N-1))125 SUM=0.0DO 130 M=1,N SUM=SUM + EXPIN(M)PARTN(M)=CEXP(CIA*SUM/WC(J)) PARTP(M) = CEXP((-1.0*CIA*SUM)/WC(J))130 CART(1) = (1.0.0.0)DO 12 NN=1.NNO DO 135 M=2.N CART(M)=CEXP(-2.0*CIA*(-NNT + NNO)*PI*(M-1)/((N-1)*1.0) 135 IF (N-2*(N/2)) 181.182.181 SUMAXP=VX(1)*PARTP(1)*CART(1) + VX(N-1)*PARTP(N-1)*CART(N-1)182 SUMAYF=VY(1)*PARTP(1)*CART(1) + VY(N-1)*PARTP(N-1)*CART(N-1)SUMAXN = VX(1) * PARTN(1) * CART(1) + VX(N-1) * PARTN(N-1) * CART(N-1)SUMAYN=VY(1)*PARTN(1)*CART(1) + VY(N-1)*PARTN(N-1)*CART(N-1)SUMBXP = (0.0.0.0)SUMBYP = (0.0.0.0)SUMBXN = (0.0, 0.0)SUMBYN = (0.0.0.0)KN = N - 2DO 146 L=2,KN,2 SUMBXP=SUMBXP + VX(L)*PARTP(L)*CART(L) SUMBYP=SUMBYP + VY(L)*PARTP(L)*CART(L) SUMBXN=SUMBXN + VX(L)*PARTN(L)*CART(L) 146 SUMBYN=SUMBYN + VY(L)*PARTN(L)*CART(L) SUMCXP = (0.0, 0.0)

SUMCYP = (0.0.0.0)SUMCXN = (0.0.0.0)SUMCYN = (0.0, 0.0)KN = N - 3DO 147 L=3,KN,2 SUMCXP=SUMCXP + VX(L)*PARTP(L)*CART(L) SUMCYP=SUMCYP + VY(L)*PARTP(L)*CART(L) SUMCXN = SUMCXN + VX(L) * PARTN(L) * CART(L)SUMCYN=SUMCYN + VY(L)*PARTN(L)*CART(L) 147 SUMDXP = (VX(N-1)*PARTP(N-1)*CART(N-1)+VX(N)*PARTP(N)*CART(N))/(2.0*(N-1))SUMDYP = (VY(N-1)*PARTP(N-1)*CART(N-1)+VY(N)*PARTP(N)*CART(N))/(2.0*(N-1))SUMDXN = (VX(N-1)*PARTN(N-1)*CART(N-1)+VX(N)*PARTN(N)*CART(N))/(2.0*(N-1))SUMDYN = (VY(N-1) * PARTN(N-1) * CART(N-1) + VY(N) * PARTN(N) * CART(N))/(2.0*(N-1))VNXP = (SUMAXP + 4.0*SUMBXP + 2.0*SUMCXP)/(3.0*(N-1))+(SUMDXP)VNYP = (SUMAYP + 4.0*SUMBYP + 2.0*SUMCYP) / (3.0*(N-1)) + (SUMCYP)VNXN=(SUMAXN + 4.0*SUMBXN +2.0*SUMCXN)/(3.0*(N-1))+(SUMDXN) VNYN = (SUMAYN + 4.0*SUMBYN + 2.0*SUMCYN)/(3.0*(N-1)) + (SUMDYN)GO TO 60 SUMAXP=VX(1)*PARTP(1)*CART(1) + VX(N)*PARTP(N)*CART(N)181 SUMAYP=VY(1)*PARTP(1)*CART(1) + VY(N)*PARTP(N)*CART(N)SUMAXN = VX(1) * PARTN(1) * CART(1) + VX(N) * PARTN(N) * CART(N)SUMAYN = VY(1) * PARTN(1) * CART(1) + VY(N) * PARTN(N) * CART(N)SUMBXP = (0.0.0.0)SUMBYP = (0.0.0.0)SUMBXN = (0.0, 0.0)SUMBYN = (0.0.0.0)KN = N - 1DO 141 L=2.KN.2 SUMBXP= SUMBXP + VX(L)*PARTP(L)*CART(L) SUMBYP= SUMBYP + VY(L)*PARTP(L)*CART(L) SUMBXN = SUMBXN + VX(L)*PARTN(L)*CART(L)SUMBYN= SUMBYN + VY(L)*PARTN(L)*CART(L) 141 SUMCXP = (0.0.0.0)SUMCYP = (0.0, 0.0)SUMCXN = (0.0, 0.0)SUMCYN = (0.0.0.0)KN=N-2DO 142 L=3,KN,2 SUMCXP = SUMCXP + VX(L) * PARTP(L) * CART(L)SUMCYP=SUMCYP + VY(L)*PARTP(L)*CART(L) SUMCXN=SUMCXN + VX(L)*PARTN(L)*CART(L) 142 SUMCYN=SUMCYN + VY(L)*PARTN(L)*CART(L)VNXP = (SUMAXP + 4.0*SUMBXP + 2.0*SUMCXP)/(3.0*(N-1))VNYP = (SUMAYP + 4.0*SUMBYP + 2.0*SUMCYP)/(3.0*(N-1))VNXN = (SUMAXN + 4.0 * SUMBXN + 2.0 * SUMCXN) / (3.0 * (N-1))

VNYN = (SUMAYN + 4.0*SUMBYN + 2.0*SUMCYN)/(3.0*(N-1))60 EN=NN $DENOMP = (1 \cdot 0/T) + CIA * (F * AVGVZ(J) - W - ((-NNT * 1 \cdot 0 + EN) * WC(J)))$ DENOMN = (1.0/T) - CIA*(F*AVGVZ(J)+W+((-NNT*1.0+EN)*WC(J)))VNXPP(NN)=VNXP VNYPP(NN)=VNYP VNXNN(NN)=VNXN VNYNN(NN)=VNYN DENP(NN)=DENOMP 12 DENM(NN)=DENOMN CALCULATION OF SUMS FOR SIGXX AND SIGXY CSUMA = (0.0.0.0)CSUMB = (0.0.0.0)CSUMC = (0.0.0.0)CSUMD = (0.0, 0.0)DO 10 NN=1.NNO CSUMA=CSUMA + VNXNN(NN)*CONJG(VNXNN(NN))/DENM(NN) CSUMB=CSUMB + VNXNN(NN)*CONJG(VNYNN(NN))/DENM(NN) IF (J-1) 275,276,275 275 CSUMC=CSUMC + VNXPP(NN)*CONJG(VNXPP(NN))/DENP(NN) CSUMD=CSUMD + VNXPP(NN)*CONJG(VNYPP(NN))/DENP(NN) 276 CONTINUE 10 CONTINUE CVYXX(JS+1-J,K)=CSUMACVYXY(JS+1-J,K) = CSUMBCVYXX(JS-1-J,K) = CSUMCCVYXY(JS-1-J,K) = CSUMD600 CONTINUE 265 CONTINUE REWIND 8 DO 700 K=IPEG.IPED F=(K*PI)/QSAXX=ARMC(JS)*CVYXX(1,K) + ARMC(JS)*CVYXX(2*JS-1,K) SAXY=ARMC(JS)*CVYSY(1,K) + ARMC(JS)*CVYXY(2*JS-1,K) SBXXB=ARMC(1)*CVYXX(JS,K) SBXYB=ARMC(1)*CVYXY(JS,K) SBXXA = (0.0, 0.0)SBXYA = (0.0, 0.0)DO 285 L=2,JS-1,2 SBXXA=SBXXA + ARMC(JS+1-L)*CVYXX(L,K)+ARMC(L)*CVYXX(JS-1+L,K) 285 SBXYA=SBXYA + ARMC(JS+1-L)*CVYXY(L,K)+ARMC(L)*CVYXY(JS-1+L,K) DO 290 L=3, JS-2,2 SBXXB=SBXXB + ARMC(JS+1-L)*CVYXX(L,K)+ARMC(L)*CVYXX(JS-1+L,K) SBXYB=SBXYB + ARMC(JS+1-L)*CVYXY(L,K)+ARMC(L)*CVYSY(JS-1+L,K) 290 SIGXX = (SAXX + 4.0*SBXXA + 2.0*SBXXB)/3.0SIGXY = (SAXY + 4.0*SBXYA + 2.0*SBXYB)/3.0ANSRX=REAL(SIGXX)*A ANSRY=REAL(SIGXY)*A ANSIX=AIMAG(SIGXX)*A

```
ANSIY=AIMAG(SIGXY)*A
      V=(4.0*PI*W*(ANSRX + ANSIY))/(D**2)
      S=(F**2)-((W*4.0*PI*(ANSIX-ANSRY))/(D**2))
      DET = (S^{**2}) + (V^{**2})
      IF (K-2*(K/2)) 15,16,15
  16
      PNO=-1.0
      GO TO 500
  15
      PNO = 1.0
      DARTR=PNO*V/DET
 500
      DARTI=PNO*S/DET
      IF (K-IPEG) 801,800,801
 800
      BETA = 0.0
      GAMMA = 0.0
 801
      BETA=BETA + DARTR
      GAMMA=GAMMA + DARTI
 700
      CONTINUE
      ETEII = (4.0 \text{W} \text{GAMMA}) / (Q \text{D})
      ETEIR=(4.0*W*BETA)/(Q*D)
      WRITE(6,900) ETEIR, ETEII, B
      FORMAT(1X,6HETEIR=,D14.3,4X,6HETEII=,D14.3,4X,2HB=,F10.2)
 900
1000
      CONTINUE
      REWIND 8
      END
```

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