Tunneling of correlated electrons in ultra high magnetic field

Shan-Wen Tsai\textsuperscript{1}, D. L. Maslov\textsuperscript{1} and L. I. Glazman\textsuperscript{2}

\textsuperscript{1)Institute for Fundamental Theory and}
Department of Physics, University of Florida, Gainesville, FL 32611
\textsuperscript{2)Theoretical Physics Institute, University of Minnesota, Minneapolis, MN 55455}

(Received: January 9, 2022)

Effects of the electron-electron interaction on tunneling into a metal in ultra-high magnetic field (ultra-quantum limit) are studied. The range of the interaction is found to have a decisive effect both on the nature of the field-induced instability of the ground state and on the properties of the system at energies above the corresponding gap. For a short-range repulsive interaction, tunneling is dominated by the renormalization of the coupling constant, which leads eventually to the charge-density wave instability. For a long-range interaction, there exists an intermediate energy range in which the conductance obeys a power-law scaling form, similar to that of a 1D Luttinger liquid. The exponent is magnetic-field dependent, and more surprisingly, may be positive or negative, i.e., interactions may either suppress or enhance the tunneling conductance compared to its non-interacting value. At energies near the gap, scaling breaks down and tunneling is again dominated by the instability, which in this case is an (anisotropic) Wigner crystal instability.

PACS numbers: 71.10Pm, 72.15Gd, 72.15

Low-dimensional systems exhibit “zero-bias anomalies” in tunneling (non-linearities of the current-voltage characteristics at small biases), which reflect the renormalization of the density of states by the electron-electron interaction. In particular, tunneling into a one-dimensional (1D) metal (Luttinger liquid) is characterized by a power-law suppression of the tunneling conductance, $g_T$. A three-dimensional metal placed in a strong magnetic field that depopulates all but one Landau levels (ultra-quantum limit, UQL) provides an example of a very special quasi-1D system. It is well-known\textsuperscript{1} that repulsive interactions in the UQL lead to a charge-density wave (CDW) or a Wigner crystal instability of the ground state. This has been confirmed, for example, by experiments on graphite in high magnetic fields.\textsuperscript{4} The most complete analysis of this instability for the case of a short-range interaction was performed in Ref. \textsuperscript{1} by solving the renormalization-group (RG) equation for the interaction vertex. On the other hand, it has recently been shown that for the case of a long-range (Coulomb) interaction $g_T$ exhibits a power-law, Luttinger-liquid-like behavior at sufficiently high energies.\textsuperscript{6} This result was obtained in two ways – via a perturbative Hartree-Fock procedure, which results in the RG equation for the transmission amplitude, and via bosonization in the coherent state basis – neither of which took into account the renormalization of the interaction vertex. In this paper, we combine the approaches of Refs. \textsuperscript{1} and \textsuperscript{5} to study the behavior of the tunneling conductance in the whole energy interval from the Fermi energy down to the energy associated with the instability. This is accomplished by solving a coupled system of the RG equations for both the interaction vertex and transmission amplitude. We find that for a long-range interaction there is a parametrically wide energy interval in which the flow of the transmission amplitude already results in its power-law scaling with energy, whereas the renormalization of the vertex is not yet important. In this interval, the system behaves as a Luttinger liquid and the results of Ref.\textsuperscript{5} are applicable. At energies close to the gap, the power-law scaling breaks down and tunneling is dominated by the electron properties in the vicinity of the critical point. For a short-range interaction, the critical point dominates tunneling at all energies, and there is no Luttinger-liquid behavior. We also find that above a certain magnetic field, when the lowest Landau level is strongly depopulated, the interaction enhances the tunneling conductance of the Luttinger-liquid regime above its free-electron value. Such an unusual behavior receives a natural explanation in terms of electron scattering from the Friedel oscillation near the tunneling barrier.

First we review the procedure of finding the effect of the electron-electron interaction on the tunneling conductance of a metal in the UQL.\textsuperscript{4} Let the magnetic field be perpendicular to the contact plane ($z = 0$) that separates two metallic sides. We consider both the symmetric configuration, where both sides are in the UQL, and the asymmetric one, where one of the sides is made of a high-carrier concentration and/or dirty metal so that, in the first approximation, the magnetic field does not affect that side. The transmission and reflection amplitudes for non-interacting electrons, $t_0$ and $r_0$, are assumed to be known. We choose the Landau basis for the free electron wavefunctions (non-interacting electrons in the UQL and in the presence of the barrier):

$$\psi^{(0)}_{p_z, p_x}(r) = \psi_{p_z}(z) \chi_{p_z}(x, y),$$

where

\begin{align}
\psi_{p_z}(z) &= \theta(z) t_0 e^{i p_z z} + \theta(-z) (e^{i p_z z} + r_0 e^{-i p_z z}) \\
\psi_{-p_z}(z) &= \theta(z) (e^{-i p_z z} + r_0 e^{i p_z z}) + \theta(-z) t_0 e^{-i p_z z}. \quad (1)
\end{align}

$\chi_{p_z}(x, y)$
Backscattering of magnetically quantized electrons at the barrier gives rise to a Friedel oscillation in the charge density, whose amplitude decays away from the barrier as $z^{-1}$. Correspondingly, the exchange, $V_{ex}(r, r')$, and Hartree, $V_H(r)$, potentials calculated using exact (in the presence of the barrier) but otherwise free wavefunctions, also exhibit rapid $2k_F$-oscillations and decay as $z^{-1}$ away from the barrier. The first-order correction to the wavefunction due to interaction is given by

$$\delta \Psi_{p_x,p_y}(r) = \int dr' \int dr'' G(r, r'; E) [V_{ex}(r', r'')] + \delta(r' - r'') V_H(r'')] \Psi^{(0)}_{p_x,p_y}(r''),$$

(2)

where $G(r, r'; E)$ is the Green’s function in the presence of the magnetic field. When the momentum transfer $(2p_z)$ matches the wavevector of the Friedel oscillation $(2k_F)$, a slow $(z^{-1})$ decay of the oscillations in $V_{ex}$ and $V_H$ leads to a logarithmic singularity in $\delta \Psi(r)$, and consequently, in $dt$. For electron’s energy close to the Fermi energy, we have

$$t = t_0 \left[1 - |r_0|^2 c \Gamma_0 \ln (W/E) \right]$$

(3)

where $c = 1(1/2)$ for symmetric (asymmetric) configuration, $W \sim E_F$ is the effective bandwidth, $\Gamma_0 \equiv \Gamma(q_\perp = 0)$, and $\Gamma(q_\perp)$ is expressed via the Fourier transform of the interaction potential $U_0(q_z, q_\perp)$:

$$\Gamma(q_\perp) = \frac{1}{(2\pi)^2 v_F} \int \frac{d^2 k_\perp}{2\pi} e^{i q_\perp \cdot k_\perp - k_\perp^2/2} U_0(0, k_\perp)$$

$$- U_0(2k_F, q_\perp) e^{-q_\perp^2/2}.$$  

(4)

The interaction potential, $U_0$, can be both of long- and short-range. The first situation corresponds to a single-band metal, in which there are no more charge carriers besides those already in the UQL. In this case, $U_0(0, q_\perp) = 4\pi e^2/(q^2 + \kappa^2)$ (for $q \ll k_F$), where $\kappa = \omega_p/v_F$ and $\omega_p$ is the plasma frequency at $B = 0$. The perturbation theory works if $c^2/v_F \approx \kappa^2 \ll 1$, which means that the interaction is long range. As the magnetic field reduces the phase-space available for the motion along the field, $k_F$ and $v_F$ decrease with the field (in the UQL, $k_F, v_F \propto B^{-1}$). Therefore, the perturbation theory breaks down at sufficiently strong field so that $\kappa \ll 1$. The case of a short-range interaction may correspond to a situation when a small pocket of the Fermi surface is in the UQL, whereas other parts of the Fermi surface still contain many Landau levels. Screening of the interaction among the magnetically quantized carriers is then mostly due to the “external” carriers, and the interaction may be of short-range.

The first and second terms in Eq. (4) are the exchange and Hartree contributions, respectively. Due to the Pauli principle, they enter Eq. (4) with opposite signs. Note that exchange contribution involves integration over the transverse momentum $k_\perp$, whereas the Hartree contribution enters $\text{Eq. (4)}$ at $q_z = 0$. This is due to a simple fact that the Friedel oscillations in charge density, and hence in the Hartree potential, is translationally-invariant along the barrier, whereas the exchange potential involves the density matrix and is thus non-local. If the interaction potential is peaked strongly at $q = 0$ (e.g., Coulomb potential) the Hartree contribution is enhanced due to $q_z = 0$ and may in fact dominate over the exchange one. In this case $\Gamma_0$ in Eq. (3) is negative and thus the transmission amplitude is enhanced by the interaction. Such an unusual behavior should be contrasted, e.g., to a strictly 1D case, in which $\Gamma_0 = U_0(0) - U_0(2k_F)$ and is thus positive for any “realistic” interaction potential, i.e., such that $U_0(q)$ is a monotonically decreasing function of $q$. In our case, due to the 3D nature of the problem, the exchange and Hartree contributions are not simply given by the corresponding forward and backscattering amplitudes but also involve averaging over the transverse direction. As a result, $\Gamma_0$ can take either sign.

At the next order, $V_{ex}$ and $V_H$ are re-calculated using the corrected wavefunctions and are substituted back into Eq. (3). The higher order corrections to $\delta \Psi(r)$ are higher powers of logs. Similarly to tunneling through a barrier of weakly interacting electrons in 1D, summation of the most divergent corrections to $t$ in all orders of the perturbation theory can be performed via an RG equation

$$\frac{dt}{d\xi} = -c \Gamma_0 t \left[1 - |t|^2 \right].$$

(5)

where $\xi \equiv \ln (W/E)$ and $t(\xi = 0) = t_0$. The solution of Eq. (5) is

$$t = \frac{t_0 (E/W)^{c \Gamma_0}}{\sqrt{|r_0|^2 + |t_0|^2 (E/W)^{2c \Gamma_0}}}.$$  

(6)

For small values of $|t_0|$ (high barrier), Eq. (6) reduces to a power-law form $t = t_0 (E/W)^{c \Gamma_0}$. The tunneling conductance is related to the transmission amplitude via the Landauer formula.

The power-law solution for $t$ is obtained using the bare interaction vertex. However, the vertex is also subject to renormalization. The flow of $\Gamma(q_\perp, \xi)$ is described by the integro-differential RG equation (4), which has a more transparent meaning when written for the Fourier transform of $\gamma(r_\perp, \xi) = 1/(2\pi)^2 \int d^2 q, \Gamma(q_\perp, \xi)e^{-i q_\perp \cdot r_\perp}$$

$$\frac{d \gamma(r_\perp, \xi)}{d \xi} = \int dr' \gamma(r', \xi) \gamma(r_\perp - r', \xi) \times \left(1 - e^{i r_\perp \cdot r'} \right),$$  

(7)

where $a \wedge b = a_x b_y - a_y b_x$. The initial condition for Eq. (4) is given by $\Gamma(q_\perp, \xi = 0) = \Gamma(q_\perp)$, where $\Gamma(q_\perp)$ is defined by Eq. (4). The flow of the vertex affects the RG equation for $t$, in which now the bare vertex has to be replaced by the renormalized one:
\[ \frac{dt}{d\xi} = -e\Gamma(q_\perp = 0, \xi) t (1 - |t|^2). \]  
\[ (8) \]

Eqs. (3), (7) and (8) provide a full description of tunneling into a three-dimensional metal in the UQL.

Few words about the general features of Eq. (8) are now in place. Physically, \( \gamma(r_\perp, 0) \) plays the role of an interaction potential between two electrons whose guiding centers are separated by a “distance” \( |r_\perp| \) across the magnetic field. (Note however that \( \gamma(r_\perp, 0) \) is related to the Fourier transform of the bare interaction, cf. Eq. (4).)

The vertex diverges most rapidly at \( q = 0 \), the CDW critical point dominates the physical words, the CDW critical point dominates the physical.

In contrast to Eq. (7), the full solution shows that the higher order correction to \( t(\xi) \) from Eq. (8), have already summed up into a power-law form but the renormalization of \( \Gamma \) is not significant yet. \( E_{PL} \) may be defined by the condition that the lowest-order correction to \( t_0 \) becomes equal to \( t_0 \):

\[ \xi_{PL} = \ln \left( W/E_{PL} \right) = 1/c\Gamma_0. \]

For a screened Coulomb potential, Eq. (10) together with the condition \( \kappa \ll k_F \leq 1 \) yields

\[ \Gamma_0 = \left( e^2/\pi v_F \right) \left[ \ln |\kappa| - (4k_F^2)^{-1} \right]. \]  
\[ (12) \]

A simple estimate for \( E_{PL} \) can be obtained if the first (exchange) term in Eq. (12) is larger than the second (Hartree) one, in which case \( E_{PL} \approx W \exp(-\pi v_F/ce^2) \ln |\kappa| \). At the same time, for \( k_F \approx 1 \), the gap is estimated as \( \Delta = W \exp(-a_{eF}/e^2) \), where \( a \approx 1 \). We see that \( E_{PL} \gg \Delta \) due to the presence of a large logarithm under the exponential.

As it has already been pointed out, \( \Gamma_0 \) is not necessarily positive: if \( k_F \) is sufficiently small (the lowest Landau level is strongly depleted), the Hartree (backscattering) term dominates and \( \Gamma_0 = 0 \). The dependence of \( \Gamma_0 \) on \( k_F \) is shown in the inset of Fig. 4 for a range of \( \kappa \). Recalling that both \( k_F \) and \( \kappa \) are magnetic-field-dependent, one can express the field \( B_0 \) at which \( \Gamma_0 \) changes sign in terms of the nominal field \( B_Q \) at which the UQL is achieved, and the familiar gas parameter, \( r_s \), as \( B_0 \approx 2B_Q/\ln^{1/3}C r_s^{-1} \), where \( C \approx 1 \). For \( r_s \approx 1 \) \( B_0 \approx 2B_Q \). The gap as a function of \( k_F \) and \( \kappa \) was obtained by solving the RG equation (4) numerically with Eq. (8) as an initial condition. The ratio \( \xi_e/\xi_{PL} \equiv \ln W/\Delta/\ln(W/E_{PL}) \) is plotted in Fig. 1. All numerical results presented so far are for a symmetric contact, i.e., \( e = 1 \). For a sufficiently long-range interaction (cf., filled circles for \( \kappa = 0.1 \) in Fig. 1) there is a wide range of \( k_F \) (and thus of the magnetic field) in which \( \xi_e/\xi_{PL} > 1 \). Once this condition is satisfied, there exists an energy interval \( \Delta \ll E \ll E_{PL} \xi_{PL} \ll \xi \ll \kappa_e \) in which the system effectively behaves as a 1D Luttinger liquid, exhibiting a characteristic power-law scaling of the tunneling conductance. (For \( E_{PL} \ll E \ll W \) the power-law reduces to the first log-correction.) Near the gap \( \Delta \), the power-law behavior crosses over to the threshold behavior [Eq. (11)]. If the interaction is
not sufficiently long-ranged, there is no energy interval in which $\xi_c/\xi_{PL} > 1$ is satisfied. This is illustrated by the open circles in Fig. 1 ($\kappa = 0.6$). (We did not plot the points corresponding to $k_F < \kappa$, as it would have been inconsistent with our weak-coupling approximation.)

Inset: $\Gamma (\text{open circles})$. The dashed line corresponds to the value $1$. Dashed lines: solutions of Eq. (8) with $\Gamma$ the open circles in Fig. 1 (in which $t$ is not sufficiently long-ranged, there is no energy interval in which $\xi_c/\xi_{PL} > 1$ is satisfied. This is illustrated by the open circles in Fig. 1 ($\kappa = 0.6$). (We did not plot the points corresponding to $k_F < \kappa$, as it would have been inconsistent with our weak-coupling approximation.)

FIG. 1. Ratio $\xi_c/\xi_{PL}$ for $\kappa = 0.1$ (filled circles) and $\kappa = 0.6$ (open circles). The dashed line corresponds to the value $1$. Inset: $\Gamma_0/(e^2/\pi v_F)$ as function of $k_F$ for various values of $\kappa$ (from top to bottom: $\kappa = 0.1, 0.2, 0.3, 0.4, 0.5, 0.6$).

FIG. 2. Solid lines: solution of the full system of RG equations Eqs. (8, 7, 4) for the screened Coulomb potential. Dashed lines: solutions of Eq. (8) with $\Gamma_0$ given by its bare value. In the main plot $t_0 = 0.1, k_F = 0.8$ and the values of $\kappa$ are indicated in the figure. Inset: $k_F = 0.4$. The curves, from bottom to top, correspond to $\kappa = 0.1, 0.2, 0.3$ and $0.4$.

A full solution of Eqs. (8, 7, 4) illustrating various regimes of the $t(E)$-dependence for a screened Coulomb potential is represented by solid lines in Fig. 2 for $t_0 = 0.1$. Dashed lines are solutions of Eq. (8) only with $\Gamma_0$ given by its bare value [Eq. (8)], i.e., without taking into account the renormalization of the coupling constant. In the main panel, $k_F = 0.8$. The region of small $\xi$ shows a good agreement between the full solution (solid lines) and the Luttinger-liquid solution (dashed lines). As $E$ approaches $\Delta$ ($\xi \rightarrow \xi_c$), scaling breaks down and $t(E)$ vanishes at $E = \Delta$. In the inset of Fig. 2, $k_F = 0.4$. The rise of $t$ with $\xi$ corresponds to the negative value of the bare vertex, $\Gamma_0$. This corresponds to an enhancement of $g\xi$ above its non-interacting value. Had the vertex renormalization been absent, $t$ would have reached unity in the limit $E \rightarrow 0$ ($\xi \rightarrow \infty$). However, the vertex renormalization curbs the rise in $t$ and eventually brings it down to zero at $E = \Delta$. At the instability, the system becomes an insulator. For long-range interaction, we find that the vertex diverges most rapidly at finite $q_\perp$ (for $\kappa \rightarrow 0, q_\perp \rightarrow 0.7$). This is consistent with previous works and suggests that in this case the instability is of the Wigner-crystal type, with charge modulation both along the field and in the transverse directions.

In conclusion, we found that the range of the electron-electron interaction plays a crucial role both in the nature of a field-induced instability and in tunneling at energies above the corresponding gap. For a short-range interaction, there is a CDW instability which dominates tunneling in the whole energy interval. For a long-range interaction, depending on the depletion of the lowest Landau level, tunneling may exhibit two types of characteristic behavior. For not too strong depletion, the tunneling conductance decreases as a power-law of the relevant energy scale at higher energies and vanishes in a threshold-like manner near the gap (which for the long-range case is the Wigner-crystal gap). For a stronger depletion, the conductance first increases as a power of energy, then goes through a maximum and finally vanishes at the gap. These predictions are amenable to a direct experimental verification by tunneling into low-carrier-density materials, e.g., graphite.

We are grateful to E. V. Sukhoroukov and V. M. Yakovenko for very instructive discussions. D. L. M. acknowledges support from the NHMFL In-House Research Program, NSF Grant No. DMR-9703388 and Research Corporation (RI0082). L. I. G. acknowledges support from NSF Grant No. DMR-9731756.

1 S. A. Brazovskii, Zh. Eksp. Teor. Fiz. 61, 2401 (1971) [Sov. Phys. JETP 34, 1286 (1972)].
2 H. Fukuyama, Solid State Comm. 26, 783 (1978).
3 V. M. Yakovenko, Phys. Rev. B 47, 8851 (1993).
4 Y. Iye, et al., Phys. Rev B 25, 5478 (1982); Y. Iye and G. Dresselhaus, Phys. Rev. Lett. 54, 1182 (1985); H. Yaguchi and J. Singleton, Phys. Rev. Lett. 81, 5193 (1998).
5 C. Biagini, D. L. Maslov, M. Yu. Reizer and L. I. Glazman, Europhys. Lett. 55, 383 (2001).
6 D. Yue, L. I. Glazman and K. A. Matveev, Phys. Rev. B 49, 1966 (1994).