





Superconductivity in the uniform electron gas: Irrelevance of the Kohn-Luttinger mechanismXiansheng Cai ¹, Tao Wang ¹, Nikolay V. Prokof'ev ^{1,*}, Boris V. Svistunov^{1,2,†} and Kun Chen ^{3,‡}¹*Department of Physics, University of Massachusetts, Amherst, Massachusetts 01003, USA*²*Wilczek Quantum Center, School of Physics and Astronomy, Shanghai Jiao Tong University, Shanghai 200240, China*³*Center for Computational Quantum Physics, Flatiron Institute, 162 5th Avenue, New York, New York 10010, USA*

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We study the Cooper instability in jellium model in the controlled regime of small to intermediate values of the Coulomb parameter $r_s \leq 2$. We confirm that superconductivity naturally emerges from purely repulsive interactions described by the Kukkonen-Overhauser vertex function. By employing the implicit renormalization approach and the discrete Lehmann representation we reveal that even in the small- r_s limit, the dominant mechanism behind Cooper instability is based on dynamic screening of the Coulomb interaction—accurately captured by the random phase approximation, whereas the Kohn-Luttinger contribution is negligibly small and, thus, not relevant.

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Introduction. Conventional BCS theory predicts Cooper instability at low temperature in a Fermi liquid with weak short range attractive coupling between the low-energy electrons, typically originating from the electron-phonon interaction (EPI). With the discovery of unconventional superconducting systems, such as *d*-wave cuprate superconductors [1–4], s^{+-} Fe-based superconductors [5–8], multilayer graphene systems [9–12], etc., alternative theories of Cooper instability have been drawing great research interest. In the past decades, a number of different mechanisms were proposed for the pairing instability in systems where the EPI alone was not sufficient to explain the data [13–17]. Some of them are based on purely repulsive bare electron-electron interactions [13–15]. In this work, we revisit superconducting properties of the uniform electron gas (jellium model) where pairing instability originates from purely repulsive Coulomb interparticle interactions and Fermi energy is the only relevant energy scale (all other energy scales are emergent). Our prime focus is the quantitative study of two canonical scenarios emerging from renormalized interactions: the Kohn-Luttinger (KL) mechanism based on the $2k_F$ singularity (where k_F is the Fermi momentum) and the dynamic screening mechanism.

In 1965, Kohn and Luttinger argued that for any weak short-range repulsive interaction, the two-particle effective interaction induced by many-body effects always becomes attractive at large enough orbital momenta $\ell \gg 1$, and could lead to Cooper instability [13]. They used the same analysis to estimate an effective Cooper channel coupling for the static screened Coulomb interaction. The KL mechanism has motivated a series of theoretical attempts to explain unconventional superconducting systems [18–21], and is widely

believed to be the dominant mechanism leading to superconductivity in jellium at small r_s . However, a thorough investigation of the dominant mechanism leading to superconductivity in this system is still missing: (i) the KL approach is based on the static screened potential, which is an uncontrolled approximation; (ii) we are not aware of any precise numerical study of the KL mechanism because it only emerges at large ℓ and leads to extremely small values of T_c ; (iii) whenever Cooper instability is observed in the simulation, one needs to differentiate between different scenarios behind it by revealing and evaluating their contributions separately. Therefore, whether the KL mechanism ever becomes dominant in the uniform electron gas in the high-density limit $r_s \rightarrow 0$ is still an unsolved fundamental question.

It has been known for decades that dynamic screening of the Coulomb interaction could also induce the pairing instability in jellium. Early work by Tolmachev and Bogoliubov [22] demonstrated that even if the Cooper channel coupling is repulsive at all frequencies, after its high-frequency part is renormalized to a smaller value the net result might be an attractive low-frequency effective potential. Later, Takada and others calculated the critical temperatures T_c of jellium numerically using various forms of dynamically screened Coulomb interaction [14,15,23,24]. Subsequent studies also reported that dynamic screening plays an important role in the superconductivity of metallic hydrogen and alkali metals [25,26] as well as in the dilute electron gas [16]. The superconducting phase diagram produced by Takada [24] stated that jellium enters a normal phase at $r_s < 2.0$, in contradiction with the KL prediction. Apparently, the values of T_c at large ℓ were too small to be resolved and, therefore, were ignored for practical purposes. Because of this limitation, the KL and dynamic screening mechanisms have never been quantitatively compared to each other, despite their coexistence in the uniform electron gas.

To determine which mechanism is dominant and under what conditions, we study the Cooper instability in the con-

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trolled regime of small to intermediate values of $r_s \leq 2$. The particle-particle irreducible four-point vertex is approximated with the Kukkonen-Overhauser (KO) ansatz [27], which becomes exact in the high-density limit $r_s \rightarrow 0$. We compare contributions from both mechanisms in two different ways. In the first protocol, we compute the largest eigenvalues, $\lambda(T)$, of the gap equation down to $T/E_F = 10^{-6}$ (below we use Fermi energy E_F as the unit of energy). We then remove the $q = 2k_F$ singularity in the polarization function, on which the KL mechanism is based, and measure the relative change of the eigenvalue, $\eta(T) = \delta\lambda/\lambda$. Finally, we estimate the magnitude of $\eta(T_c)$, which represents the relative contribution of the KL mechanism. This perturbative treatment is justified if $\eta(T)$ is small, which turns out to be always the case. The second protocol is based on the implicit renormalization (IR) approach [28]. By integrating out the high-frequency/energy degrees of freedom the IR approach solves a new eigenvalue problem for which the largest eigenvalue, $\bar{\lambda}(T)$, is also equal to unity at $T = T_c$. The crucial advantage of looking at $\bar{\lambda}(T)$ instead of $\lambda(T)$ is that its temperature dependence is a linear function of $\ln(T)$ for a properly chosen energy separation scale and, thus, can be accurately extrapolated to T_c from $T \gg T_c$. Computational costs are further dramatically reduced by employing the discrete Lehmann representation (DLR) [29]. This combination of methods is what allows us to determine the superconducting channel ℓ_c with the highest value of T_c at $r_s = 0.33, 0.5, 1, 2$; otherwise, the problem cannot be solved using standard techniques. By computing the critical values of orbital channels, $\ell_{KL}(r_s)$, when the KL mechanism first induces an attractive Cooper channel coupling, and comparing the asymptotic behavior of ℓ_c and ℓ_{KL} at small r_s we determine what mechanism is dominating in the high-density limit.

Within the first protocol we find the KL mechanism contribution remains extremely small, and therefore irrelevant, for any value of r_s when the transition temperatures exceed 10^{-10} . The second protocol reveals that the Cooper instability in jellium takes place at all values of r_s tested. More importantly, the dominant channel ℓ_c increases much slower than the critical channel of the KL mechanism, ℓ_{KL} , as r_s decreases, indicating that in the high-density limit superconductivity is induced by the dynamic screening effects accurately captured by the random phase approximation, well before the KL mechanism could have any impact.

Model. Jellium model is defined by the Hamiltonian

$$H = \sum_{\mathbf{k}\sigma} \epsilon_{\mathbf{k}} a_{\mathbf{k}\sigma}^\dagger a_{\mathbf{k}\sigma} + \frac{1}{2} \sum_{\substack{\mathbf{q} \neq 0 \\ \mathbf{q}, \mathbf{k}, \mathbf{k}', \sigma, \sigma'}} V_{\mathbf{q}} a_{\mathbf{k}+\mathbf{q}\sigma}^\dagger a_{\mathbf{k}'-\mathbf{q}\sigma}^\dagger a_{\mathbf{k}\sigma} a_{\mathbf{k}'\sigma}, \quad (1)$$

with $a_{\mathbf{k}\sigma}^\dagger$ the creation operator of an electron with momentum \mathbf{k} and spin $\sigma = \uparrow, \downarrow$, dispersion $\epsilon_{\mathbf{k}} = \frac{k^2}{2m_e} - \mu$, and Coulomb potential $V_{\mathbf{q}} = \frac{4\pi e^2}{q^2}$. The dimensionless coupling parameter (the Wigner-Seitz radius) is given by $r_s = \frac{1}{a_0} \left(\frac{3}{4\pi n} \right)^{\frac{1}{3}}$, where n is the number density and a_0 is the Bohr radius. The gap function equation reads

$$\lambda(T) \Delta_{\omega_n, \mathbf{k}} = -T \sum_m \int \frac{d\mathbf{p}}{(2\pi)^d} \Gamma_{\omega_m, \mathbf{p}}^{\omega_n, \mathbf{k}} G_{\omega_m, \mathbf{p}} G_{-\omega_m, -\mathbf{p}} \Delta_{\omega_m, \mathbf{p}}. \quad (2)$$

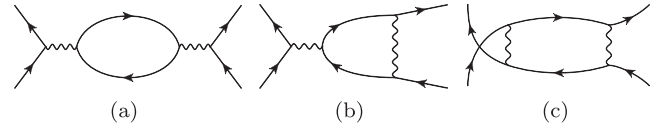


FIG. 1. Second-order diagrams contributing to the particle-particle irreducible four-point vertex Γ in the KL analysis. Only the first bubble diagram is resummed within the RPA approach, while all three diagrams are resummed in the KO formulation. In the small- r_s limit, the contributions of the diagrams (b) and (c) might prove appreciable only if the Kohn-Luttinger mechanism is the leading channel of Cooper instability; otherwise, the leading contribution is accurately captured by RPA.

Here Γ is the particle-particle irreducible four-point vertex, G is the single particle Green's function, Δ is the gap function, and λ is its eigenvalue. The key approximation used in this work is the Kukkonen-Overhauser ansatz [27,30] for Γ :

$$V_{\vec{\sigma}\vec{\sigma}'}^{KO}(\omega, q) = V_+ + V_+(q)^2 Q_+(\omega, q) + V_-(q)^2 Q_-(\omega, q) \vec{\sigma} \cdot \vec{\sigma}', \quad (3)$$

with

$$Q_{\pm}(\omega, q) = -\frac{\Pi_0(\omega, q)}{1 + V_{\pm}(\omega, q)\Pi_0(\omega, q)}, \quad (4)$$

$$V_+ = (1 - G_+)V, \quad V_- = -G_-V. \quad (5)$$

It is defined in terms of the polarization function Π_0 (based on the convolution of bare Green's functions) and local field factors $G_{\pm}(q)$. When $G_{\pm}(q)$ are set to zero, the KO interaction reduces to the random phase approximation (RPA). The local field factors encode the many-body exchange and correlation effects beyond RPA. For direct comparison with previous work by Takada we adopt the same ansatz for $G_{\pm}(q)$ [23,24] and take the functional form of Π_0 to be that at $T = 0$:

$$\begin{aligned}
 \Pi_0(q, \omega) &\simeq \frac{mk_F}{2\pi^2} P\left(\frac{q}{2k_F}, \frac{m\omega}{qk_F}\right), \quad (6) \\
 P(z, u) &= 1 + \frac{1 - z^2 + u^2}{4z} \ln \frac{(1+z)^2 + u^2}{(1-z)^2 + u^2} \\
 &\quad - u \tan^{-1} \frac{2u}{u^2 + z^2 - z}. \quad (7)
 \end{aligned}$$

This is justified by the smallness of the critical temperature. The gap equation is decomposed into different orbital channels, ℓ , which are solved independently. The critical temperature T_c in each channel corresponds to the point where the largest eigenvalue $\lambda(T)$ equals unity. For every choice of the vertex function considered in this work the single particle self-energy was computed within the G_0W_0 approximation as in [24].

Kohn-Luttinger mechanism. Cooper instability in the KL theory is induced by the logarithmic singularity in the static effective interaction at $q = 2k_F$ [13]. Within the second-order perturbation theory it arises from diagrams shown in Fig. 1. After projecting static Γ to the ℓ th orbital channel,

$$W_{\ell}(k, p, \omega = 0) = \int_{-1}^1 P_{\ell}(\chi) \Gamma(k, p, \chi, \omega = 0) d\chi, \quad (8)$$

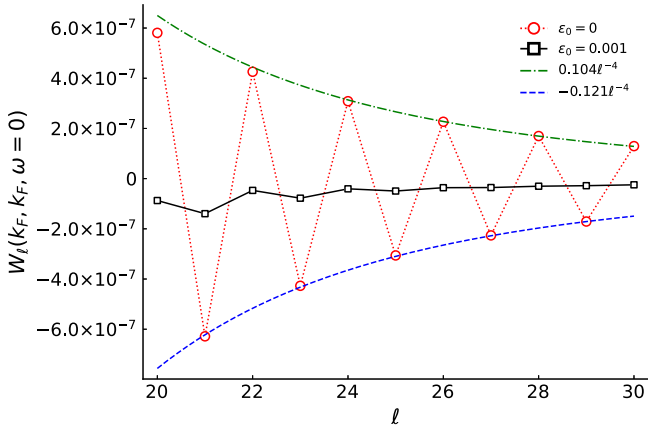


FIG. 2. Static KO vertex function on the Fermi surface $W_\ell(k_F, k_F, \omega = 0)$ at $r_s = 1.0$. Dotted red line with circles represents the original W_ℓ with clear KL oscillations decaying as ℓ^{-4} (dot-dash green and dashed blue fits). The solid black line with squares represents the regularized W_ℓ .

where $\chi = \cos \theta$, with θ the angle between momenta \mathbf{k} and \mathbf{p} , and $P_\ell(\chi)$ are Legendre polynomials, Kohn and Luttinger found that in the large- ℓ limit, $W_\ell(k_F, k_F, \omega = 0)$ decays as ℓ^{-4} and oscillates between the odd and even values of ℓ . The attractive effective coupling at large enough odd ℓ could then give rise to Cooper instability.

The KL treatment silently ignores the dynamic nature of screening despite the fact that at any finite frequency the Coulomb potential cannot be screened at small momenta $q \ll \omega/v_F$. Moreover, the singular nature of the Coulomb potential at small q calls for proper resummation of diagrams shown in Fig. 1 beyond the second-order perturbation theory. Clearly, a more controlled analysis is necessary to quantitatively evaluate the relative importance of KL and dynamic screening effects in the jellium model.

Resummation of diagrams shown in Fig. 1 is achieved within the KO vertex function that provides an excellent framework for thorough investigation of competing mechanisms. To quantify the contribution of the $q = 2k_F$ singularity, we introduce regularized $\Pi_\epsilon(q, \omega)$ that differs from $\Pi_\epsilon(q, \omega)$ by the replacement $(1-z)^2 + u^2 \rightarrow (1-z)^2 + u^2 + \epsilon(z)$ with $\epsilon(z) = \epsilon_0 e^{-4(z-1)^2}$. This modification is limited to the vicinity of the $2k_F$ singularity and ensures that changes in λ are attributed to the KL mechanism. The value of $\epsilon_0 = 0.001$ was chosen by establishing when $W_\ell(k_F, k_F, \omega = 0)$ for the regularized KO interaction has its odd-even channel oscillations suppressed; see Fig. 2. Next, we study the effect of ϵ_0 on the gap equation eigenvalues λ_ℓ .

Implicit renormalization approach. It is impossible to solve the gap equation directly when the critical temperature is extremely low because the number of required Matsubara frequency points is too large. Thus reliable extrapolation from $T \gg T_c$ is essential for determining where $\lambda(T) = 1$. Such an extrapolation can hardly be done for frequency-dependent vertices because $\lambda(T)$ turns out to be an unknown nonlinear function of $\ln T$. The implicit renormalization (IR) approach proposed in Ref. [28] offers a solution to this problem. The idea is to decompose the gap function into two complementary

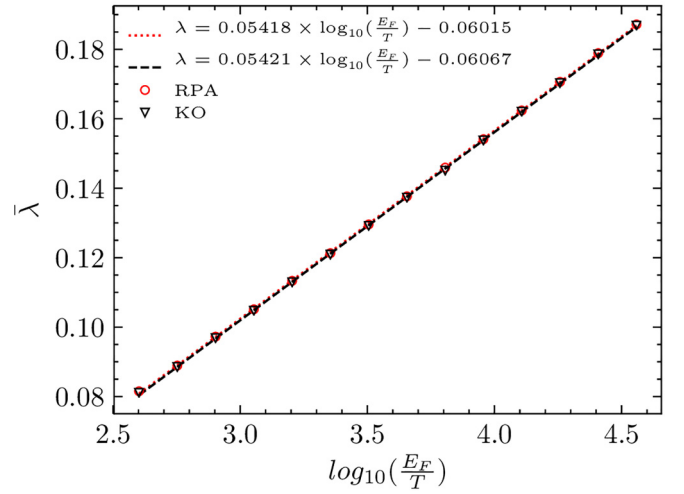


FIG. 3. Temperature dependence of the eigenvalues $\tilde{\lambda}(T)$ for RPA (red circles) and KO (black triangles) vertex functions at $r_s = 2.0$ and $\ell = 3$. The linear fits of the RPA (red dotted line) and KO (black dashed line) data are almost identical. The extrapolated value of $T_c = 2.71 \times 10^{-20} E_F$ is extremely small.

parts, $\Delta = \Delta^{(1)} + \Delta^{(2)}$, with $\Delta_n^{(1)} = 0$ for $|\omega_n| > \Omega_c$ and $\Delta_n^{(2)} = 0$ for $|\omega_n| < \Omega_c$, and solve an eigenvalue problem for the low-energy part $\Delta_n^{(1)}$ only. (The integration of high-energy degrees of freedom with the IR protocol is achieving the same goal as the pseudopotential theory.) The new eigenvalue $\tilde{\lambda}(T)$ is expected to have a nearly perfect linear dependence on $\ln T$ for a properly chosen energy scale separation. The IR approach allows us to accurately determine T_c as low as $10^{-20} E_F$ by extrapolating $\tilde{\lambda}(T)$ from the $10^{-5} < T/E_F < 10^{-3}$ interval. Once the Tolmachev-McMillan logarithm [15,22,31] is accounted for, the linear flow of $\tilde{\lambda}(T)$ illustrated in Fig. 3 provides direct access to the Coulomb pseudopotential μ^* in a given orbital channel ℓ . Note that the difference between μ^* of the KO and RPA vertex functions is smaller than a few percent for odd ℓ at $r_s \leq 2.0$ and even ℓ for $r_s \leq 1.0$. This indicates that higher-order vertex corrections for considered values of r_s are negligible.

Discrete Lehmann representation. Even within the IR protocol, solving the gap equation at exponentially low temperature faces technical challenges because the vertex function is a multidimensional object with nontrivial structure in momentum and frequency. The key step is optimization of frequency grids to store just enough information for accurate interpolation of functions. Fortunately, for a given ultraviolet cutoff ω_{\max} and numerical accuracy ϵ , the required grids are provided by the recently developed discrete Lehmann representation [29,32], or DLR. The number of grid points scales as $O(\ln(\omega_{\max}/T) \ln(\frac{1}{\epsilon}))$ and only 65 frequencies are required to achieve accuracy $\epsilon = 10^{-10}$ at $\omega_{\max}/T = 10^5$. Fast implementation of all key operations including Fourier transforms, interpolation, and convolution are available within the DLR [30]. By dramatically reducing memory and computational costs, the DLR grids allow us to simulate much lower temperatures and compute extremely small T_c . Specifically, we were able to determine the dominant superconducting orbital channel ℓ_c at $r_s \geq 0.33$.

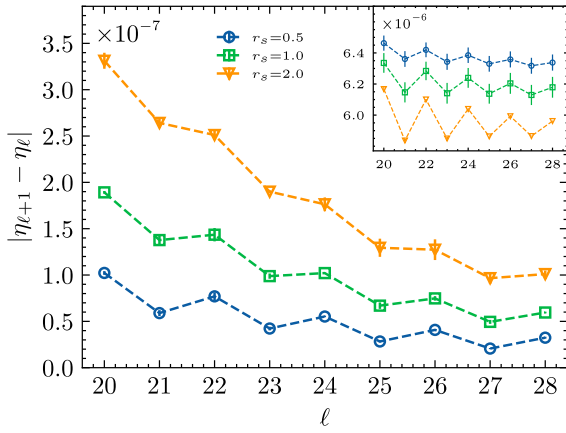


FIG. 4. Amplitudes of the relative eigenvalue oscillations, $|\eta_{\ell+1} - \eta_\ell|$, induced by regularization of the KO vertex at $T = 10^{-5}E_F$ and $r_s = 0.5$ (blue circles), $r_s = 1$ (green squares), and $r_s = 2$ (yellow triangles). Inset: relative eigenvalue changes η_ℓ . Data in the main figure and inset involve scaling factors of 10^{-7} and 10^{-6} , respectively.

Results. In the Fig. 4 inset we show the relative change of eigenvalues $\eta = \delta\lambda/\lambda$ at $T = 10^{-5}E_F$ when the KO function is regularized. We observe that KL oscillations are superimposed on a slowly decaying background originating from the vertex modification in finite vicinity of the $q = 2k_F$ point. The contribution from singularity is best characterized by the oscillation amplitude or the difference between the η_ℓ and $\eta_{\ell+1}$ values plotted in the lower panel of Fig. 4. We observe that (i) the relative contribution of the KL mechanism is $\sim 10^{-7}$ and (ii) the oscillation is less pronounced for smaller r_s . Both facts indicate that the KL mechanism is irrelevant for superconducting properties of jellium.

Our conclusions do not change with temperature as evidenced by simulation results shown in Fig. 5. The KL contributions $\eta_{\ell+1} - \eta_\ell$ increase when temperature decreases, but the rate is tiny and the curves tend to saturate. Even if $\eta_{\ell+1} - \eta_\ell$ were to grow at a constant rate beyond $T/E_F = 10^{-6}$, they would be smaller than a few percent at $T/E_F = 10^{-10}$, which is “zero” for all practical purposes.

Academically speaking, i.e., regardless of how small T_c is in the $r_s \rightarrow 0$ limit, the KL mechanism is not ruled out by results presented above because we cannot determine whether $\eta_{\ell+1} - \eta_\ell$ ultimately saturate to small values that decrease with r_s . However, recall that the KL mechanism is attractive only at $\ell \geq \ell_{KL}$, where $\ell_{KL}(r_s)$ is the first channel with $W_\ell(k_F, k_F, \omega = 0) < 0$. It has to be compared with the IR solution for the dominant superconducting channel, $\ell_c(r_s)$, for dynamic KO vertex function (this can be done for $1/3 \leq r_s \leq 2$). The comparison presented in Fig. 6 demonstrates that $\ell_{KL} > \ell_c$ for all r_s and the difference keeps growing when $r_s \rightarrow 0$. Thus superconductivity in the ℓ_c channel is induced by the dynamical screening well before the KL mechanism becomes viable, including the $r_s \rightarrow 0$ limit.

Conclusions. We studied superconductivity in the jellium model for $r_s \leq 2$ by solving the gap equation based on the Kukkonen-Overhauser vertex function. We find that superconductivity emerges from repulsive Coulomb interactions due to

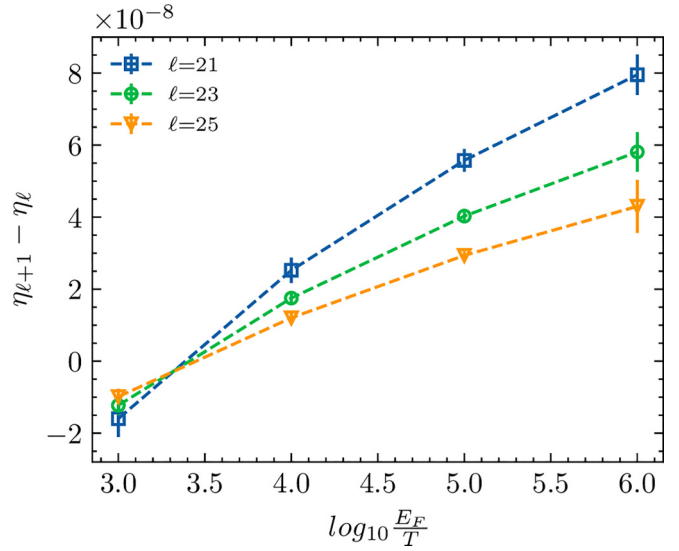


FIG. 5. Oscillation amplitudes as functions of $\log_{10}(E_F/T)$ for $r_s = 0.5$ and $\ell = 21$ (blue squares), $\ell = 23$ (green circles), and $\ell = 25$ (yellow triangles). When T/E_F is reduced by one order of magnitude, the amplitudes increase by less than 3×10^{-8} for $\ell \geq 21$ at $T \sim 10^{-6}E_F$.

dynamic screening effects. The Kohn-Luttinger mechanism, often assumed to be the prime reason behind Cooper instability in the high-density limit, is not relevant for two reasons: (i) for $r_s \leq 2$, it contributes to the Cooper instability only at orbital momenta $\ell \geq \ell_{KL}$ much larger than the dominant superconducting channel ℓ_c selected by the dynamic screening effect; (ii) for $\ell \geq \ell_{KL}$, the relative contribution of the KL mechanism is extremely small numerically and can be safely ignored. Since at $r_s \leq 1$ the physics of Cooper instability is accurately captured by RPA, one may attempt to find an analytic solution in the $r_s \rightarrow 0$ limit.

We solve the fundamental problem of Cooper instability in jellium in the high-density limit and revise a popular, yet incorrect, belief that the $q = 2k_F$ singularity is the key

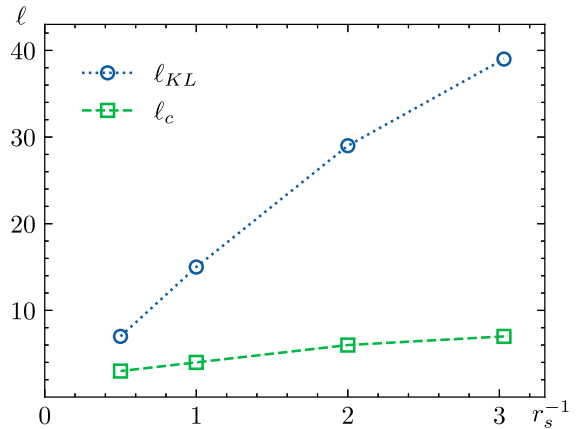


FIG. 6. Dominant superconducting channel ℓ_c (green squares connected by the dashed line) as a function of r_s , along with the critical channel of the KL mechanism ℓ_{KL} (blue circles connected by the dotted line).

reason. Our approach offers a systematic way for studies of the Cooper instability in other correlated electronic systems.

The symmetrized discrete Lehmann representation algorithm is a registered Julia package [32,33].

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