

BCS-BEC crossover in a quasi-two-dimensional Fermi gas

Andrea M. Fischer and Meera M. Parish

*Cavendish Laboratory, Department of Physics, University of Cambridge, JJ Thomson Avenue, Cambridge CB3 0HE, United Kingdom
and London Centre for Nanotechnology, Gordon Street, London WC1H 0AH, United Kingdom*

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We consider a two-component gas of fermionic atoms confined to a quasi-two-dimensional (quasi-2D) geometry by a harmonic trapping potential in the transverse direction. We construct a mean-field theory of the BCS-BEC crossover at zero temperature that allows us to extrapolate to an infinite number of transverse harmonic oscillator levels. In the extreme BEC limit, where the confinement length exceeds the dimer size, we recover 3D dimers confined to two dimensions with weak repulsive interactions. However, even when the interactions are weak and the Fermi energy is less than the confinement frequency, we find that the higher transverse levels can substantially modify fermion pairing. We argue that recent experiments on pairing in quasi-2D Fermi gases [Y. Zhang *et al.*, *Phys. Rev. Lett.* **108**, 235302 (2012)] have already observed the effects of higher transverse levels.

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I. INTRODUCTION

Recent experimental advances have made it possible to study quasi-two-dimensional (quasi-2D) atomic Fermi gases in a very controlled manner [1–8]. Such simple quasi-2D systems may provide useful insights into structurally complicated unconventional superconductors such as the cuprates, where superconductivity originates in the copper oxide planes [9]. Quasi-2D Fermi gases are also of fundamental interest since they are the marginal case of the Mermin-Wagner theorem and thus have modified superfluid properties [10,11]. In addition, the quasi-2D geometry can strongly affect fermion pairing within the superfluid, as we investigate in this paper.

In cold-gas experiments, atoms may be confined to one or more quasi-2D “pancake” structures using a 1D optical lattice. The system can then be tuned from three to two dimensions by increasing the confining lattice potential. For sufficiently strong lattices, the confining potential for a single quasi-2D layer can be modeled as a harmonic oscillator potential $V(z) = \frac{1}{2}m\omega_z^2 z^2$ in the transverse z direction, where m is the atom mass. When the temperature $k_B T \ll \hbar\omega_z$ and the Fermi energy $\varepsilon_F \ll \hbar\omega_z$, the atoms will reside in the lowest harmonic oscillator level in the absence of interactions and the gas is considered to be kinematically two dimensional. The short-range interatomic interactions are also renormalized to yield an effective 2D s -wave scattering amplitude and associated two-body bound state with a binding energy ε_B that depends on both the 3D scattering length and the confinement [12,13]. Thus, by varying the ratio $\varepsilon_B/\varepsilon_F$, one can explore the crossover in two dimensions from weak BCS pairing ($\varepsilon_B/\varepsilon_F \ll 1$) to the Bose-Einstein condensation (BEC) of dimers ($\varepsilon_B/\varepsilon_F \gg 1$) [14,15]. However, the interactions also mix in higher harmonic levels: For instance, in the BEC limit, dimers will be smaller than the confinement length $l_z = \sqrt{\hbar/m\omega_z}$ once $\varepsilon_B > \hbar\omega_z$, so they essentially become 3D bosons confined to quasi-two-dimensions. Here we are interested in how the confinement can impact pairing and lead to a departure from 2D behavior throughout the BCS-BEC crossover.

We focus on zero temperature, where there is a well-defined condensate in two dimensions, and we construct a

mean-field theory that generalizes the 2D results of Randeria *et al.* [14] to quasi-two-dimensions. We expect the mean-field approximation to be reasonable since it appears to be consistent with recent experiments in the 2D limit ($\varepsilon_F \ll \hbar\omega_z$) [7]. The BCS regime of the quasi-2D Fermi gas was previously studied in Ref. [16] using a mean-field Bogoliubov–de Gennes calculation that included the lowest three harmonic oscillator levels. Our approach, however, allows us to extrapolate to an infinite number of levels and thus explore the entire BCS-BEC crossover. In the limit $\varepsilon_B \gg \hbar\omega_z$, we find that our calculation recovers weakly repulsive 3D bosons confined to quasi-two-dimensions. However, even for weak interactions $\varepsilon_B \ll \hbar\omega_z$, we find that higher harmonic levels can substantially modify fermion pairing as we perturb away from pure two dimensions and ε_F approaches $\hbar\omega_z$. We determine the radio frequency (RF) spectra for the quasi-2D Fermi gas and show that recent measurements of pairing in the quasi-2D regime $\varepsilon_F \sim \hbar\omega_z$ [8] are consistent with effects due to higher transverse levels.

II. METHODOLOGY

We consider a two-component (\uparrow, \downarrow) Fermi gas interacting close to a broad s -wave Feshbach resonance in three dimensions. Under harmonic confinement in the z direction, the motion of each atom can be parametrized by its 2D momentum \mathbf{k} in the x - y plane and the harmonic oscillator quantum number n in the transverse direction. The many-body Hamiltonian is thus (setting the system volume to 1)

$$\hat{H} = \sum_{\mathbf{k}, n, \sigma} (\epsilon_{\mathbf{k}n} - \mu) c_{\mathbf{k}n\sigma}^\dagger c_{\mathbf{k}n\sigma} + \sum_{\substack{\mathbf{k}, n_1, n_2 \\ \mathbf{k}', n_3, n_4 \\ \mathbf{q}}} \langle n_1 n_2 | \hat{g} | n_3 n_4 \rangle c_{\mathbf{k}n_1 \uparrow}^\dagger c_{\mathbf{q}-\mathbf{k}n_2 \downarrow}^\dagger c_{\mathbf{q}-\mathbf{k}'n_3 \downarrow} c_{\mathbf{k}'n_4 \uparrow}, \quad (1)$$

where $\epsilon_{\mathbf{k}n} = k^2/2m + n\omega_z$ are the single-particle energies relative to the zero-point energy of the $n = 0$ state (we now set $\hbar = 1$) and μ is the chemical potential. Note that we assume that the mass and chemical potential (and thus the particle density) are the same for each spin σ .

Since the short-range interactions only depend on the relative motion, we obtain the interaction matrix elements $\langle n_1 n_2 | \hat{g} | n_3 n_4 \rangle$ by switching to relative and center-of-mass harmonic oscillator quantum numbers ν and N , respectively. This yields

$$\begin{aligned} \langle n_1 n_2 | \hat{g} | n_3 n_4 \rangle &= g \sum_N f_\nu \langle n_1 n_2 | N \nu \rangle f_{\nu'} \langle N \nu' | n_3 n_4 \rangle \\ &\equiv g \sum_N V_N^{n_1 n_2} V_N^{n_3 n_4}, \end{aligned} \quad (2)$$

where $f_\nu = \sum_{k_z} \tilde{\phi}_\nu(k_z)$, with $\tilde{\phi}_\nu$ the Fourier transform of the ν th harmonic oscillator eigenfunction. It is easily seen that $f_{2\nu+1} = 0$ and $f_{2\nu} = \frac{(-1)^\nu}{\nu!} \left(\frac{m\omega_z}{2\pi}\right)^{1/4} \sqrt{\frac{(2\nu)!}{2^{2\nu}}}$. The change of basis coefficients are given in Ref. [17]. Since $\langle n_1 n_2 | N \nu \rangle \sim \delta_{N+\nu, n_1+n_2}$ and ν and ν' are even, $n_1 + n_2$ must equal, modulo 2, $n_3 + n_4$ to obtain a nonzero interaction matrix element. The 3D contact interaction g can be written in terms of the binding energy ε_B of the two-body bound state, which always exists in the quasi-2D geometry:

$$-\frac{1}{g} = \sum_{\mathbf{k}, n_1, n_2} \frac{f_{n_1+n_2}^2 |\langle n_1 n_2 | 0 \ n_1 + n_2 \rangle|^2}{\epsilon_{\mathbf{k}n_1} + \epsilon_{\mathbf{k}n_2} + \varepsilon_B}. \quad (3)$$

Here we simply take $N = 0$ since ε_B is independent of the center-of-mass motion. One can also determine ε_B as a function of the 3D scattering length a_s [12,13] from Eq. (3) using $\frac{1}{g} = \frac{m}{4\pi} \left(\frac{1}{a_s} - \frac{2\Lambda}{\pi}\right)$, where Λ is a UV cutoff for the 3D momentum that can be sent to infinity at the end of the calculation.

Now if we define the superfluid order parameter

$$\Delta_{\mathbf{q}N} = g \sum_{\mathbf{k}, n_1, n_2} V_N^{n_1 n_2} \langle c_{\mathbf{q}-\mathbf{k}n_2\downarrow} c_{\mathbf{k}n_1\uparrow} \rangle \quad (4)$$

and assume fluctuations around this are small, we obtain the mean-field Hamiltonian

$$\begin{aligned} \hat{H}_{\text{MF}} &= \sum_{\mathbf{k}, n, \sigma} (\epsilon_{\mathbf{k}n} - \mu) c_{\mathbf{k}n\sigma}^\dagger c_{\mathbf{k}n\sigma} \\ &+ \sum_{\mathbf{q}, N} \left(\Delta_{\mathbf{q}N} \sum_{\mathbf{k}, n_1, n_2} V_N^{n_1 n_2} c_{\mathbf{k}n_1\uparrow}^\dagger c_{\mathbf{q}-\mathbf{k}n_2\downarrow}^\dagger \right. \\ &\left. + \Delta_{\mathbf{q}N}^* \sum_{\mathbf{k}', n_3, n_4} V_N^{n_3 n_4} c_{\mathbf{q}-\mathbf{k}'n_3\downarrow} c_{\mathbf{k}'n_4\uparrow} - \frac{|\Delta_{\mathbf{q}N}|^2}{g} \right). \end{aligned} \quad (5)$$

We further assume that the ground state has a uniform order parameter without nodes so that $\Delta_{\mathbf{q}N} = \delta_{\mathbf{q}0} \delta_{N0} \Delta_0$. In this case, Eq. (5) only contains a single unknown parameter Δ_0 , so it can be diagonalized to yield

$$\hat{H}_{\text{MF}} = \sum_{\mathbf{k}, n} (\epsilon_{\mathbf{k}n} - \mu - E_{\mathbf{k}n}) - \frac{\Delta_0^2}{g} + \sum_{\mathbf{k}, n, \sigma} E_{\mathbf{k}n} \gamma_{\mathbf{k}n\sigma}^\dagger \gamma_{\mathbf{k}n\sigma}, \quad (6)$$

where $E_{\mathbf{k}n}$ are the quasiparticle excitation energies. The quasiparticle creation and annihilation operators are respectively

given by

$$\gamma_{\mathbf{k}n\uparrow}^\dagger = \sum_{n'} (u_{\mathbf{k}n'n} c_{\mathbf{k}n'\uparrow}^\dagger + v_{\mathbf{k}n'n} c_{-\mathbf{k}n'\downarrow}), \quad (7)$$

$$\gamma_{-\mathbf{k}n\downarrow} = \sum_{n'} (u_{\mathbf{k}n'n} c_{-\mathbf{k}n'\downarrow} - v_{\mathbf{k}n'n} c_{\mathbf{k}n'\uparrow}^\dagger), \quad (8)$$

where the real amplitudes u and v only depend on the magnitude $k \equiv |\mathbf{k}|$ and satisfy $\sum_{n'} (|u_{\mathbf{k}n'n}|^2 + |v_{\mathbf{k}n'n}|^2) = 1$. Note that while they have a well-defined spin and momentum, they involve a superposition of different harmonic oscillator levels. The corresponding BCS wave function $|\Psi_{\text{MF}}\rangle \propto \prod_{\mathbf{k}n\sigma} \gamma_{\mathbf{k}n\sigma} |0\rangle$, where $|0\rangle$ is the vacuum state for the bare operators $c_{\mathbf{k}n\sigma}$. We then minimize $\langle \hat{H}_{\text{MF}} \rangle = \sum_{\mathbf{k}, n} (\epsilon_{\mathbf{k}n} - \mu - E_{\mathbf{k}n}) - \frac{\Delta_0^2}{g}$ with respect to Δ_0 at fixed μ to obtain the ground state. The value of μ is chosen to keep the density of particles $\rho = 2 \sum_{\mathbf{k}, n', n} |v_{\mathbf{k}n'n}|^2$ and thus the Fermi energy ε_F constant throughout the crossover.¹

A. Two-level calculation

It is first instructive to consider the case of only two levels $n = 0, 1$. Here Eq. (6) is greatly simplified since there is no pairing between atoms in the $n = 0$ and 1 levels. The quasiparticle dispersions are then $E_{\mathbf{k}n} = \sqrt{(\epsilon_{\mathbf{k}n} - \mu)^2 + (V_0^{nn} \Delta_0)^2}$. One can now minimize $\langle \hat{H}_{\text{MF}} \rangle$ by simply using $\partial \langle \hat{H}_{\text{MF}} \rangle / \partial \Delta_0 = 0$. Combining this with Eq. (3) yields the implicit equation

$$\left(\frac{\sqrt{\Delta^2 + \mu^2} - \mu}{\varepsilon_B} \right)^4 = \frac{\varepsilon_B + 2\omega_z}{\omega_z - \mu + \sqrt{\frac{\Delta^2}{4} + (\omega_z - \mu)^2}}, \quad (9)$$

where we have defined $\Delta = \Delta_0 V_0^{00}$, the pairing gap in the lowest level. Also, the density $\rho = -\partial \langle \hat{H}_{\text{MF}} \rangle / \partial \mu$ and in the regime $\varepsilon_F \leq \omega_z$ where $\varepsilon_F = \pi\rho/m$, this gives

$$2\varepsilon_F = \sqrt{\Delta^2 + \mu^2} + \sqrt{\Delta^2/4 + (\omega_z - \mu)^2} + 2\mu - \omega_z. \quad (10)$$

We see that Eqs. (9) and (10) reduce to the 2D mean-field equations [14] in the limit $\omega_z \rightarrow \infty$, as expected. They also yield the lowest-order correction to the 2D result due to confinement ($\varepsilon_F/\omega_z \neq 0$): In the BCS regime $\varepsilon_B/\varepsilon_F \ll 1$, we have

$$\frac{\Delta}{\varepsilon_F} \simeq \sqrt{\frac{2\varepsilon_B}{\varepsilon_F}} \left(1 + \frac{\varepsilon_F}{8\omega_z} \right), \quad \frac{\mu}{\varepsilon_F} \simeq 1 - \frac{\varepsilon_B}{2\varepsilon_F} \left(1 + \frac{\varepsilon_F}{4\omega_z} \right).$$

Thus Δ is enhanced by the confinement in this regime while μ is suppressed. This trend is also observed in the full calculation involving many levels (see Figs. 1 and 2).

B. Multiple levels

For more accurate results, we must include multiple levels of the confinement, especially when perturbing away from the 2D limit $\varepsilon_F, \varepsilon_B \ll \omega_z$. In general, \hat{H}_{MF} must be diagonalized

¹We define ε_F to be the chemical potential of an ideal Fermi gas with the same density ρ .

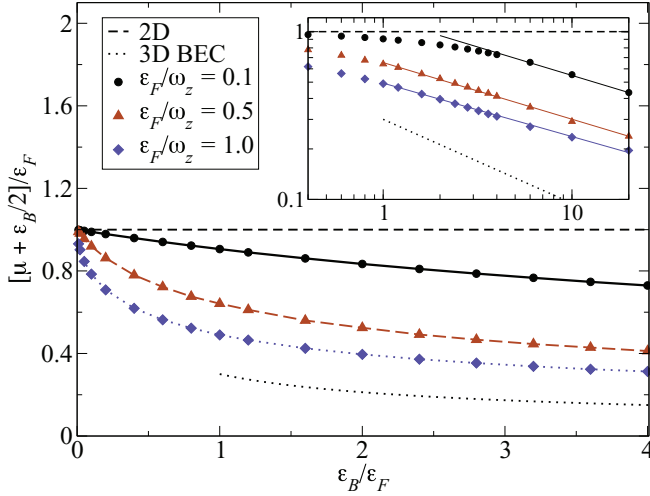


FIG. 1. (Color online) Chemical potential μ measured with respect to half the binding energy ε_B for several values of the Fermi energy ε_F/ω_z . The dashed line is the 2D mean-field result [14], while the dotted curve is the 3D BEC result, $\mu + \frac{\varepsilon_B}{2} \simeq -\frac{2\sqrt{2}\varepsilon_F}{3\pi} \sqrt{\frac{\varepsilon_F}{\varepsilon_B}}$. The inset shows the asymptotic behavior in the BEC regime $\varepsilon_B/\varepsilon_F > 1$ plotted on a logarithmic scale. Once $\varepsilon_B/\hbar\omega_z > 1$, we obtain 3D bosons confined to quasi-two-dimensions. The solid straight lines are straight line fits to the data with gradient $-1/3$. Error bars for the numerical data are within symbol size.

numerically to obtain E_{kn} and the quasiparticle amplitudes for a given μ and Δ . Equivalently, one can solve the Bogoliubov–de Gennes equations self-consistently, but it is considerably faster to minimize the energy $\langle \hat{H}_{MF} \rangle$ directly and it also allows us to take into account up to 100 levels. Indeed, we find that higher harmonic levels are important even for weak interactions once ε_F/ω_z shifts away from zero. For the values of ε_B and ε_F considered in this paper, we can in fact extrapolate the results for μ and Δ to an infinite number of harmonic levels since we find that they both scale linearly with the inverse of the number of levels in this limit. We emphasize that the extrapolated values and those obtained using only a few levels

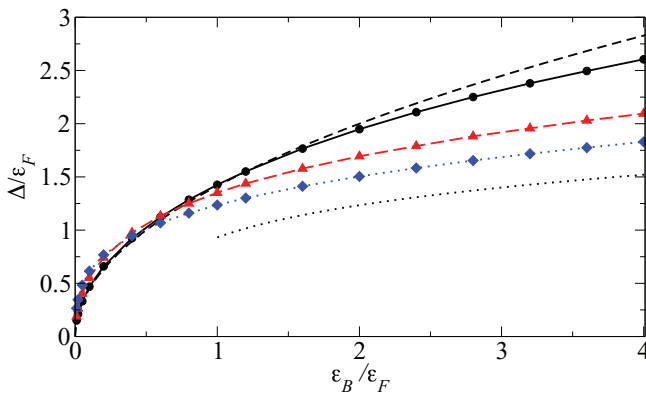


FIG. 2. (Color online) Behavior of the order parameter Δ throughout the BCS-BEC crossover for different values of ε_F/ω_z . The dashed curve is the 2D mean-field result $\Delta = \sqrt{2\varepsilon_B\varepsilon_F}$ [14], while the dotted curve is the 3D BEC result $\Delta \simeq \varepsilon_F \sqrt{\frac{16}{3\pi} \left(\frac{\varepsilon_B}{2\varepsilon_F}\right)^{1/4}}$. The key for the numerical data is the same as in Fig. 1.

can differ by tens of percent, even in the BCS regime (see the Appendix).

III. RESULTS

A. Order parameter and chemical potential

By incorporating an infinite number of levels, we can determine the evolution of the chemical potential μ and order parameter Δ throughout the BCS-BEC crossover in a quasi-2D Fermi gas, as depicted in Figs. 1 and 2. The lowest density we consider ($\varepsilon_F/\omega_z = 0.1$) corresponds approximately to the experiments of Refs. [3–6]. Note that in the extreme BCS and BEC limits, we must have $\mu \rightarrow \varepsilon_F$ and $\mu \rightarrow -\varepsilon_B/2$, respectively, a feature that holds across all dimensions. In the BCS regime, the behavior is in qualitative agreement with the two-level calculation: μ is suppressed and Δ is enhanced with respect to the 2D result, with the deviation from two dimensions being increased with increasing ε_F/ω_z . However, multiple levels are required to correctly capture the dependence on $\varepsilon_B/\varepsilon_F$ as ε_F shifts away from zero. We also see in Fig. 1 that the relative chemical potential $\mu + \varepsilon_B/2$ exhibits a very steep gradient as $\varepsilon_B \rightarrow 0$ for $\varepsilon_F/\omega_z \gtrsim 0.5$. This illustrates how higher levels can lead to a strong deviation from two dimensions even when $\varepsilon_F < \omega_z$ and $\varepsilon_B/\varepsilon_F \ll 1$.

For larger $\varepsilon_B/\varepsilon_F$, Δ eventually becomes suppressed compared to the 2D result and approaches the 3D mean-field curve in the BEC limit (Fig. 2). The chemical potential, in contrast, always remains lower than the 2D result and has a behavior in the BEC regime that is intermediate between 2D and 3D mean field. In the limit $\varepsilon_B/\varepsilon_F \gg 1$, the pairing gap is no longer given by Δ and the properties of the Bose superfluid are instead encoded in μ . In particular, the relative quantity $\mu + \varepsilon_B/2$ yields the mean-field energy for the repulsion between dimers. Referring to Fig. 1 (inset), we see that it tends to zero as a power law with increasing $\varepsilon_B/\varepsilon_F$, similarly to the 3D and in contrast to the 2D mean-field result. This is consistent with a dimensional crossover to 3D dimers once $\varepsilon_B \gtrsim \omega_z$. In this case, we expect to have weakly interacting bosons confined to quasi-two-dimensions with a mean-field energy that scales as $a_s/l_z \sim \sqrt{\omega_z/\varepsilon_B}$ [18], i.e., it behaves as a power law with exponent $-1/2$ like in the 3D BEC regime. However, our quasi-2D calculation gives a power of $-1/3$ rather than $-1/2$. This discrepancy is due to the fact that our mean-field approximation does not allow for the scattering of dimers in the transverse direction since they are constrained to be in the $N = 0$ center-of-mass mode. Removing a spatial degree of freedom means we only recover $2/3$ of the full exponent. The fact that we recover a power-law dependence at all for the dimer-dimer interaction is because we can extrapolate to an infinite number of levels. Note that $N > 0$ scattering between dimers is not expected to be significant in the BCS regime. The repulsion between dimers has also been discussed in the context of a two-channel model for the quasi-2D system [19], although no explicit dependence on ε_B was given.

B. Experimental probes

Deviations from 2D behavior will also be apparent in experimental probes of the quasi-2D superfluid. Typically,

investigations of pairing have exploited RF spectroscopy [20], where atoms in one hyperfine spin state (e.g., \downarrow) are transferred via an RF pulse to another hyperfine state that is initially unoccupied. In the ideal scenario where the final state is noninteracting, the mean-field transition rate or RF current is given by

$$I_{RF}(\omega) \propto \sum_{\mathbf{k}, n', n} |v_{\mathbf{k}n'n}|^2 \delta(\epsilon_{\mathbf{k}n'} - \mu + E_{\mathbf{k}n} - \omega), \quad (11)$$

where ω is the frequency shift relative to the bare transition frequency between hyperfine states. Here the onset frequency of the RF spectrum corresponds to $E_{\mathbf{k}=0, n=0} - \mu$ and is associated with the pairing gap of the superfluid. In the 2D case, (11) reduces to $I_{RF}(\omega) \propto \frac{\Delta^2}{\omega^2} \Theta(\omega - \epsilon_B)$ and thus the RF pairing gap is simply ϵ_B , as noted by Sommer *et al.* [7]. Perturbing away from two dimensions, we find that the RF spectrum can be substantially modified by the higher confinement levels. In Fig. 3 we see that the strongest effects are in the BCS regime, where the pairing gap is initially enhanced compared to ϵ_B . Such an enhancement is not surprising given that, in three dimensions, a pairing gap can exist even when there is no two-body bound state. However, once $\epsilon_F \gtrsim \omega_z$, the pairing gap drops below ϵ_B and even becomes *negative* for small enough ϵ_B/ϵ_F . This is because the coupling between $n = 0$ and 2, and the associated level repulsion (see inset of Fig. 3), reduces the energy $E_{\mathbf{k}, n=0}$. In this case, the lowest-energy quasiparticle contains a smaller fraction of the $n = 0$ harmonic level and the RF peak is instead dominated by the $n = 2$ quasiparticle. Thus the onset frequency is no longer an accurate measure of pairing, as we can see in Fig. 4. Note that the deep lattices used in Ref. [7] correspond to $\epsilon_F/\omega_z \approx 0.03$ and thus the pairing gap will lie very close to the 2D result, as was observed.

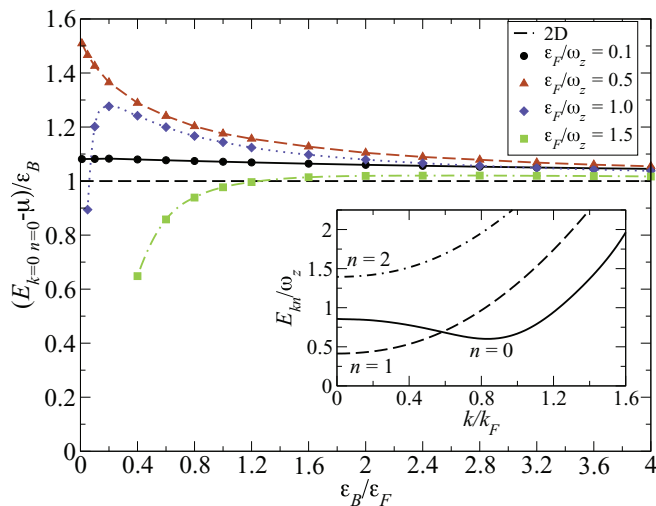


FIG. 3. (Color online) Pairing gap measured using RF spectroscopy as a function of ϵ_B/ϵ_F for a range of different ϵ_F/ω_z . In the BCS regime, it deviates substantially from the 2D result with increasing ϵ_F/ω_z . In the BEC limit, $E_{\mathbf{k}=0, n=0} - \mu$ must always approach ϵ_B , regardless of ϵ_F/ω_z . The inset shows the lowest quasiparticle dispersions for $\epsilon_F/\omega_z = 1$, $\epsilon_B/\epsilon_F = 0.1$. Note that there is level repulsion between $n = 0$ and 2, but no avoided crossing between the $n = 0$ and 1 dispersions.

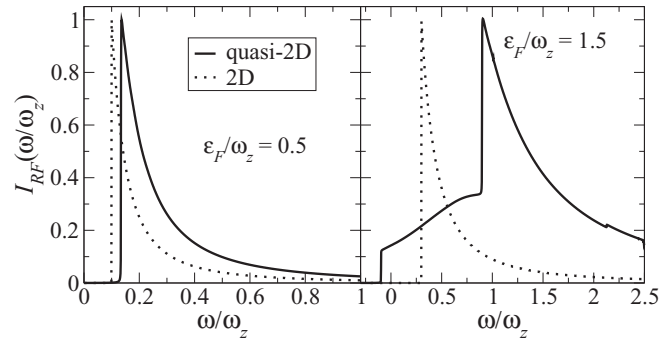


FIG. 4. Radio frequency spectra of the quasi-2D Fermi gas in the BCS regime with $\epsilon_B/\epsilon_F = 0.2$ for $\epsilon_F/\omega_z = 0.5, 1.5$. The RF current I_{RF} is scaled so that the peak value is always 1. The peak is strongly shifted towards higher frequencies compared to the 2D result when $\epsilon_F/\omega_z \simeq 1.5$. Note that the spectra are calculated assuming that the final state is noninteracting.

In contrast, the experiments in Ref. [8] correspond to $\epsilon_F/\omega_z \simeq 1.5$ and thus the RF spectrum in the BCS regime will be strongly affected by the confinement. In particular, we see in Fig. 4 that the RF peak is shifted to higher frequencies compared to the 2D case and develops more structure at lower frequencies. Furthermore, the pairing gap in the BEC regime appears to be less sensitive to confinement and closer to the 2D result (Fig. 3) since it is dominated by two-body physics. These features are all consistent with the experimental observations [8]. A direct comparison with these experiments is not straightforward in the BCS regime because of strong final-state interactions (which we have neglected). In addition, thermal effects are expected to smear out the fine structure in the RF spectra in Fig. 4, leaving an RF peak that more closely resembles that observed. It has even been suggested that finite temperature plays a crucial role here; see Ref. [21] for an alternative explanation based on fermionic polarons. However, the shift due to confinement appears to be substantial at $T = 0$, with a direction that is consistent with experiment, and therefore it cannot be disregarded. Indeed, it has also been shown that effects due to confinement can be significant in spin-polarized quasi-2D Fermi gases [22].

IV. CONCLUSION

To conclude, we have constructed a mean-field theory for the quasi-2D Fermi gas that is able to capture the deviations from 2D behavior resulting from confinement. We expect it to provide a benchmark for further investigations into quasi-2D Fermi systems. In the future, it would be interesting to explore the effects of finite temperature and the Berezinskii-Kosterlitz-Thouless transition in our model.

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APPENDIX: EFFECT OF HIGHER LEVELS

Our approach of directly minimizing the mean-field energy, enabled us to do calculations for up to 100 levels. Indeed, changing the number of levels by one does not change the results by more than a few percent, as noted in Ref. [16]. However, there is an incremental change if one continues to add more and more levels, resulting in an overall significant difference. This is to be expected in the BEC regime, but surprisingly also holds in the BCS regime. Upon plotting μ and Δ as a function of the inverse number of levels used in the calculation, we observe a linear behavior (see Fig. 5). This enables us to extrapolate to the origin, corresponding to an infinite number of harmonic levels. For $\varepsilon_F/\omega_z = 1$, as shown in Fig. 5, Δ changes by 32% for $\varepsilon_B/\omega_z = 0.01$ and 17% for $\varepsilon_B/\omega_z = 0.1$, when the extrapolated values are compared to those obtained in a three-level calculation. For μ , the differences are 3% for $\varepsilon_B/\omega_z = 0.01$ and 13% for $\varepsilon_B/\omega_z = 0.1$. These results are for the BCS regime. For the BEC regime, the percentage changes between a few levels and the extrapolated values are considerably larger.

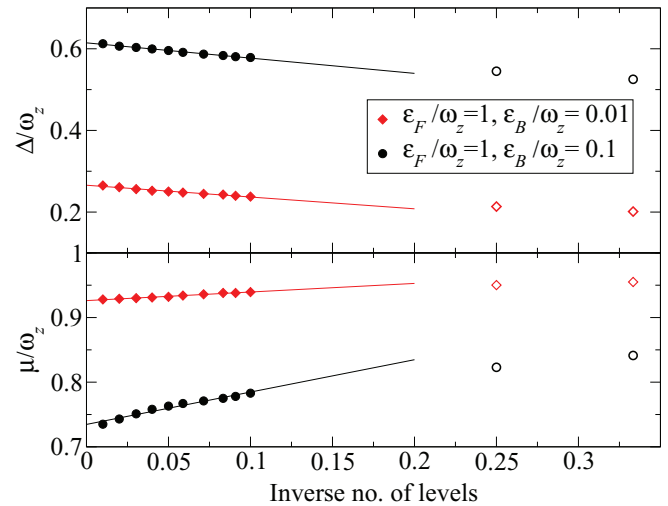


FIG. 5. (Color online) Order parameter Δ (upper panel) and chemical potential μ (lower panel) versus inverse number of levels for $\varepsilon_F/\omega_z = 1$ and $\varepsilon_B/\omega_z = 0.01$ (red diamonds) $\varepsilon_B/\omega_z = 0.1$ (black circles). The solid lines are linear fits to the data for ten or more levels. The open symbols, corresponding to the three- and four-level cases, have the same values of $\varepsilon_B, \varepsilon_F$ as their solid counterparts, but were not included in the fit.

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