Necessary and sufficient condition for quantum adiabaticity in a driven one-dimensional impurity-fluid system

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We study under what conditions the quantum adiabaticity is maintained in a closed many-body system consisting of a one-dimensional fluid and an impurity particle dragged through the latter by an external force. We employ an effective theory describing the low-energy sector of the system to derive the time dependence of the adiabaticity figure of merit: the adiabatic fidelity. We find that in order to maintain adiabaticity in a large system the external force \( F_N \) should vanish with the system size \( N \) as \( 1/N \) or faster. This improves the necessary adiabatic condition \( F_N = O(1/ \ln N) \) obtained for this system previously [O. Lychkovskiy et al., in Fourth International Conference on Quantum Technologies (ICQT-2017), AIP Conf. Proc. No. 1936 (AIP, New York, 2018), p. 020024]. The experimental implications of this result and its relation to the quasi-Bloch oscillations of the impurity are discussed.

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I. INTRODUCTION

The quantum adiabatic theorem is a fundamentally important result in the theory of quantum systems with time-dependent Hamiltonians. In essence, it states that a system initially prepared in an instantaneous eigenstate of a Hamiltonian remains arbitrarily close to the (time-evolving) instantaneous eigenstate provided the ramp rate (i.e., the rate of change of the Hamiltonian) is slow enough [1,2]. When it comes to applying the adiabatic theorem in practice, the key question to be addressed is how slow is slow enough. While this question can be exhaustively answered for a simple two-level system [3,4], it becomes complicated for many-body systems and/or for continuous quantum systems with infinite-dimensional Hilbert spaces. Although numerous sufficient conditions for adiabaticity are known (see the pioneering work in [5] and the review in [6]), they often prove to be inapplicable for continuous quantum systems due to the divergence of operator norms entering these conditions. Recently, a necessary condition for adiabaticity was proved [7] which is free from this shortcoming and is well suited for applying to many-body systems. Anyway, any adiabatic condition, whether sufficient or necessary, provides only a bound on the driving rate, without an indication of how tight this bound is.

In the present paper we derive a necessary and sufficient condition for quantum adiabaticity in a many-body system consisting of a one-dimensional fluid and an impurity particle dragged through the latter by a constant external force. This system exhibits a spectacular phenomenon predicted in Refs. [8–10] and experimentally observed in Ref. [11]: quasi-Bloch oscillations of the impurity’s velocity and position. These oscillations are somewhat reminiscent of the Bloch oscillations of a single particle in a periodic potential. Their root cause is the nontrivial spectral edge of the impurity-fluid system in one dimension, which is periodic in the thermodynamic limit as a function of the total momentum, with the period \( 2 \pi \rho \) determined by the density of the fluid \( \rho \equiv N/L \) (here \( N \) is the number of particles of the fluid, \( L \) is the linear dimension of the system, and we set \( \hbar = 1 \) throughout the paper). There are two important features, however, distinguishing quasi-Bloch oscillations from the conventional Bloch oscillations. First, intriguingly, quasi-Bloch oscillations occur in the translation-invariant system, in the absence of any external periodic potential. Second, quasi-Bloch oscillations are a genuinely many-body phenomenon.

The exact conditions for the occurrence of the quasi-Bloch oscillations is a matter of a controversy [8–10,12–17]. However, it is undisputable that the many-body adiabaticity is a sufficient (although, in general, not necessary) condition for the quasi-Bloch oscillations [9]. This is the reason for our interest in conditions for adiabaticity in the one-dimensional impurity-fluid system.

Recently, we have applied a necessary adiabatic condition of Ref. [7] to the impurity-fluid system, with the result that in order to maintain adiabaticity the driving force \( F_N \) should vanish with the system size (with the density of the fluid being fixed) at least as fast as \( O(1/ \ln N) \) [18]. This result demonstrates that the adiabaticity does not survive in the thermodynamic limit, as is expected on general grounds for a gapless many-body system [19]. However, if the \( O(1/ \ln N) \) scaling were a true scaling of the maximal force tolerated by adiabaticity, it would be very possible to observe the adiabatic

1One could imagine that both sufficient and necessary conditions are available and provide bounds which are close to each other. In practice, however, such fortunate occurrences are rare, if not nonexistent, at least in the many-body context.
evolution in state-of-the-art cold-atom experimental settings with moderately large \( N \), e.g., with \( N \sim 100 \) like in the experiment of Ref. [11]. This observation motivated us to search for a necessary and sufficient adiabatic condition in the impurity-fluid system.

Here we report such a condition obtained in the framework of an effective theory describing an impurity slowly moving in a one-dimensional (1D) quantum fluid [20–22]. The condition has the form \( F_N < O(1/N) \), which is a dramatic quantitative difference from the logarithmic scaling obtained previously. This result implies that maintaining many-body adiabaticity in the impurity-fluid system is a very challenging experimental task.

This paper is organized as follows. After a general discussion of the notion of adiabaticity in Sec. II, we introduce the impurity-fluid system and its effective description in Sec. III. The diagonalization of the effective Hamiltonian at a given moment of time is reviewed in Sec. IV. The solution of the full dynamical problem is presented in Sec. V. In Sec. VI the results are presented, and their immediate experimental implications are discussed. In Sec. VII we summarize our results and make a couple of concluding remarks. Most of the technicalities are reserved for the appendixes.

II. ADIABATICITY: FIGURE OF MERIT

We start from introducing the notion of adiabaticity in quantitative terms. Consider a parameter-dependent Hamiltonian \( H_Q \), with \( Q \) being for a moment an abstract parameter. We introduce time dependence in this Hamiltonian by assuming that \( Q \) linearly varies with time \( t \),

\[ Q = F_N t. \]

At the moment \( F_N \) is treated merely as another abstract parameter quantifying the driving rate. The subscript \( N \) in \( F_N \) indicates that the driving rate may, in general, scale with the system size. For each \( Q \) one can define an instantaneous ground state \( \Phi_Q \), which is the lowest-eigenvalue solution to Schrödinger’s stationary equation,

\[ \hat{H}_Q \Phi_Q = E_Q \Phi_Q. \]  

Here \( E_Q \) is the instantaneous ground-state energy. We assume that the ground state is nondegenerate for any \( Q \).

The dynamics of the system is governed by the Schrödinger equation, which can be written in a rescaled form as

\[ (\hat{H}_Q - i F_N \partial_Q)\Psi_Q = 0. \]  

Here \( \Psi_Q \) is the state vector of the system, which depends on time through the time-dependent parameter \( Q \). Initially, at \( t = 0 \) (or, equivalently, at \( Q = 0 \)), the system is prepared in the instantaneous ground state:

\[ \Psi_0 = \Phi_0. \]  

The evolution is called adiabatic as long as the state of the system \( \Psi_Q \) stays close to the instantaneous ground state \( \Phi_Q \). To what extent adiabaticity is preserved during the evolution is quantified by the adiabatic fidelity \( \mathcal{F}_Q \), which is the probability that the evolved state coincides with the instantaneous ground state,

\[ \mathcal{F}_Q \equiv |\langle \Phi_Q | \Psi_Q \rangle|^2. \]

Perfect adiabaticity would imply \( \mathcal{F}_Q = 1 \). Calculating \( \mathcal{F}_Q \) for the driven impurity-fluid system is the main goal of our study.

Another useful quantity is the adiabatic mean free path \( Q_* \), which quantifies how far the system can travel in the parameter space for a given driving rate before the adiabaticity breaks down. We define this quantity as the smallest positive solution of the equation

\[ \ln \mathcal{F}_Q_* = -1. \]

III. IMPURITY-FLUID SYSTEM

A. Preliminary considerations

An object of our study is a one-dimensional many-body system, consisting of a quantum fluid and an impurity particle. A constant external force \( F_N \) is exerted upon the impurity. The fluid consists of \( N \) identical particles, either fermions or bosons. The particles of the fluid interact with the impurity and, in general, with each other.

As a preliminary step, we discuss how to describe a one-body problem of a noninteracting impurity particle pulled by an external force. This can be done conveniently by introducing a time-dependent Hamiltonian,

\[ \hat{H}_Q^{\text{imp}} = \left( \frac{\hat{P} + Q}{2m} \right)^2, \]

where \( m \) is the mass of the impurity, \( \hat{P} \equiv -i \partial / \partial X \) is the canonical momentum of the impurity, and \( X \) is the coordinate of the impurity. Periodic boundary conditions with some period \( L \) are implied. In this context \( Q = F_N t \) is the impulse of the force.

The interacting impurity-fluid system is described by the microscopic Hamiltonian

\[ \hat{H}_Q^{\text{micro}} = \hat{H}_Q^{\text{imp}} + \hat{H}_f + \hat{H}_{\text{if}}, \]

where \( \hat{H}_f \) is the Hamiltonian of the fluid which includes the kinetic term and the pairwise interactions between the particles of the fluid and \( \hat{H}_{\text{if}} \) is the impurity-fluid coupling. We do not specify microscopic Hamiltonians \( \hat{H}_f \) and \( \hat{H}_{\text{if}} \) explicitly since our analysis will be based on the effective low-energy model described in the next section.

A general feature of translation-invariant one-dimensional systems described by the Hamiltonian (7) is that eigenenergies as a function of \( Q \) is periodic in thermodynamic limit [23]. The period is determined by the number density of particles and, in our case, is given by \( 2k_F \) up to finite-size corrections, where \( k_F \equiv \pi N / L = \pi \rho \). The latter quantity sets the typical momentum scale of the problem. For the fluid consisting of noninteracting fermions, \( k_F \) coincides with the Fermi momentum. It should be emphasized, however, that in the general case we do not ascribe any “fermionic” meaning to \( k_F \). In particular, we consider bosonic and fermionic fluids on equal footing. Note that Fermi statistics plays no role in the above-mentioned periodicity of eigenenergies, which is present for bosons as well.
B. Effective Hamiltonian

Under fairly general conditions the low-lying excitations of a one-dimensional quantum fluid can be treated by means of an effective Luttinger liquid theory [24]. This theory can be extended to describe the low-energy sector of the one-dimensional impurity-fluid system (7) [20–22,25]. This extension is valid for a sufficiently small absolute value of the velocity of the impurity \(v_Q\) (below we will discuss this condition in more detail). The corresponding effective Hamiltonian reads [20–22,25]

\[
\hat{H} = v_Q \hat{P} + v_s \sum_q |q| \hat{a}^\dagger_q \hat{a}_q - \frac{1}{\sqrt{2\pi L}} \sum_q \sqrt{q} \delta^\dagger_Q \Delta v_Q \left( \hat{a}^\dagger_q e^{-iqX} + \hat{a}_q e^{iqX} \right).
\]

(8)

Here \(v_s\) is the sound velocity of the fluid, \(\hat{a}^\dagger_q, \hat{a}_q\) are creation and annihilation operators of bosonic excitations of the fluid carrying momentum \(q\).

\[
\Delta v_Q^0 \equiv v_s - v_Q \text{ sgn} q,
\]

(9)

and

\[
\delta^\dagger_Q = \begin{cases} 
\delta^\dagger, & q > 0, \\
0, & q = 0, \\
\delta, & q < 0,
\end{cases}
\]

(10)

is the scattering phase which is determined by the impurity-fluid interaction.

In Eq. (8) and throughout the paper the increment in the sums over \(q\) equals the momentum quantum \(\delta k = 2\pi/ L\). For definiteness, we employ the ultraviolet cutoff equal to \(k_F\), although the exact cutoff value does not enter the final results.

The total canonical momentum

\[
\hat{P}_{\text{tot}} \equiv \hat{P} + \sum_q q \hat{a}^\dagger_q \hat{a}_q
\]

(11)

commutes with the Hamiltonian \(H\) and therefore is conserved. Its eigenvalue is quantized in units of \(\delta k\). The canonical momentum should not be confused with the total kinetic momentum, \(\hat{P}_{\text{kin}} = \hat{P}_{\text{tot}} + Q\), which is not quantized and grows linearly with time due to the action of the external force.

In fact, the effective model (8) is well defined only in a subspace of the Hilbert space corresponding to some eigenvalue \(\Lambda_{\text{tot}}\) of the total canonical momentum. In this subspace the kinetic momentum also has a well-defined, although varying with time, eigenvalue \(P_{\text{kin}}\). Effective “constants” \(v_Q, \delta_Q^0\), and \(\delta_Q\) are, in fact, functions of the kinetic momentum, \(P_{\text{kin}} = P_{\text{tot}} + Q\) (and therefore the notations \(v_{\text{kin}}, \delta_{\text{kin}}^0, \delta_{\text{kin}}\) would be more consistent). We, however, choose to fix \(P_{\text{tot}}\) and refrain from referring to it explicitly throughout the paper, except in the present section and Appendix A. The only importance of the precise value of \(P_{\text{tot}}\) is to fix \(v_Q, \delta_Q^0, \delta_Q\) at \(Q = 0\).

The range of validity of the effective Hamiltonian (8) is a somewhat subtle issue. This Hamiltonian is designed to describe a low-energy sector of the Hilbert space, i.e., an energy shell of a width \(\Delta E_{\text{kin}}\) above the ground state. The subscript in \(\Delta E_{\text{kin}}\) indicates that the range of validity of the effective model varies with \(P_{\text{kin}}\). In particular, this is the case for the integrable model solved by McGuire [28], which is discussed in Appendix E. In other cases, however, the effective model (8) breaks down in a finite portion of the Brillouin zone. In particular, this happens for a sufficiently light impurity weakly interacting with a one-dimensional fluid [12,14]. In general, one expects that \(\Delta E_{\text{kin}}\) is nonzero as long as the impurity moves with the velocity below the generalized critical velocity \(v_s \leq v_c\), which ensures the absence of the Cherenkov-like radiation [14]. The latter critical velocity is typically on the order of \(v_s\). To summarize, to be on the safe side, one can assume

\[
v_Q \ll v_s,
\]

(12)

although the actual range of the validity of the effective Hamiltonian (8) can be much wider.

IV. INSTANTANEOUS GROUND STATE

The first ingredient required for calculating the adiabatic fidelity \(F_Q\) is the instantaneous ground state \(\Phi_Q\) of the Hamiltonian (8). The latter can be diagonalized exactly [20–22]. We describe and discuss the diagonalization procedure and identification of the ground state in Appendix A. Here we give the final result, which reads

\[
e^{\hat{W}_{\text{Q}}} \hat{H}_{\text{Q}} e^{-\hat{W}_{\text{Q}}} = \hat{H}_{\text{Q}}^d + C,
\]

(13)

where

\[
\hat{H}_{\text{Q}} = v_Q \hat{P} + v_s \sum_q |q| \hat{a}^\dagger_q \hat{a}_q
\]

(14)

and \(C\) is a constant which is omitted in what follows.4 The anti-Hermitian operator \(W\) reads

\[
\hat{W}_{\text{Q}} = \sum_q (\alpha_q^\dagger \hat{a}^\dagger_q e^{-iqX} - \alpha_q \hat{a}_q e^{iqX}),
\]

(15)

where

\[
\alpha_q^\dagger = -\frac{\delta_q^0}{\sqrt{2\pi L q}}
\]

(16)

4In fact, a \(Q\)-dependent \(c\) number responsible for reproducing the correct ground-state energy is already omitted in the definition (8) of \(H\). Predicting the ground-state energy is beyond the scope of the effective low-energy model.
and the overbar in $\bar{\alpha}_q$ and elsewhere refers to complex conjugation.

The ground state of $\hat{H}_Q$ for a fixed total canonical momentum $P_{\text{tot}}$ reads

$$\Phi_Q = e^{-\hat{\mathcal{H}}_Q} |\text{vac}, \, P_{\text{tot}}\rangle,$$  \hspace{1cm} (17)

where $|\text{vac}, \, P_{\text{tot}}\rangle \equiv |\text{vac}\rangle \otimes |P_{\text{tot}}\rangle$ is a product state, with $|\text{vac}\rangle$ being a Fock vacuum with respect to bosonic operators $\hat{a}_q$ and $|P_{\text{tot}}\rangle$ being the state of impurity with the momentum $P_{\text{tot}}$.

V. DYNAMICS

A. Dynamical diagonalization

The second ingredient for calculating $F_Q$ is the dynamical state vector $\Psi_Q$ evolving according to the Schrödinger equation (2). Remarkably, the dynamics described by this equation is integrable, in the sense that the operator $(\hat{H}_Q - i F \partial_Q)$ can be diagonalized by a unitary transformation analogous to the transformation (13):

$$e^{\hat{\mathcal{H}}_Q} (\hat{H}_Q - i F \partial_Q) e^{-\hat{\mathcal{H}}_Q} = \hat{H}_Q^d - i F \partial_Q + C'.$$  \hspace{1cm} (18)

Here $C'$ is a $c$ number which will be omitted in what follows,

$$\hat{Y}_Q = \sum_q \left( \beta_q^d \hat{a}_q e^{-i q X} - \beta_{q}^d \overline{\hat{a}}_q e^{i q X} \right),$$  \hspace{1cm} (19)

and the coefficients $\beta^d_q$ satisfy the differential equation

$$i F \partial_q \beta^d_q - |q| \Delta v^d_q \beta^d_q - \frac{\delta^d_q}{\sqrt{2 \pi L}} \overline{\sqrt{|q|}} \Delta v^d_q = 0.$$  \hspace{1cm} (20)

The key insight behind the dynamical diagonalization is that $[Y_Q, \partial_Q Y_Q]$ is a $c$ number, and therefore

$$-i F e^{\hat{\mathcal{H}}_Q} (\partial_Q e^{-\hat{\mathcal{H}}_Q}) = i F \left( \partial_Q Y_Q + \frac{1}{2} [Y_Q, \partial_Q Y_Q] \right)$$  \hspace{1cm} (21)

is linear in boson operators and has the same structure as $Y_Q$ and the interaction term of the Hamiltonian (8).

In order to satisfy the initial condition of the Schrödinger equation (3) we supplement the differential equation (20) with the initial condition

$$\beta^d_0 = \alpha^d_0.$$  \hspace{1cm} (22)

Equation (20) can be solved in quadratures, with the result

$$\beta^d_q = \alpha^d_q - \int_0^Q \partial_Q \alpha^d_q \exp \left( -i \frac{|q|}{F_N} \int_0^Q d\tilde{Q} \Delta v^d_q \right) dQ'.$$  \hspace{1cm} (23)

Equation (18) entails that the dynamical state of the system evolves according to

$$\Psi_Q = e^{-\hat{\mathcal{H}}_Q} |\text{vac}, \, P_{\text{tot}}\rangle,$$  \hspace{1cm} (24)

where an irrelevant $c$-number phase factor is omitted.

B. Dynamics of adiabatic fidelity

With $\Phi_Q$ and $\Psi_Q$ in hand, we are prepared to proceed to calculating the adiabatic fidelity. Substituting Eqs. (17) and (24) into the definition (4) and applying the Baker-Campbell-Hausdorff formula, we obtain

$$F_Q = \exp \left( - \sum_q |\beta^d_q - \alpha^d_q|^2 \right).$$  \hspace{1cm} (25)

VI. RESULTS

A. Adiabatic fidelity and adiabatic mean free path

In principle, Eq. (25) along with Eqs. (16) and (23) allows one to calculate $F_Q$ for any values of parameters within the validity range of the model (8). However, a further asymptotic analysis is required to reveal the scaling properties of $F(Q)$. Here we present the main results of such an analysis, the details can be found in Appendix B.

The asymptotic behavior of the adiabatic fidelity in the limit of large system size depends on how the force $F_N$ scales with $N$. Before turning to a general case, we consider an important special case of the force independent of the system size. In this case we obtain

$$\ln F_Q = - \xi^2 \left( \frac{Q}{k_F} \right)^2 \ln N,$$  \hspace{1cm} (26)

with

$$\xi^2 = k_F^2 \left[ \left( \frac{\partial \delta Q}{\partial Q} \right)^2 + \left( \frac{\partial \delta Q}{\partial Q} \right)^2 \right].$$  \hspace{1cm} (27)

This is the leading term of the double asymptotic expansion of $\ln F_Q$ in the limit of

$$N \gg 1, \quad N/L \text{ fixed,}$$  \hspace{1cm} (28)

$$Q \ll k_F,$$  \hspace{1cm} (29)

where the limit (28) is taken first.

Quite remarkably, the force does not enter Eq. (26). In fact, the adiabatic fidelity follows the orthogonality catastrophe overlap, $F_Q \sim C_Q \equiv |\langle \Phi_Q | \Phi_Q \rangle|^2$, in agreement with the scenario described in Ref. [7]. One can obtain Eq. (26) without solving the Schrödinger equation (2) by calculating the orthogonality overlap $|\langle \Phi_Q | \Phi_Q \rangle|^2$ and applying the general result of Ref. [7], as detailed in Appendix C. However, this method is applicable in a superficially narrow range of $Q$ which shrinks when $N$ grows, $Q = o(1)$. Solving the dynamical problem shows that the validity range of the approximation $F_Q \sim C_Q$ appears to be larger than that guaranteed by the rigorous result of Ref. [7]; analogous conclusions have been made on the basis of explicit solutions of other models [7].

Now we turn to a more general case when the force $F_N$ can vanish with the system size but not faster than $1/N$. In this case one needs to consider separately two ranges of $Q$:

$$\ln F_Q = - \xi^2 \left( \frac{Q}{k_F} \right)^2 \ln N;$$  \hspace{1cm} (29)

$$Q \ll F_N,$$  \hspace{1cm} (30)

$$Q \ll k_F,$$  \hspace{1cm} (31)

where $\xi$ is given by Eq. (27).

One can see from the first line of Eq. (30) that for small momenta the result coincides with those of Eq. (26) and, in

$$\ln F_Q = - \xi^2 \left( \frac{Q}{k_F} \right)^2 \ln N;$$  \hspace{1cm} (29)

$$Q \ll F_N,$$  \hspace{1cm} (30)

$$Q \ll k_F,$$  \hspace{1cm} (31)

where $\xi$ is given by Eq. (27).

One can see from the first line of Eq. (30) that for small momenta the result coincides with those of Eq. (26) and, in
fact, again can be obtained by the method of Ref. [7] without considering dynamics (see Appendix C). Solving the time-dependent Schrödinger equation explicitly is mandatory for obtaining the expression for larger momenta [the second line of Eq. (30)]. Observe that the latter expression manifestly depends on the force.

One can easily find the adiabatic mean free path \( Q_\ast \) from Eq. (30):

\[
\frac{Q_\ast}{k_F} = \frac{1}{\xi} \left\{ (\ln N)^{-\frac{1}{2}}, \quad \frac{F_N}{k_F v_F} \gg \frac{1}{\sqrt{\ln N}}, \quad \frac{1}{\sqrt{\ln N}} \ll \frac{F_N}{k_F v_F} \ll \frac{1}{\ln N}. \right. 
\]

(31)

Again, solving the time-dependent Schrödinger equation is essential for obtaining the second line of this expression, while the first line is obtained in Ref. [18] without considering dynamics.

Equation (31) implies that

\[
F_N = O(1/N)
\]

(32)

is a necessary asymptotic condition for the many-body adiabaticity. While the scaling \( F_N = O(1/N) \) is beyond the range of validity of Eq. (31), the second line of this equation indicates that the condition (32) can also be sufficient for adiabaticity. This is indeed the case, as proved in Appendix D. Thus, we establish Eq. (32) as a necessary and sufficient condition for adiabaticity in the driven impurity-fluid system in the large-system-size limit.

At this point it is worth commenting on the results of Ref. [18], where we have studied the adiabaticity breakdown in the integrable impurity-fluid system using the method of Ref. [7]. Integrability allowed us to calculate \( C_Q \) explicitly and establish the necessary adiabatic condition \( F_N \leq O(1/\ln N) \). Obviously, the result (32) of the present paper is much stronger. However, the analysis of a microscopic integrable model retains its value since it underpins the effective model (8) and illustrates how the phenomenological parameters of the effective model are related to microscopic parameters. We present this analysis in Appendix E. A particularly interesting conclusion from this analysis is that the results of the calculation of \( C_Q \) within the integrable model and within the effective model (8) coincide for all \( Q \), not just within the conservative validity range (12). This indicates that the actual validity range of the effective treatment can be larger.

B. Experimental implication

Damped oscillations of an impurity particle driven through a 1D quantum fluid were observed in a recent experiment [11] where a quantum fluid consisted of \( N \approx 60 \) cesium atoms and the force was equal to 1/3 of the gravitational force. In this experiment the adiabaticity faded away on a timescale smaller than one period of the oscillations. This was clear from the direct numerical simulation of the experimental setting, which showed the increase of the energy of the system above the ground-state energy already during the first half period of oscillations [11]. We confirm this conclusion by calculating the adiabatic mean free path according to Eq. (31), which appears to be around \( k_F \) throughout the whole range of experimental conditions.

One may wonder whether it is possible to maintain the many-body adiabaticity for at least a few cycles of oscillations by, e.g., applying a smaller force. Our analysis indicates that this can be challenging. The reason is that in practice the spatial amplitude of oscillations is limited by the size of the quasi-1D optical cigar-shaped trap. It is easy to see that this amplitude is inversely proportional to the force. This constrains the force to satisfy \( F_N \geq k_F^2/(2\pi m_s N) \), where \( m_s \) is the effective mass of the impurity in the fluid. This can be only marginally consistent with the adiabatic condition (32). Therefore maintaining adiabaticity for several periods of the oscillations would require an extremely careful choice of the driving force: not too high to sustain adiabaticity but not too low to keep the spatial amplitude of the oscillations within the trap size.

VII. SUMMARY AND CONCLUDING REMARKS

To summarize, we have analyzed the dynamics of the many-body adiabatic fidelity \( F_Q \) in a one-dimensional impurity-fluid system where a force applied to the impurity pulls the latter through the fluid. We have employed an effective low-energy theory which enabled us to find explicit expressions for \( F_Q \) and the adiabatic mean free path \( Q_\ast \) in terms of the size of the system and the effective parameters of the fluid and the impurity-fluid coupling. Our results imply that the state vector of the impurity-fluid system completely departs from the instantaneous ground state already for acquired momenta which are vanishingly small in thermodynamic limit, unless the force scales down with the system size as \( 1/N \) or faster. This dramatically improves the necessary adiabatic condition \( F_N \leq O(1/\ln N) \) obtained previously [18].

It is remarkable that quantum adiabaticity breaks down already at small acquired momenta. When the acquired momentum reaches the vicinity of \( \pi \rho \), our quantitative results may be inapplicable since the vicinity of \( Q = \pi \rho \) can be beyond the range of validity of the effective model employed. However, the adiabaticity is gone anyway at this point [unless \( F_N \leq O(1/N) \)]. It is worth noting that crossing the \( Q = \pi \rho \) point is potentially the most dangerous part of the oscillation cycle with respect to preserving adiabaticity [15,16]. This means that the adiabatic condition can become only more stringent beyond the range of validity of our approach.

It should be emphasized that while the many-body adiabaticity would be sufficient to ensure quasi-Bloch oscillations in an arbitrary one-dimensional impurity-fluid system, adiabaticity breakdown is not necessarily fatal for oscillations. This has been confirmed in the experiment of Ref. [11]. On the theoretical side, there is a consensus that quasi-Bloch oscillation of a sufficiently heavy impurity can occur in the thermodynamic limit for a force which is finite in the thermodynamic limit but sufficiently small compared to other intensive (i.e., independent of the system size) quantities of the system [8–10,12,14]. The latter condition defines a notion of a thermodynamic adiabaticity, as contrasted by the genuine quantum many-body adiabaticity studied in the present paper.5

5An instructive demonstration of the subtlety of the interplay between the two concepts was recently presented in Ref. [27], where
It is a matter of debate whether thermodynamic adiabaticity is sufficient for quasi-Bloch oscillations of light impurity [8–10,12–17]. The present paper does not contribute to this debate.

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APPENDIX A: DIAGONALIZATION FOR A FIXED Q

Here we present the details of diagonalization of the Hamiltonian $H(Q)$ at a given $Q$.

1. Unitary transformation

Unitary transformation generated by $W$ defined by Eq. (15) with arbitrary coefficients $\alpha_q^0$ yields

$$e^{\hat{W}_0} \hat{\alpha}_q e^{-\hat{W}_0} = \hat{\alpha}_q - \alpha_q^0 e^{-i\pi X}$$

(A1)

and

$$e^{\hat{W}_0} \hat{P} e^{-\hat{W}_0} = \hat{P} + \sum_q q (\alpha_q^0 \hat{\alpha}_q e^{-i\pi X} + \text{H.c.}) - \Delta P,$$

(A2)

where the constant $\Delta P$ reads

$$\Delta P = \sum_q q |\alpha_q^0|^2.$$  

(A3)

As a consequence,

$$e^{\hat{W}_0} \hat{P}_{\text{tot}} e^{-\hat{W}_0} = \hat{P}_{\text{tot}}$$

(A4)

and

$$e^{\hat{W}_0} \hat{H} e^{-\hat{W}_0} = \hat{H}_d + C - \sum_q q |\alpha_q^0| \left(\alpha_q \Delta v^q + \frac{\delta_q^q}{\sqrt{2\pi L|q|}} \Delta v_q^q \right) \times \hat{\alpha}_q e^{-i\pi X} + \text{H.c.},$$

(A5)

with $\hat{H}_d$ given by Eq. (14) and the constant $C$ given by

$$C = v_s \sum_q q |\alpha_q^0|^2$$

$$+ \frac{1}{\sqrt{2\pi L}} \sum_q \sqrt{|q|} \delta_q^q \Delta v^q \left(\alpha_q^0 + \alpha_q^0\right).$$  

(A6)

The last term in Eq. (A5) vanishes when one chooses coefficients $\alpha_q^0$ according to (16). In this case

$$\Delta P = \frac{k_F}{2\pi} \left[\delta_q^+ - \delta_q^-\right]^2.$$  

(A7)

an impurity-fluid system with a time-dependent coupling constant (but in the absence of a force) was considered. It appeared that the outcomes of the thermodynamically adiabatic and genuinely quantum adiabatic evolutions could be identical or dramatically different in the same system, depending on the choice of the initial state.

2. Ground state

Eigenstates of the diagonalized Hamiltonian (14) do not depend on $Q$ and read

$$|\{n_q\},K\rangle \equiv |\{n_q\}\rangle \otimes |K\rangle,$$

(A8)

where $|\{n_q\}\rangle$ is an eigenstate of the oscillator part of (14) with $n_q$ bosons for each $q$, while $|K\rangle$ is the state of the impurity with momentum $K$. The eigenvalues of the Hamiltonian and the total momentum read, respectively,

$$E_{[n_q],K} = v_Q K + v_s \sum_q q |n_q + C,$$

(A9)

$$P_{\text{tot}[n_q],K} = K + \sum_q q n_q.$$  

(A10)

We wish to find the minimal eigenenergy $E_{[n_q],K}$ for a given total momentum $P_{\text{tot}[n_q],K}$. To this end we introduce

$$P_{\pm} \equiv \sum_{\Delta q > 0} q |n_q$$

(A11)

and rewrite Eqs. (A9) and (A10) as, respectively,

$$E_{[n_q],P_{\text{tot}}} = v_Q K + v_s (P_+ + P_-) + C,$$

(A12)

$$P_{\text{tot}[n_q],K} = K + P_+ - P_-.$$  

(A13)

This leads to

$$E_{[n_q],K} = v P_{\text{tot}} + C + (v_s - v)P_+ + (v_s + v)P_-.$$  

(A14)

Since the last two terms in this equation are nonnegative while $C$ is the same for all $|n_q\rangle$ and $K$, the right-hand side (rhs) of Eq. (A14) has a minimum at $P_+ = P_-$. Hence the ground state of $\hat{H}_d$ for a given value of the total momentum reads $|\text{vac}, P_{\text{tot}}\rangle$, where $|\text{vac}\rangle$ is $|\{n_q\}\rangle$ with all $n_q = 0$, and the ground state $\Phi_Q$ of $\hat{H}_Q$ is given by Eq. (17).

We note that

$$\langle \Phi_Q | \hat{P} | \Phi_Q \rangle = P_{\text{tot}} - \Delta P.$$  

(A15)

APPENDIX B: ASYMPTOTIC ANALYSIS

From Eqs. (25) and (23) one gets

$$\ln \mathcal{F}(Q) = - \int_0^Q \int_0^Q dQ^d Q^d \frac{\partial \phi^d}{\partial Q^d}$$

$$\times \frac{\partial Q^d}{\partial \phi^d} \Sigma_+(Q'^d - Q^d)$$

$$- \{\delta^d_+ \rightarrow \delta_-, \Sigma_+ \rightarrow \Sigma_-\},$$  

(B1)

with

$$\Sigma_{\pm} \equiv \sum_{n=1}^{M-1} \frac{1}{n} e^{-\delta n \phi_{\pm}} = - \frac{1}{2} \ln(2 - 2 \cos \phi_{\pm}) + Ci(M \phi_{\pm})$$

(B2)

and

$$\phi_{\pm} \equiv \frac{\delta k}{F_N} \int_{Q^d} d\tilde{Q} (v_+ \mp v_Q).$$  

(B3)
Here Ci stands for the cosine integral function, and we use an arbitrary momentum cutoff
\[ u \equiv (M - 1)\delta k \sim k_F. \tag{B4} \]
If one takes \( u = k_F \), as we do in the rest of the paper, then \( M = (N + 1)/2 \). Note that
\[ \phi_\pm \ll 1 \tag{B5} \]
since we assume that \( N F_N \gg 1 \).

We analyze the asymptotic behavior of the rhs of Eq. (B2) in two opposite limits. The first one is
\[ M\phi_\pm \ll 1, \tag{B6} \]
which is equivalent to \( Q \ll F_N/(v_\kappa k_F) \). In this case one gets
\[ \Sigma_\pm = \ln M + \gamma_E + O((\phi_\pm M)^2) + O\left( \frac{1}{N} \right) \]
\[ = \ln N + \gamma_E + \ln \frac{M}{N} + O\left( \left( \frac{Qv_\kappa k_F}{F_N} \right)^2 \right). \tag{B7} \]
Observe that only the leading term is cutoff independent. The second limit is
\[ M\phi_\pm \gg 1, \tag{B8} \]
which is equivalent to \( Q \gg F_N/(v_\kappa k_F) \). In this case the cutoff-dependent part of the rhs of Eq. (B2) is vanishingly small,
\[ \text{Ci}(M\phi_\pm) = O\left( \frac{1}{M\phi_\pm} \right). \tag{B9} \]
We find it convenient to further expand the cutoff-independent part of Eq. (B2) in \( \phi_\pm \) to obtain
\[ \Sigma_\pm = -\ln \phi_\pm + O(\phi_\pm^2) + O\left( \frac{1}{N\phi_\pm} \right). \tag{B10} \]

Now we can substitute the asymptotic expansions (B7) and (B10) into Eq. (B1). Since the adiabaticity breaks down already for \( Q/k_F = o(1) \), it is reasonable to further expand all functions in the integrand in Eq. (B1) over \( Q \) and \( Q' \). This finally leads to Eq. (30).

APPENDIX C: ORTHOGONALITY CATASTROPHE AND ADIABATICITY

1. Orthogonality catastrophe

From Eqs. (17) and (15) one can calculate the orthogonality overlap
\[ C_Q \equiv |\langle \Phi_0 | \Phi_Q \rangle|^2 = \exp \left( -\sum_q |\alpha_q^+ - \alpha_0^q|^2 \right). \tag{C1} \]
This is done with the help of the relation
\[ e^{(a - a')\hat{a}^\dagger} e^{-(a - a')\hat{a}} = e^{(a - a')\hat{a}^\dagger} e^{-(a - a')/2 - i \Im a\hat{a}}, \tag{C2} \]
which is valid for arbitrary \( a \) and \( a' \).

We calculate \( \ln C_Q \) explicitly by substituting the expression (16) for \( \alpha_0^q \) in Eq. (C1), expanding the latter in \( Q \) and performing the sum over \( q \). In this way we obtain in the leading order the rhs of Eq. (26).

2. Relation between the adiabaticity and the orthogonality catastrophe

We wish to establish that \( C_Q \simeq F_Q \) in the limit of large \( N \).

The rigorous form of this relation was proven in Ref. [7]. In the considered case it reads
\[ |C_Q - F_Q| \leq \frac{1}{F_N} \int_0^Q dQ' \sqrt{\langle \Psi_0 | \hat{H}_Q^2 | \Psi_0 \rangle - \langle \Psi_0 | \hat{H}_Q | \Psi_0 \rangle^2}. \tag{C3} \]
The integral in this equation is cutoff dependent but does not diverge with \( N \). Therefore the rhs of the inequality can be made vanishingly small, i.e., \( o(1) \) in the limit \( N \to \infty \), whenever \( Q = o(F_N) \). In this way we reproduce Eq. (26) and the first line of Eq. (30) without solving the Schrödinger equation, with the use of a shortcut introduced in Ref. [7]. The price is a superficially reduced range of validity of the results, \( Q = o(F_N) \) instead of \( Q = O(F_N) \) obtained by solving the dynamical problem.

APPENDIX D: PROOF THAT \( F_N = O(N) \) IS SUFFICIENT FOR ADIABATICITY

For the sake of such a proof we take the integral in Eq. (23) by parts. Importantly, this produces the factor \( F_N/q \), which eventually does the job. In this way for \( q > 0 \) one estimates
\[ |\beta_Q^q - \alpha_Q^q|^2 \leq \frac{F_N^2}{2\pi q^4 L} \left( A_Q^\pm \right)^2, \tag{D1} \]
where
\[ \frac{A_Q^\pm}{k_F v_\kappa} \equiv \left| \frac{\partial Q^\pm}{\partial Q} \right| \left| \frac{\partial Q^\pm}{\partial Q} \right|_{Q=0} + \int_0^Q dQ' \partial Q^\pm \left( \delta^\pm_{Q} \right) \left( \frac{dQ}{v_\kappa} \right) \tag{D2} \]
does not depend on \( q \) and is finite in the thermodynamic limit. Here the sign in \( A_Q^\pm \) and \( \delta^\pm_{Q} \) is the same as the sign of \( q \).

The sum over positive \( q \) in Eq. (25) is bounded from above according to
\[ \sum_{q>0} |\beta_Q^q - \alpha_Q^q|^2 \leq \xi(3) \left( \frac{A_Q^+ F_N N}{4\pi k_F^2 v_\kappa} \right)^2, \tag{D3} \]
and the sum over negative \( q \) is constrained analogously. Here \( \xi(3) \) is the Riemann zeta function. This proves that whenever
\[ F_N \leq \frac{\epsilon}{N} \frac{4\pi k_F^2 v_\kappa}{\sqrt{\xi(3) (A_Q^+)^2 + (A_Q^-)^2}}, \tag{D4} \]
the adiabatic fidelity is bounded from below,
\[ F_Q \geq e^{-\epsilon^2}. \tag{D5} \]

APPENDIX E: CONSISTENCY CHECK FOR THE INTEGRABLE MODEL

Here we consider a microscopic integrable impurity-fluid model. The model consists of \( N \) fermions and a single impurity particle with a mass \( m \) equal to the mass of a fermion. Fermions
do not interact with each other but couple to the impurity via the repulsive contact potential. The Hamiltonian reads
\[ \hat{H}_Q = \frac{-i}{\hbar} \nabla + Q \nabla \nabla + \sum_{j=1}^{N} \frac{1}{2m} \hat{a}_{j}^{\dagger} + g \sum_{j=1}^{N} \delta(x_j - x) \],
(E1)
where \( Q = F t \) is the impulse of the force, \( X \) and \( x_j \) are the coordinates of the impurity and the \( j \)-th fermion, respectively, and \( g > 0 \) is the impurity-fluid coupling.

For a fixed \( Q \) the model (E1) is integrable, as shown by McGuire [28]. In fact, this model is one of the simplest models solvable via the Bethe ansatz: Its eigenfunction can be expressed through (8) and better understand its validity range. In what follows the technique of Ref. [7] and relate the adiabaticity breakdown to the orthogonality catastrophe in Ref. [18]. Here we focus on the relation between the microscopic model (E1) and the adiabatic breakdown to the orthogonality catastrophe in Ref. [18]. Here we focus on the relation between the microscopic model (E1) and the adiabatic breakdown to the orthogonality catastrophe in Ref. [18]. Here we focus on the relation between the microscopic model (E1) and the adiabatic breakdown to the orthogonality catastrophe in Ref. [18]. Here we focus on the relation between the microscopic model (E1) and the adiabatic breakdown to the orthogonality catastrophe in Ref. [18]. Here we focus on the relation between the microscopic model (E1) and the adiabatic breakdown to the orthogonality catastrophe in Ref. [18]. Here we focus on the relation between the microscopic model (E1) and the adiabatic breakdown to the orthogonality catastrophe in Ref. [18]. Here we focus on the relation between the microscopic model (E1) and the adiabatic breakdown to the orthogonality catastrophe in Ref. [18].

First, we would like to relate the effective scattering phases in Eq. (8) to the microscopic scattering phases of the Bethe ansatz [27],
\[ \delta_{BA}^\pm = \frac{\pi}{2} - \arctan \left( \frac{\Lambda}{2 v_s} \right), \]
(E2)
where \( v_s \) is the Fermi velocity in the present context and the parameter \( \Lambda \) can be found from the equation [27]
\[ \frac{Q}{k_f} = \frac{g}{2 \pi v_s} \left[ \left( \Lambda + \frac{2 v_s}{g} \right) \arctan \left( \Lambda + \frac{2 v_s}{g} \right) - \left( \Lambda - \frac{2 v_s}{g} \right) \arctan \left( \Lambda - \frac{2 v_s}{g} \right) + \frac{1}{2} \ln \left( 1 + (2 v_s/g - \Lambda)^2 \right) \right], \]
(E3)
To do this we consider the overlap
\[ \tilde{C}_Q = |\langle \Phi_Q | \Phi_Q \rangle|^2 \]
(E4)
between the ground state \( \Phi_Q \) and the noninteracting ground state \( \Phi_0 \) of the impurity-fluid system. \( \tilde{C}_Q \) can be calculated in both the microscopic model (E1) [31] and the effective model (8), with the leading order results
\[ \ln \tilde{C}_Q = - \left[ \frac{\delta_{BA}^+}{\pi} + \frac{\delta_{BA}^-}{\pi} \right]^2 \ln N \]
(E5)
and
\[ \ln \tilde{C}_Q = - \left[ \left( \frac{\delta_{Q}^+}{2 \pi} \right)^2 + \left( \frac{\delta_{Q}^-}{2 \pi} \right)^2 \right] \ln N, \]
(E6)
respectively. These two equations are compatible when
\[ \delta_{Q}^\pm = 2 \delta_{BA}^\pm. \]
(E7)
Now we turn to the orthogonality overlap of interest, \( C_Q \). We have calculated it within the McGuire model in a manner similar to that in Ref. [31], with the result
\[ \ln C_Q = - \left[ \left( \frac{\delta_{BA}^+ - \delta_{BA}^-}{\pi} \right)^2 + \left( \frac{\delta_{BA}^+ - \delta_{BA}^-}{\pi} \right)^2 \right] \times \left[ \ln N + O(1) \right], \]
(E8)
where \( O(1) \) refers to the limit of \( N \to \infty \).

After accounting for Eq. (E7) and expanding in small \( Q \) this result agrees with Eq. (26). It should be emphasized that this agreement is not limited to small \( Q \) but takes place in the whole Brillouin zone, \( -k_F < Q < k_F \). This indicates that the actual range of validity of the effective model can span well beyond the conservative condition (12). This agreement also suggests that the boson operators \( a_q, a_q^\dagger \) in fact do not depend on \( Q \) (see footnote 3), at least in the impurity-fluid system described by the microscopic Hamiltonian (E1).

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