Article

Nonadiabatic Energy Fluctuations of Scale-Invariant Quantum Systems in a Time-Dependent Trap

Mathieu Beau and Adolfo del Campo

1 Department of Physics, University of Massachusetts, Boston, MA 02125, USA; mathieu.beau.89@gmail.com
2 Donostia International Physics Center, E-20018 San Sebastián, Spain
3 IKERBASQUE, Basque Foundation for Science, E-48013 Bilbao, Spain
4 Theory Division, Los Alamos National Laboratory, MS-B213, Los Alamos, NM 87545, USA
* Correspondence: adolfo.delcampo@dipc.org

Received: 14 April 2020; Accepted: 29 April 2020; Published: 30 April 2020

Abstract: We consider the nonadiabatic energy fluctuations of a many-body system in a time-dependent harmonic trap. In the presence of scale-invariance, the dynamics becomes self-similar and the nondiabatic energy fluctuations can be found in terms of the initial expectation values of the second moments of the Hamiltonian, square position, and squeezing operators. Nonadiabatic features are expressed in terms of the scaling factor governing the size of the atomic cloud, which can be extracted from time-of-flight images. We apply this exact relation to a number of examples: the single-particle harmonic oscillator, the one-dimensional Calogero-Sutherland model, describing bosons with inverse-square interactions that includes the non-interacting Bose gas and the Tonks-Girardeau gas as limiting cases, and the unitary Fermi gas. We illustrate these results for various expansion protocols involving sudden quenches of the trap frequency, linear ramps and shortcuts to adiabaticity. Our results pave the way to the experimental study of nonadiabatic energy fluctuations in driven quantum fluids.

Keywords: quantum fluids; trapped systems; scale invariance; quantum dynamics; energy fluctuations; Calogero-Sutherland model; unitary Fermi gas; sudden quenches; shortcuts to adiabaticity

1. Introduction

Non-equilibrium quantum phenomena in many-body systems are often hard to describe by numerical methods due to the growth of entanglement with the time of evolution. In some cases, exact solutions can be found by analytical methods. This is often the case when the dynamics is characterized by the presence of dynamical invariants, also known as invariants of motion. Such conserved quantities can give rise to symmetries manifested in the dynamics. A prominent example is the presence of scale-invariance in the dynamics of a broad family of quantum fluids in time-dependent traps, which is exploited with routine in ultracold atom laboratories [1,2]. In this context, time-of-flight imaging techniques rely on the scale-invariant dynamics of a suddenly released atomic cloud, that expands freely [3–5]. Scale invariance is then manifested in local correlation functions, such as the density profile or the density-density correlation function, that become self-similar. One can then relate the correlations measured at a given time after the expansion with those of the initial trapped system.

Scale invariance is not restricted to free dynamics and it is also common in the presence of interactions. Indeed, in harmonically trapped systems it is always realized when the interparticle potential has the same scaling dimension as the kinetic energy. As an example, Bose-Einstein condensates with s-wave contact interactions exhibit it exactly in two-spatial dimensions as a result of the Pitaevskii-Rosch symmetry [6]. Similarly, it is present in the one-dimensional (rational) Calogero-Sutherland gas
describing bosons with inverse-square interactions [7]. Scale invariance is also manifested in systems with divergent interactions. In one-dimensional harmonically-confined systems it is respected exactly in the presence of hard-core interactions. It thus rules the dynamics of spin-polarized fermions (free fermions, where the Pauli exclusion principle gives rise to an effective hard-core interaction) as well as bosons in the Tonks-Girardeau regime [8–10]. In three spatial dimensions it governs the dynamics of a unitary Fermi gas [11,12]. Finally, scale-invariance can also be invoked effectively in some regimes. Ultracold bosons well-described by mean-field theory exhibit this symmetry in arbitrary spatial dimensions in the Thomas-Fermi regime, where the contribution of the kinetic energy can be neglected [3–5].

Beyond the use of time-of-flight imaging techniques, scale invariance has applications in a wide variety of contexts. It allows for the engineering of fast control protocols known as shortcuts to adiabaticity [13–17], as demonstrated in the laboratory [18–23]. It makes possible an exact treatment of the finite-time thermodynamics [24] and its optimization [25–28], as well as the suppression of quantum friction in quantum machines (like engines and refrigerators) that use a scale-invariant quantum fluid as a working substance [28–30]. Exploiting scale invariance, frictionless expansion and compression strokes have been implemented in a unitary Fermi gas [31]. In addition, the study of scale invariance in the presence of perturbations can be used as a probe for quantum anomalies [32], the finite amplitude of the interaction strength [33], and the breakdown of effective descriptions based on dimensional reduction [34]. In foundations of physics, it can shed light on regimes that would be challenging to explore experimentally or numerically. For instance, it can be used to describe exactly the quantum decay of quantum states at arbitrarily long times [35].

While local correlations functions are simply related at different times via scale invariance, other observables exhibit a more intricate dependence. A prominent one is the mean energy of the system that has been studied both in single-particle [36] and many-particle [24,28] quantum systems and that is crucial for applications in quantum thermodynamics.

In this work, we focus on the characterization of energy fluctuations in many-particle quantum systems exhibiting scale invariance. Energy fluctuations are known to govern characteristic time scales in the dynamics in view of the Mandelstam-Tamm time-energy uncertainty relation [37]. More generally, they play a crucial role in the formulation of speed limits, that establish the minimum time for a state to evolve into a distinct state, in both classical [38,39] and quantum [40] systems. Energy fluctuations can also be used to quantify the cost of thermodynamic processes [41,42], including shortcuts to adiabaticity, and their minimization paves the way to the finding of optimal protocols, e.g., the solution of the quantum brachistochrone problem [43,44].

2. Exact Many-Body Dynamics under Scale Invariance

The description of the exact many-body dynamics under scale invariance can be simplified using a sequence of symmetry transformations. To this end, we follow closely Lohe [36] and in particular the formulation for many-body systems in [24,28]. Consider a N-particle quantum system in D spatial dimensions and let us denote by \( \vec{r}_i \) and \( \vec{p}_i \) the position and momentum of the i-th particle, respectively. We consider a confined many-particle system described by the time-dependent Hamiltonian

\[
H(t) = \sum_{i=1}^{N} \left[ \frac{\vec{p}_i^2}{2m} + \frac{1}{2}m\omega(t)^2 \vec{r}_i^2 \right] + \sum_{i<j} V(\vec{r}_i - \vec{r}_j),
\]  

(1)

where the interparticle interaction potential exhibits the scaling property \( V(\lambda \vec{r}) = \lambda^{-2} V(\vec{r}) \). The dynamical group for the Hamiltonian (1) is \( SU(1,1) \) [45]. Given an arbitrary energy eigenstate
at $t = 0$, the exact dynamics generated by the Hamiltonian (1) for any modulation of the trapping frequency $\omega(t)$ is described by a scaling law \cite{15,16,46},

$$\Psi_n(\vec{r}_1, \ldots, \vec{r}_N, t) = b^{-\frac{N(D)}{2}} \exp \left[ i \frac{mb}{2\hbar b} \sum_{i=1}^{N} \vec{r}_i^2 - i \int_0^t \frac{E_n(0)}{\hbar \omega(t)} dt' \right] \Psi_n \left( \frac{\vec{r}_1}{b}, \ldots, \frac{\vec{r}_N}{b}, t = 0 \right),$$  \(2\)

where $n = \{ n_i : i = 1, \ldots, DN \}$ is a shorthand for a complete set of quantum numbers that labels the quantum state as well as the corresponding energy eigenvalue $E_n(0) = E_n[\omega(0)]$. The scaling factor $b = b(t) > 0$ as a function of time is the solution of the Ermakov equation

$$\ddot{b} + \omega(t)^2 b = \omega_0^2 b^{-3},$$  \(3\)

with the initial boundary conditions $b(0) = 1$ and $\dot{b}(0) = 0$, such that the time-dependent state (2) reduces to the initial eigenstate at the beginning of the process $t = 0$.

The time-dependent state (2) can be obtained by acting with a sequence of unitary transformations on the initial state. We follow Lohe \cite{36} and Jaramillo et al. \cite{24} to describe the time-evolution in terms of the action of two unitary transformations, elements of SU(1, 1). To this end, we introduce

$$T_{dil} = \exp \left[ -i \frac{\log b}{2\hbar} \sum_{i=1}^{N} (\vec{r}_i \cdot \vec{p}_i + \vec{p}_i \cdot \vec{r}_i) \right],$$  \(4\)

$$T_r = \exp \left[ i \frac{mb}{2\hbar b} \sum_{i=1}^{N} \vec{r}_i^2 \right].$$  \(5\)

We note that $T_{dil}$ implements scaling transformations by acting on a function in the coordinate representation $f(\vec{r}_1, \ldots, \vec{r}_N)$ according to $T_{dil} f(\vec{r}_1, \ldots, \vec{r}_N) T_{dil}^\dagger = f(\vec{r}_1/b, \ldots, \vec{r}_N/b)$. Similarly, acting on a function in the momentum representation $g(\vec{p}_1, \ldots, \vec{p}_N)$, one finds $T_{dil} g(\vec{p}_1, \ldots, \vec{p}_N) T_{dil}^\dagger = g(\vec{p}_1/b, \ldots, \vec{p}_N/b)$. By contrast, $T_r$ acts as a multiplicative factor in the coordinate representation while it generates translations in momentum space acting on a function $g(\vec{p}_1, \ldots, \vec{p}_N)$ as $T_r g(\vec{p}_1, \ldots, \vec{p}_N) T_r^\dagger = g(\vec{p}_1 - mb\vec{r}_1/b, \ldots, \vec{r}_N - mb\vec{r}_N/b)$.

Denoting $T = T_r T_{dil}$ and the Hamiltonian (1) at $t = 0$ by $H_0$, we identify the invariant of motion

$$\mathcal{I} = TH_0 T^\dagger$$

$$= \frac{N}{2} \left[ \frac{b^2 \vec{p}_i^2}{2m} + \frac{1}{2} m \vec{r}_i^2 \left( \frac{\omega(0)^2}{b^2} + b^2 \right) - \frac{bb}{2} (\vec{r}_i \cdot \vec{p}_i + \vec{p}_i \cdot \vec{r}_i) \right] + b^2 \sum_{i<j} V(\vec{r}_i - \vec{r}_j),$$  \(6\)

as an instance of the many-body invariants of motion known under scale-invariant driving \cite{16,17}. Equation (6) is indeed a generalization of the Lewis-Riesenfeld invariant of motion known for a single-particle driven harmonic oscillator \cite{47}, satisfies

$$\frac{\partial}{\partial t} \mathcal{I} = \frac{1}{\hbar} [H(t), \mathcal{I}],$$  \(7\)

and has time-dependent eigenfunctions $\Phi_n(\vec{r}_1, \ldots, \vec{r}_N, t) = T \Psi_n(\vec{r}_1, \ldots, \vec{r}_N, 0)$ and time-independent eigenvalues $E_n(0)$. In terms of the eigenvectors of $\mathcal{I}$ the time-evolving state (2) can be written as

$$\Psi_n(\vec{r}_1, \ldots, \vec{r}_N, t) = e^{i\alpha_t} T \Psi_n(\vec{r}_1, \ldots, \vec{r}_N, 0),$$  \(8\)

where $\alpha_t = - \int_0^t \frac{E_n(0)}{\hbar \omega(t)} dt'$ is a time-dependent dynamical phase.
3. Exact Nonadiabatic Mean Energy

Knowledge of the exact time-dependent dynamics under scale invariance makes possible to find the exact nonadiabatic mean energy of the system, as shown in [24]. To simplify the notation, we introduce the square position and squeezing operators

\[ Q = \sum_{i=1}^{N} \vec{r}_i^2, \quad C = \frac{1}{2} \sum_{i=1}^{N} (\vec{r}_i \cdot \vec{p}_i + \vec{p}_i \cdot \vec{r}_i) = \frac{1}{2} \sum_{i=1}^{N} \{\vec{r}_i, \vec{p}_i\}, \]

and note that the time-dependent Hamiltonian (1) can be rewritten as

\[ H(t) = \frac{1}{\hbar^2} \mathcal{I} + i\hbar \frac{\partial T}{\partial t} T^t, \]

where

\[ i\hbar \frac{\partial T}{\partial t} T^t = \left[ -\frac{1}{2} m \left( \frac{\dot{b}}{b} + \frac{\dot{b}^2}{b^2} \right) Q + \frac{b}{b} C \right]. \]

Using Equation (8), the expectation value of the time-dependent Hamiltonian in the time-evolving state reads

\[ \langle \Psi_n(t)|H(t)|\Psi_n(t)\rangle = \langle \Psi_n(0)|T^t H(t) T|\Psi_n(0)\rangle = \langle \Psi_n(0)| \left( \frac{H_0}{\hbar^2} + i\hbar T^t \frac{\partial T}{\partial t} \right)|\Psi_n(0)\rangle, \]

with

\[ i\hbar T^t \frac{\partial T}{\partial t} = \left[ m \left( \frac{\dot{b}}{b} - \frac{\dot{b}}{b} \right) Q + \frac{b}{b} C \right]. \]

As a result, the mean nonadiabatic energy is given by

\[ \langle \Psi_n(t)|H(t)|\Psi_n(t)\rangle = \frac{E_n(0)}{\hbar^2} - \frac{m}{2} \left( \dot{b} \dot{b} - \dot{b}^2 \right) \langle Q(0) | Q(0) \rangle + \frac{b}{b} \langle \Psi_n(0)|C|\Psi_n(0)\rangle. \]

As this expression holds for an arbitrary eigenstate, the expectation value can be also carried out with respect to an arbitrary state diagonal in the energy eigenbasis, such as an equilibrium thermal state. Specifically, the nonadiabatic mean-energy following a change of \( \omega(t) \) is found to be

\[ \langle H(t) \rangle_t = \frac{1}{\hbar^2} \langle H(0) \rangle + \frac{m}{2} \left( \dot{b} \dot{b} - \dot{b}^2 \right) \langle Q(0) | Q(0) \rangle + \frac{b}{b} \langle C(0) \rangle, \]

where \( \langle \bullet \rangle \equiv \operatorname{tr} (\bullet \rho_0) \) denotes the mean value at the time \( t = 0 \) with respect to a state \( \rho_0 \) diagonal in the eigenbasis of \( H(0) \). Similarly, \( \langle \bullet \rangle_t \equiv \operatorname{tr} (\bullet \rho_t) \) denotes the mean value at the time \( t \) with respect to the state \( \rho_t = U(t) \rho_0 U(t)^\dagger \), where \( U(t) \) is the time evolution operator. Under scale invariance it is thus possible to relate the evolution of the mean nonadiabatic energy to the initial expectation values of the Hamiltonian, square position and squeezing operators.

For an initial state at equilibrium, diagonal in the eigenbasis of \( H(0) \), the mean squeezing vanishes \( \langle C(0) \rangle = 0 \). In addition, the initial expectation value of \( Q \) is proportional to the mean energy of the system [24],

\[ \langle Q(0) \rangle = \frac{\langle H(0) \rangle}{m\omega_0^2}. \]

Using this property and the Ermakov Equation (15), the nonadiabatic mean energy reads

\[ \langle H(t) \rangle_t = Q^*(t) \frac{\omega(t)}{\omega_0} \langle H(0) \rangle = Q^*(t) \langle H(t) \rangle_{ad}, \]
where we have introduced the adiabatic mean energy \( \langle H(t) \rangle_{\text{ad}} = \langle H(0) \rangle \omega(t)/\omega(0) \) as well as the nonadiabatic factor \( Q^*(t) \) that accounts for the amount of energy excitations over the adiabatic dynamics

\[
Q^*(t) = \frac{\omega_0}{\omega(t)} \left[ \frac{1}{b(t)^2} + \frac{\dot{b}(t)^2 - b(t)\ddot{b}(t)}{2\omega_0^2} \right] = \frac{\omega_0}{\omega(t)} \left[ \frac{1}{2b(t)^2} + \frac{\omega(t)^2\dot{b}(t)^2}{2\omega_0^2} + \frac{\dot{b}(t)^2}{2\omega_0^2} \right]. \tag{18}
\]

Equations (17) and (18) provide the exact nonadiabatic mean energy of the family of systems governed by Hamiltonian (1). We emphasize that the nonadiabatic factor \( Q^*(t) \) is simply the ratio between the nonadiabatic and adiabatic mean energies. As it can be expressed in terms of the scaling factor, that governs the size of the atomic cloud, it is an experimentally accessible quantity \([23,31]\).

Summarizing these background results, reported in \([24]\), it suffices to know the evolution of the nonadiabatic and adiabatic mean energies. As it can be expressed in terms of the scaling factor, that accounts for the amount of energy excitations over the eigenstate, or any initial state \( |n\rangle \) that corresponds to the zero time which are model dependent (we have not been able to relate them to properties of the Hamiltonian via symmetries).

4. Nonadiabatic Energy Fluctuations

We next turn to the focus of our work, the characterization of the energy fluctuations. To this end, we consider the expectation value of the second momentum of the Hamiltonian at time \( t > 0 \) in the corresponding time-evolving state

\[
\langle \Psi_n(t)|H(t)|\Psi_n(t)\rangle = \langle \Psi_n(0)|T^tH(t)\Psi_n(0)\rangle = \langle \Psi_n(0)|(T^tH(t)T)^2|\Psi_n(0)\rangle. \tag{19}
\]

We recall that

\[
T^t H(t) T = \frac{H_0}{b^2} + i\hbar T^t \frac{\partial T}{\partial t} = \frac{H_0}{b^2} + \alpha Q + \beta C, \tag{20}
\]

where for compactness we define the time-dependent coefficients

\[
\alpha = \frac{m}{2} \left( b^2 - \dot{b}b \right), \quad \beta = \frac{\dot{b}}{b}. \tag{21}
\]

We thus find

\[
(T^t H(t) T)^2 = \frac{H_0^2}{b^4} + \alpha^2 Q^2 + \beta^2 C^2 + \alpha \frac{\dot{b}}{b^2} \{H_0, Q\} + \frac{\beta^2}{b^2} \{H_0, C\} + \alpha\beta \{Q, C\}. \tag{22}
\]

Evaluation of the expectation value of the fourth, fifth and the sixth terms in (22) over an energy eigenstate, or any initial state \( \rho_0 \) commuting with \( H_0 \), is straightforward and yields

\[
\langle \{H_0, Q\} \rangle = \frac{2\langle H_0^2 \rangle}{m\omega_0^2}, \tag{23a}
\]

\[
\langle \{H_0, C\} \rangle = 2\langle H_0 \rangle/\langle C \rangle = 0, \tag{23b}
\]

\[
\langle \{Q, C\} \rangle = \frac{i\hbar}{2} \langle \{Q, [H_0, Q] \} \rangle = \frac{i\hbar}{2} \langle [H_0, Q^2] \rangle = 0, \tag{23c}
\]

as \( \langle Q \rangle = \langle H_0 \rangle/m\omega_0^2 \) from equipartition theorem and \( \langle C \rangle = 0 \) in an equilibrium eigenstate. Although the mean value of the squeezing operator vanishes, the analogue quantity for its square is
where we remind that \( \langle \bullet \rangle \equiv \text{tr} (\bullet \rho_0) \) denotes the mean value at time \( t = 0 \). This is the main result of this work. As for the mean energy, the nonadiabatic energy fluctuations can be related to the initial expectation value of the second moment of the Hamiltonian, square position and squeezing operators. Once these quantities are known, nonadiabatic effects are simply encoded in the time-dependent scaling factor \( b(t) \), as the coefficients \( \alpha(t) \) and \( \beta(t) \) are directly determined by it.

At this stage we are ready to consider the adiabatic evolution, corresponding to the limit \( \dot{\omega}/\omega^2 \ll 1 \) \cite{Raghu2010}. At the level of the Ermakov equation, one can set \( \beta \approx 0 \) and \( \dot{b} \approx 0 \) and find that the scaling factor in the adiabatic limit is set by \( b_{\text{ad}}(t) = (\omega(0)/\omega(t))^{1/2} \). As a result, \( \alpha \approx \beta \approx 0 \) and the adiabatic evolution yields

\[
\langle H(t)^2 \rangle_t = \frac{\langle H(0)^2 \rangle}{b_{\text{ad}}^4} = \frac{\omega(t)^2}{\omega(0)^2} \langle H(0)^2 \rangle.
\]

5. Nonadiabatic Energy Fluctuations: Explicit Examples

5.1. Single-Particle Time-Dependent Quantum Harmonic Oscillator

Equation (24) determines exactly non-equilibrium energy fluctuation as a function of the initial expectation values at time \( t = 0 \). However, it is not possible in general to obtain a simple expression similar to (17) for the mean value of the non-equilibrium energy. In (23b) we use that the mean expectation values at time \( t \) operators, one finds

\[
H \rightarrow \text{evaluate } \langle \text{explicit computation model dependent and challenging, in general. To illustrate this observation, let us similar to (17) for the mean value of the non-equilibrium energy. In (23b) we use that the mean expectation values at time } t \text{ operators, one finds }
\]

\[
\alpha^2 \langle Q^2 \rangle + \beta^2 \langle C^2 \rangle
\]

\[\text{(24)}\]

where we remind that \( \langle \bullet \rangle \equiv \text{tr} (\bullet \rho_0) \) denotes the mean value at time \( t = 0 \). This is the main result of this work. As for the mean energy, the nonadiabatic energy fluctuations can be related to the initial expectation value of the second moment of the Hamiltonian, square position and squeezing operators. Once these quantities are known, nonadiabatic effects are simply encoded in the time-dependent scaling factor \( b(t) \), as the coefficients \( \alpha(t) \) and \( \beta(t) \) are directly determined by it.

At this stage we are ready to consider the adiabatic evolution, corresponding to the limit \( \dot{\omega}/\omega^2 \ll 1 \) \cite{Raghu2010}. At the level of the Ermakov equation, one can set \( \beta \approx 0 \) and \( \dot{b} \approx 0 \) and find that the scaling factor in the adiabatic limit is set by \( b_{\text{ad}}(t) = (\omega(0)/\omega(t))^{1/2} \). As a result, \( \alpha \approx \beta \approx 0 \) and the adiabatic evolution yields

\[
\langle H(t)^2 \rangle_t = \frac{\langle H(0)^2 \rangle}{b_{\text{ad}}^4} = \frac{\omega(t)^2}{\omega(0)^2} \langle H(0)^2 \rangle.
\]

\[\text{(25)}\]

The nonadiabatic expectation value of the second moment of the Hamiltonian then reads

\[
\langle \Psi_n(t)|H(t)^2|\Psi_n(t)\rangle = E_n(t)^2 + \frac{\hbar^2}{16\omega_0^2} (6n^2 + 6n + 3) \left[ \left( \frac{\hbar^2}{2b^2} + \frac{\omega_0^2}{b^2} \right)^2 + \frac{4\omega_0^2 b^2}{b^2} \right]
\]

\[\text{(26)}\]

where \( Q^* \) is given by Equation (18), \( E_n(t) = \hbar \omega(t)(n + \frac{1}{2}) \) is the instantaneous energy eigenvalue at time \( t \) and \( \langle H_1 \rangle_t = \langle \Psi_n(t)|H(t)|\Psi_n(t)\rangle = Q^* E_n(t) \) the corresponding nonadiabatic mean energy.
The noadiabatic energy variance $\Delta_n H(t)^2 \equiv \langle \Psi_n(t)|H(t)^2|\Psi_n(t)\rangle - \langle \Psi_n(t)|H(t)|\Psi_n(t)\rangle^2$ in a given eigenmode is therefore

$$\Delta_n H(t)^2 = \frac{\hbar^2 \omega(t)^2}{2}(n^2 + n + 1) \left[ (Q^*)^2 - 1 \right].$$

Using (24), our result reproduces the prediction by Husimi [48] and differs from that by Chen and Muga [49] (as the latter result does not vanish in the adiabatic limit, it is incorrect). We verify that in the adiabatic limit when $\dot{b} = 0$, $b = b_{ad}$ and $Q^* = 1$, energy fluctuations in a single driven energy mode vanish identically, as expected.

For a canonical thermal state as initial state, with density matrix $\rho_0 = e^{-\beta H(0)}/Z_0$ normalized in terms of the partition function $Z_0 = \text{tr}[e^{-\beta H(0)}] = (2 \sinh[\beta \hbar \omega_0 / 2])^{-1}$, one finds

$$\Delta H(t)^2 = \frac{\hbar^2 \omega(t)^2}{4} \frac{1}{\sinh[\beta \hbar \omega_0 / 2]^2} \left[ (Q^*)^2 + \left( (Q^*)^2 - 1 \right) \cosh[\beta \hbar \omega_0] \right].$$

5.2. Calogero-Sutherland Gas in a Time-Dependent Harmonic Trap

We next consider the nonadiabatic energy fluctuations in a driven many-body quantum system. We focus on the (rational) Calogero-Sutherland model describing one-dimensional bosons subject to inverse-square pairwise interactions [50,51]. The system Hamiltonian is

$$H(t) = \sum_{i=1}^{N} \left[ -\frac{\hbar^2}{2m} \frac{\partial^2}{\partial z_i^2} + \frac{\hbar^2}{2m} \omega(t)^2 z_i^2 \right] + \frac{\hbar^2}{2m} \sum_{i<j} g (g - 1) \frac{1}{(z_i - z_j)^2}.$$  (30)

In one dimension, interactions are inextricably linked to particle statistics. The notion of generalized exclusion statistics [52,53], complementary to the familiar exchange statistics, classifies particles according to their exclusion in Hilbert space. In this context, the Calogero-Sutherland model can be described as a non-interacting gas of particles obeying generalized exclusion statistics [54]. The constant $g \geq 0$ thus plays a dual role. While in the standard description it determines the strength of the inverse-square interactions, it also serves to parameterize generalized exclusion statistics. The value $g = 0$ corresponds to ideal bosons. The case $g = 1$ corresponds to hard-core bosons in the continuum, i.e., the so-called Tonks-Girardeau gas [55]. For $g = 1/2$ the system is sometimes referred to as composed of Haldane semions, and more generally particles are called geons for an arbitrary value $g$ of the exclusion parameter. The many-body wavefunction of the Calogero-Sutherland ground-state at $t = 0$ reads

$$\Psi_0(z_1, \ldots, z_N) = C_0 \prod_{i<k} |z_j - z_k|^g \prod_{i=1}^{N} \exp \left( -\frac{m\omega_0}{2\hbar} z_i^2 \right),$$

where $C_0$ is a known normalization constant, see e.g., [35], and the corresponding energy eigenvalue reads $E_0 = \hbar \omega_0 N[1 + g(N - 1)]/2$. Obviously, for $g = 0$ one recovers the wavefunction of an ideal Bose gas, while for $g = 1$ one recognizes the wavefunction of bosons in the Tonks-Girardeau regime [56]. The time evolution of (31) exhibits scale invariance [7,35] and takes the form of Equation (2) in one spatial dimension with $D = 1$. As the ground state is an energy eigenstate, the expectation value of the second moment of the Hamiltonian simply reads $\langle \hat{H}_0^2 \rangle = E_0^2$. For the computation of $\langle Q^2 \rangle$ and $\langle C^2 \rangle$ we note that these moments can be derived from the corresponding characteristic function, which can be written down as an overlap between ground-state wavefunctions. In particular, as the ground-state is a homogeneous function times a Gaussian term, the explicit computation of the overlap is possible [35,57].
We first proceed to the computation of \( \langle \hat{Q}^2 \rangle \) by introducing the overlap

\[
A_Q(\lambda) = \langle \Psi_0 | e^{-\lambda \hat{Q}} | \Psi_0 \rangle = \left[ 1 + \lambda \frac{\hbar}{m \omega_0} \right]^{-\lambda^2 \frac{1}{2} \left[ 1 + g(N-1) \right]},
\]

from which moments of \( \hat{Q} \) are found as

\[
\langle \Psi_0 | \hat{Q}^k | \Psi_0 \rangle = (-1)^k \frac{d^k}{d\lambda^k} A_Q(\lambda) \big|_{\lambda=0}.
\]

Their explicit expressions are given by

\[
\langle \Psi_0 | \hat{Q} | \Psi_0 \rangle = \left( \frac{\hbar}{m \omega_0} \right) c,
\]

\[
\langle \Psi_0 | \hat{Q}^2 | \Psi_0 \rangle = \left( \frac{\hbar}{m \omega_0} \right)^2 c(1+c),
\]

\[
\langle \Psi_0 | \hat{Q}^k | \Psi_0 \rangle = \left( \frac{\hbar}{m \omega_0} \right)^k \prod_{j=1}^k (c + j - 1),
\]

where for compactness we have defined \( c = \frac{N}{2} \left[ 1 + g(N-1) \right] \).

Similarly, we can evaluate the term \( \langle C^2 \rangle \). To this end, we first notice that the dilation operator acts on a many-body wavefunction as

\[
T_{\text{dil}} | \Psi(z_1, \ldots, z_N) \rangle = b^{-\frac{N}{2}} \Psi \left( \frac{z_1}{b}, \ldots, \frac{z_N}{b} \right),
\]

We next evaluate the overlap

\[
A_C(b) = \langle \Psi_0 | T_{\text{dil}} \Psi_0 \rangle = \left[ b \left( \frac{1 + \frac{1}{b^2}}{2} \right) \right]^{-c},
\]

which can be understood as the characteristic function of the eigenvalue distribution of the squeezing operator. Thus, moments of the generator can be found using the identity

\[
\langle \Psi_0 \left[ \sum_{l=1}^N \frac{1}{2} \left( z_l p_l + p_l z_l \right) \right]^k \rangle_{\Psi_0} = (-i \hbar)^k \left( b \frac{d}{db} \right)^k A_0(b) \big|_{b=1}.
\]

As \( A_C(b) \) is a real function, this implies that odd moments \( \langle \Psi_0 | \hat{C}^k | \Psi_0 \rangle \) identically vanish in the ground state \( | \Psi_0 \rangle \). By explicit evaluation, we find

\[
\langle \Psi_0 | \hat{C}^2 | \Psi_0 \rangle = \hbar^2 c,
\]

\[
\langle \Psi_0 | \hat{C}^4 | \Psi_0 \rangle = \hbar^4 c(2 + 3c),
\]

\[
\langle \Psi_0 | \hat{C}^6 | \Psi_0 \rangle = \hbar^6 c(16 + 30c + 15c^2),
\]

Thus, the second moment of the Hamiltonian is set by

\[
\langle H_2^2 \rangle = E_0^2 \left( \frac{1}{\beta^2} + \frac{2 \alpha}{m \omega_0^2 \beta^2} \right) + \hbar^2 \left( \frac{\hbar}{m \omega_0} \right)^2 c(1+c) + \beta^2 \hbar^2 c
\]

\[
= \hbar^2 \omega(t)^2 c^2 \left( Q^* \right)^2 + \hbar^2 \omega(t)^2 c \left[ (Q^*)^2 - 1 \right],
\]
where $Q^*$ is given by Equation (18) and $E_0 = \hbar \omega_0 c$. As a result, the variance simply reads

$$\Delta H(t)^2 = \hbar^2 \omega(t)^2 c \left[ (Q^*)^2 - 1 \right],$$

and scales at most quadratically with the number of particles [35].

The exact many-body wavefunctions of the Calogero-Sutherland model are known for excited states and can be written as the product of the ground-state wavefunction and a combination of Hermite polynomials, see e.g., [58]. Due to the later, the wavefunction of an arbitrary excited state is no longer a homogeneous function times a Gaussian term, precluding the computation of the energy fluctuations via the method put forward.

5.3. Unitary Fermi Gas in a Time-Dependent Harmonic Trap

We next consider a strongly interacting degenerate Fermi gas with two spin components [59]. The unitary limit is reached when the interaction strength between atoms becomes divergent, e.g., at a Feshbach resonance. To characterize the spectrum and eigenstates of the Hamiltonian, the interactions can be replaced by a contact condition in terms of a function $A$ such that

$$\psi(\vec{r}_1, \ldots, \vec{r}_N) = A(\vec{R}_{ij}, \{\vec{r}_k : k \neq i, j\}) + O(r_{ij}),$$

where $\vec{R}_{ij} = (\vec{r}_i + \vec{r}_j)/2$ is the center of mass position for particles $i$ and $j$, and the limit of vanishing relative distance between two particles $r_{ij} = |\vec{r}_i - \vec{r}_j| \rightarrow 0$ is taken. Tan, Werner and Castin have shown that the low-energy wavefunctions of a unitary Fermi gas in an isotropic trap are the product of a Gaussian factor times a homogeneous function [60,61]. This result together with the scale invariance exhibited by this system [11] will make possible the exact computation of the nonadiabatic energy fluctuations.

In three spatial dimensions, we introduce the 3N vector $\vec{X} = (\vec{r}_1, \ldots, \vec{r}_N)$. Similarly, we denote by $\vec{P} = (\vec{p}_1, \ldots, \vec{p}_N)$ the 3N vector of the conjugated momenta. Denoting the norm of $\vec{X}$ by $X = \sqrt{\sum r_j^2}$ and the unit vector $\hat{n} = \vec{X}/X$, we make use of the hyperspherical coordinates $(X, \hat{n})$. Eigenstates in a trap can be related to zero-energy eigenstates in free space. We follow Werner and Castin [61] and denote by $\psi_0^0(X)$ a zero-energy eigenstate in free space. The latter exhibits scale invariance

$$\psi_0^0(\vec{X}/b) = b^{-\nu} \psi_0^0(\vec{X}),$$

where $\nu$ is the degree of homogeneity, i.e., $\vec{X} \cdot P \psi_0 = -i\hbar \nu \psi_0$ [60]. The corresponding trapped eigenstate reads

$$\psi_\nu(X) = e^{-\frac{m \omega_0 X^2}{\hbar} \frac{E}{\bar{E}} - \frac{3N}{2} f(\hat{n})},$$

and its energy eigenvalue is given by

$$E(0) = \hbar \omega_0 (\nu + 2q - 3N/2),$$

in terms of a non-negative integer $q$. Energy eigenstates are connected by raising and lowering operators associated with the $SO(2, 1)$ Pitaevskii-Rosch symmetry. The lowest energy eigenstate of a given tower of states, that with $q = 0$ that is annihilated by the lowering operator, can be shown to take the form

$$\psi_\nu(\vec{X}) = Ce^{-\frac{m \omega_0 X^2}{\hbar}} X^{\frac{\nu}{2}} f(\hat{n}),$$

where $C$ is a normalization constant. This completes our review of [60,61].
By scale invariance, the time-dependent state following a modulation of the trapping frequency is of the form in Equation (2). As it turns out, this suffices to compute the nonadiabatic expectation value of the first and second energy moments of the Hamiltonian of a unitary Fermi gas initially prepared in the low-energy state (49). To this end, we note that the characteristic functions of \( Q = X^2 \) and \( C = (\vec{X} \cdot \vec{P} + \beta \cdot \vec{X}) / 2 \) are

\[
A_Q(\lambda) = \langle \psi_\nu | e^{-\lambda Q} | \psi_\nu \rangle = \left[ 1 + \lambda \frac{\hbar}{m \omega_0} \right]^{\frac{E(0)}{\hbar \omega_0}}, \tag{50}
\]

\[
A_C(b) = \langle \psi_\nu | T_{\text{dil}} \psi_\nu \rangle = \left[ \frac{b}{2} \left( 1 + \frac{1}{b^2} \right) \right]^{\frac{E(0)}{\hbar \omega_0}}, \tag{51}
\]

and thus

\[
\langle \psi_\nu | Q^2 | \psi_\nu \rangle = \left( \frac{\hbar}{m \omega_0} \right)^2 \frac{E(0)}{\hbar \omega_0} \left( 1 + \frac{E(0)}{\hbar \omega_0} \right), \tag{52}
\]

\[
\langle \psi_\nu | C^2 | \psi_\nu \rangle = \hbar^2 \frac{E(0)}{\hbar \omega_0}, \tag{53}
\]

whence it follows

\[
\langle H^2_t \rangle_t = E(0)^2 \left[ \frac{1}{b^4} \frac{b^2 - \bar{b} \bar{b}}{\omega_0^2 b^2} \right] + \frac{\hbar^2 (\bar{b}^2 - \bar{b} \bar{b})^2 E(0)}{4 \omega_0^2 \hbar \omega_0} \left( 1 + \frac{E(0)}{\hbar \omega_0} \right) + \frac{\hbar^2 \left( \frac{b}{\bar{b}} \right)^2}{4 \omega_0^2} \left( \bar{b}^2 + \omega(t) \bar{b}^2 - \omega_0^2 \frac{b^2}{\bar{b}^2} \right)^2. \tag{54}
\]

It is convenient to rewrite this expression in the form

\[
\langle H^2_t \rangle_t = \hbar^2 \omega(t)^2 \left( \frac{E(0)}{\hbar \omega_0} \right)^2 + \frac{\hbar^2 E(0)}{\hbar \omega_0} \left( 1 + \frac{E(0)}{\hbar \omega_0} \right) \left[ \left( \frac{b}{\bar{b}} \right)^2 + \frac{1}{4 \omega_0^2} \left( \bar{b}^2 + \omega(t) \bar{b}^2 - \omega_0^2 \frac{b^2}{\bar{b}^2} \right)^2 \right], \tag{55}
\]

in parallel with the results of the single-particle harmonic oscillator as well as that of the Calogero-Sutherland model with the corresponding ground states as initial state. This last expression can be rewritten as follows

\[
\langle H^2_t \rangle_t = \hbar^2 \omega(t)^2 \left( \frac{E(0)}{\hbar \omega_0} \right)^2 + \hbar^2 \omega(t)^2 \frac{E(0)}{\hbar \omega_0} \left( 1 + \frac{E(0)}{\hbar \omega_0} \right) \left[ (Q^*)^2 - 1 \right]
\]

\[
= \hbar^2 \omega(t)^2 \left( \frac{E(0)}{\hbar \omega_0} \right)^2 (Q^*)^2 + \hbar^2 \omega(t)^2 \frac{E(0)}{\hbar \omega_0} \left[ (Q^*)^2 - 1 \right], \tag{56}
\]

with \( Q^* \) given by Equation (18), leading to the expression of the energy variance in the form

\[
\langle H^2_t \rangle_t - \langle H_t \rangle_t^2 = \hbar^2 \omega(t)^2 \frac{E(0)}{\hbar \omega_0} \left[ (Q^*)^2 - 1 \right]. \tag{57}
\]

The expression of the results in terms of the ratio \( \frac{E(0)}{\hbar \omega_0} \) is particularly convenient and it also holds for the results derived in the previous subsections for the ground states of the harmonic oscillator and the Calogero-Sutherland model. As a result, we shall use it hereafter without specifying the system under consideration. Beyond these models, our results hold whenever the initial state is the product of a Gaussian term times a homogeneous function. We further note that

\[
\frac{E(0)}{\hbar \omega_0} = \langle \psi_\nu | \hat{Q} | \psi_\nu \rangle \left( \frac{m \omega_0}{\hbar} \right) = \frac{1}{x_0^2} \langle \psi_\nu | X^2 | \psi_\nu \rangle = \sigma^2, \tag{58}
\]
which is but the initial size of the cloud in units of \( x_0 = \sqrt{\frac{\hbar}{m\omega_0}} \), and is thus an experimentally measurable quantity.

6. Nonadiabatic Moments of the Square Position Operator

Due to the \( SL(1,1) \) dynamical symmetry group, we have related moments of the Hamiltonian \( H(t) \) to those of the operators \( H(0), Q \) and \( C \) at \( t = 0 \). For completeness, we note as well the nonadiabatic evolution of the moments of the square position operator \( Q \). To this end, we introduce

\[
A_Q(\lambda, t) = \langle \Psi_0(t)|e^{-\lambda \hat{Q}}|\Psi_0(t)\rangle = \langle \Psi_0(0)|e^{-\lambda \hat{Q}T_T}|\Psi_0(0)\rangle. \tag{59}
\]

Noticing that \( T_T^\dagger \hat{Q}T_T = Qb^2 \), one finds

\[
A_Q(\lambda, t) = \left[ 1 + \lambda \frac{h b^2}{m \omega_0} \right]^{-\frac{E(0)}{\hbar \omega_0}}. \tag{60}
\]

As expected, the \( k \)-th momentum of the position operator scales in a trivial form,

\[
\langle \Psi_0(t)|\hat{Q}^k|\Psi_0(t)\rangle = b^{2k} \langle \Psi_0(0)|\hat{Q}^k|\Psi_0(0)\rangle, \tag{61}
\]

and is proportional to the corresponding moment at \( t = 0 \), in agreement with the definition of the scaling factor \( b(t) \).

7. Nonadiabatic Moments of the Squeezing Operator

The evolution of the squeezing operator \( C \) is more involved than that of \( Q \), and as with the non-adiabatic energy, we shall only compute the first two moments explicitly. We first note that

\[
T_T^\dagger CT_T = T_T^\dagger \sum_{i=1}^{N} \left\{ \vec{r}_i, \vec{p}_i + \frac{m b}{b} \vec{r}_i \right\} T_T^\dagger = C + m b b Q. \tag{62}
\]

The first moment reads

\[
\langle \Psi_0(t)|C|\Psi_0(t)\rangle = \langle \Psi_0(0)|C + m b b Q|\Psi_0(0)\rangle = m b b \langle \Psi_0(0)|Q|\Psi_0(0)\rangle = \frac{\hbar}{\omega_0} b b \frac{E(0)}{\hbar \omega_0}. \tag{63}
\]

while the second one is given by

\[
\langle \Psi_0(t)|C^2|\Psi_0(t)\rangle = \langle \Psi_0(0)|C + m b b Q|^2|\Psi_0(0)\rangle = \langle \Psi_0(0)|C^2|\Psi_0(0)\rangle + m^2 \hbar b^2 \langle \Psi_0(0)|Q^2|\Psi_0(0)\rangle + m b b \langle \Psi_0(0)|\{C, Q\}|\Psi_0(0)\rangle = \frac{\hbar^2}{\omega_0} \frac{E(0)}{\hbar \omega_0} \left( \frac{b^2 \hbar^2}{\omega_0^2} \frac{E(0)}{\hbar \omega_0} \right), \tag{64}
\]

after using \( \langle \{Q, C\} \rangle = 0 \).

8. Driving Protocols

In what follows, we analyze the exact nonadiabatic energy fluctuations for specific instances of driving protocols of the time-dependent harmonic frequency.

8.1. Free Expansion

A standard protocol for time-of-flight imaging of ultracold atoms relies on suddenly releasing a cloud initially trapped in a harmonic potential with frequency \( \omega_0 \). For \( t > 0 \), \( \omega(t) = 0 \). For an
initial equilibrium state at \( t > 0 \), the initial conditions \( b(0) = 1 \) and \( \dot{b}(0) = 0 \) are required for the time-dependent state (2) to reduce to the initial state. The solution of the Ermakov equation is then
\[
b(t) = \sqrt{1 + \omega_0^2 t^2} \]
In this case there is no reference notion of adiabatic dynamics and \( Q^* \) is not defined. The mean energy satisfies
\[
\frac{\langle H(t) \rangle_t}{\langle H(0) \rangle} = \frac{1}{2b(t)^2} + \frac{\omega(t)^2 b(t)^2}{2\omega_0^2} + \frac{\dot{b}(t)^2}{2\omega_0^2} = 1, \tag{65}
\]
i.e., the mean total energy is preserved during time of flight. The same holds true for the expectation value of all the moments of the Hamiltonian. In turn, for an energy eigenstate \( \langle H_0^2 \rangle = \langle H_0 \rangle^2 \) and thus \( \langle H^2_1 \rangle_t - \langle H_1 \rangle_t^2 = 0 \) for all \( t \geq 0 \).

8.2. Sudden Quenches

Consider the sudden modulation of the trap frequency so that it varies from an initial value \( \omega_0 \) for \( t < 0 \) to a final value \( \omega_f \). Sudden quenches of this type have been studied in a variety of applications such as probing quantum correlations in ultracold gases, the study of their relaxation and thermalization [9,10,62], the analysis of Loschmidt echoes and work statistics [63–66], and finite-time quantum thermodynamics of thermodynamic cycles [24,67–72], among other examples.

The solution of the Ermakov equation with \( b(0) = 1 \) and \( \dot{b}(0) = 0 \) leads to the scaling factor that exhibits undamped oscillations [9]
\[
b(t) = \sqrt{1 + (\omega_0^2 / \omega_f^2) - 1} \sin^2(\omega_f t). \tag{66}
\]

The nonadiabatic factor \( Q^* \) accounting for quantum friction reads [24]
\[
Q^*(t) = \frac{\omega_0^2 + \omega_f^2}{2\omega_0 \omega_f}. \tag{67}
\]

As a result,
\[
\langle H^2_1 \rangle_t - \langle H_1 \rangle_t^2 = \hbar^2 \omega_f^2 \frac{E(0)}{\hbar \omega_0} \frac{(\omega_0^2 - \omega_f^2)^2}{2\omega_0 \omega_f^2}. \tag{68}
\]

8.3. Linear Frequency Ramp

As an example of a generic nonadiabatic driving we consider a linear ramp of the frequency such that \( \omega(t) = \omega_0 \) for \( t < 0 \), \( \omega(t) = \omega_0 + (\omega_f - \omega_0) t / t_f \) for \( 0 < t < t_f \) and \( \omega(t) = \omega_f \) for \( t \geq t_f \). Such time-dependence is nonanalytic at \( t = \{0, t_f\} \) and is thus expected to give rise to nonadiabatic dynamics [73]. The limit of \( t_f \to 0 \) naturally reduces to a sudden quench. The scaling factor can be obtained by solving numerically the Ermakov equation subject to \( b(0) = 1 \) and \( \dot{b}(0) = 0 \) and is shown in Figure 1. The nonadiabatic factor increases monotonically as a function of time with the final mean-energy being four times higher than the corresponding adiabatic value. The energy variance exhibits as well a monotonic behavior, with lower values reflecting the expansion nature of the process, during which the instantaneous trapping frequency decreases. Such features are however dependent on the parameters chosen and a non-monotonic behavior of \( b(t) \), \( Q^* \) and \( \Delta H(t)^2 \) is observed whenever \( t_f \) is of the order of the inverse of the frequencies involved in the process, e.g., \( \omega_0^{-1} \) and \( \omega_f^{-1} \).

We note that exact closed-form solutions of the Ermakov equation can be constructed as discussed in [74].
8.4. Shortcuts to Adiabaticity by Reverse Engineering

Reverse engineering provides a convenient approach for the fast driving of a system from an initial equilibrium state towards a final nonequilibrium state, without relying on adiabatic strategies. Fixing the duration of the protocol to be \( t_F \), with initial and final frequencies \( \omega_0 \) and \( \omega_F \), we proceed to impose the following boundary conditions on the scaling factor: \( b(0) = 1 \) and \( b(t_F) = b_{ad}(t_F) = \sqrt{\omega_0/\omega_F} \). Additional boundary conditions are required for the exact solution (2) to reduce to a stationary state. The vanishing of the derivatives of the scaling factor \( b(0) = \dot{b}(t_F) = 0 \) make sure that \( Q^* = 1 \) at \( t = \{0, t_F\} \). It proves convenient to further impose \( \ddot{b}(0) = 0 \). With these conditions one can then fix an interpolating polynomial ansatz for \( b(t) \) and determine the driving protocol \( \omega(t) \), using the Ermakov equation as \( \omega(t)^2 = \omega_0/b^4 - \dot{b}/b \). This is the essence to design a shortcut to adiabaticity by reverse engineering the scale-invariant dynamics [14,15,28], as experimentally demonstrated in [18,19,22]. Taking the interpolating ansatz to be \( b(t) = 1 + 10(t/t_F)^3(b(t_F) - 1) - 15(t/t_F)^4(b(t_F) - 1) + 6(t/t_F)^5(b(t_F) - 1) \), Figure 2 shows the modulation of the frequency-squared \( \omega(t)^2 \), nonadiabatic factor \( Q^* \) and nonadiabatic energy fluctuations for a protocol with \( t_F = 10/\omega_0 \) leading to an expansion with final \( b(t_F) = 4 \). The duration time is chosen such that \( \omega(t)^2 \geq 0 \) for all \( 0 \leq t \leq t_F \), preventing the inversion of the trap associated with values \( \omega(t)^2 \leq 0 \) required for shorter processes in which \( t_F \sim \omega_0^{-1} \). The nonadiabatic nature of the protocol is manifested in the values of \( Q^* \) exceeding unity for \( 0 < t < t_F \). Nonetheless, \( Q^* = 1 \) and \( \Delta H(t)^2 = 0 \) for \( t = \{0, t_F\} \) showing that at the beginning and end of the protocol excitations are cancelled out and the nonadiabatic evolution reduces to a stationary energy eigenstate.

8.5. Local Counterdiabatic Driving

Consider an arbitrary modulation of the trapping frequency \( \omega_R(t) \) as a reference, interpolating between a prescheduled initial value \( \omega_R(0) = \omega_0 \) and a final one \( \omega_R(t_F) = \omega_F \).
Such modulation is generally nonadiabatic. Shortcuts to adiabaticity engineered by local counterdiabatic driving [16,22,23,28,31] suppress nonadiabatic excitations by implementing instead the driving protocol with frequency $\omega(t)$, where

$$
\omega(t)^2 = \omega_R(t)^2 - \frac{3}{4} \left( \frac{\dot{\omega}_R(t)}{\omega_R(t)} \right)^2 + \frac{1}{2} \frac{\dot{\omega}_R(t)}{\omega_R(t)}.
$$

(69)

Driving the system with $\omega(t)$, it follows from the Ermakov Equation (3) that the scaling factor is precisely

$$
b(t) = \sqrt{\omega_0 / \omega_R(t)}.
$$

(70)

With the boundary conditions, $\dot{\omega}_R(0) = \ddot{\omega}_R(0) = 0$ and $\dot{\omega}_R(t_F) = \ddot{\omega}_R(t_F) = 0$, we choose the interpolating ansatz $\omega_R(t) = \omega_0 + 10(t/t_F)^3(\omega_F - \omega_0) - 15(t/t_F)^4(\omega_F - \omega_0) + 6(t/t_F)^5(\omega_F - \omega_0)$. It is then guaranteed that $\omega(0) = \omega_R(0) = \omega_0$ and $\omega(t_F) = \omega_R(t_F) = \omega_F$, so the local counterdiabatic protocol $\omega(t)$ still interpolates between the same initial and final frequencies considered in the reference protocol $\omega_R(t)$. Using Equations (69) and (70), the nonadiabatic factor (18) reads in this case

$$
Q^* = \frac{\omega_R(t)}{\omega(t)} \left[ 1 + \frac{1}{4} \left( \frac{\dot{\omega}_R(t)}{\omega_R(t)} - \frac{\ddot{\omega}_R(t)^2}{\omega_R(t)^4} \right) \right].
$$

(71)

This result corrects the expression found in [28,75]. The nonadiabatic energy variance takes the form

$$
\langle H_0^2 \rangle_t - \langle H_0 \rangle_t^2 = \hbar^2 \omega(t)^2 \frac{E(0)}{\hbar \omega_0} \left[ (Q^*)^2 - 1 \right],
$$

(72)

and vanishes identically at $t = \{0, t_F\}$. Figure 3 shows a polynomial modulation of the reference frequency-square $\omega_R^2(t)$ that decreases monotonically during the process from $\omega_0$ to $\omega_F$. By contrast, the modulation of $\omega(t)^2$ associated with the local counterdiabatic driving protocol is highly non-monotonic, involving a fast expansion, followed by a compression and an expansion again. This specific modulation as a function of time strongly depends on the duration of the process and is shown here for $t_F = 2/\omega_0$ and an expansion with $\omega_F = \omega_0/2$. The corresponding nonadiabatic factor $Q^*$ is shown to be greater than unity but at the beginning and end of the protocol when excitations are cancelled. The nonadiabatic energy variance grows from zero value and fluctuates during the process until its completion, when it vanishes again, as the state at $t = t_F$ is an eigenstate of the final Hamiltonian.

Figure 3. Local counterdiabatic driving of the scale-invariant dynamics. Given a reference modulation of the trapping frequency $\omega_R(t)$ between the initial and final trap configurations, a shortcut to adiabaticity by local counterdiabatic driving implements a modulation of the trapping frequency $\omega(t)$ that differs from the reference one $\omega_R(t)$. The reference modulation of the frequency square, the frequency square under the local counterdiabatic driving, the corresponding nonadiabatic factor and energy fluctuations are shown from left to right for an expansion with final frequency $\omega_F = \omega_0/2$ and duration $t_F = 2/\omega_0$. The latter is chosen long enough so that $\omega^2(t) > 0$ during the process.
9. Conclusions and Discussion

Scale invariance is an important symmetry that governs the dynamics of a variety of systems, including many harmonically-trapped ultracold atomic gases. It is not only of relevance to time-of-flight imaging techniques, but it also plays an important role in the implementation of quantum control techniques and the characterization of finite-time thermodynamics. We have shown that nonadiabatic energy fluctuations in many-body systems under scale-invariant driving can be simply expressed in terms of the equilibrium properties of the initial state. In particular, knowledge of expectation value of the second moments of the initial Hamiltonian, square position and squeezing operators suffices to characterize the exact nonadiabatic dynamics in terms of the evolution of the scaling factor, which governs the size of the atomic cloud. As the later is an experimentally measurable quantity, our results provide a way to characterize nonadiabatic energy fluctuations in the laboratory using time-of-flight imaging techniques. To illustrate our results, we have considered the single-particle quantum oscillator of relevance to experiments with trapped ions and thermal atomic clouds, as well as two instances of many-body systems: the Calogero-Sutherland gas and the unitary Fermi gas. The corresponding nonadiabatic evolution has been characterized in expansion protocols involving sudden quenches and finite-time protocols, including linear frequency ramps, as well as shortcuts to adiabaticity leading to stationary states upon completion of the process.

The exact results for the nonadiabatic energy fluctuations we have provided are expected to have widespread applications in both theoretical and experimental studies of these systems. For example, they can be used to characterize far-from-equilibrium dynamics in ultracold gases, quantify the cost of quantum control such as the use of optimal protocols \cite{68,76–78} and shortcuts to adiabaticity \cite{41,42}, characterize the performance of devices and processes in quantum thermodynamics \cite{30}, and study the ultimate speed limits \cite{37,38}, quantum decay \cite{35} and orthogonality catastrophe \cite{79} under scale-invariant quantum dynamics.

Author Contributions: Conceptualization, A.d.C.; Formal analysis, M.B. and A.d.C.; Investigation, M.B. and A.d.C.; Methodology, M.B. and A.d.C.; Visualization, A.d.C.; Writing—original draft, A.d.C.; Writing—Review & Editing, A.d.C. All authors have read and agreed to the published version of the manuscript.

Funding: This research was partially funded by the John Templeton Foundation.

Acknowledgments: It is a pleasure to acknowledge discussions with Juan Jaramillo and Doha Mesnaoui, comments on the manuscript by Fernando J. Gómez-Ruiz, and insightful clarifications by Iñigo L. Egusquiza.

Conflicts of Interest: The authors declare no conflict of interest.

References

1. Dalfovo, F.; Giorgini, S.; Pitaevskii, L.P.; Stringari, S. Theory of Bose-Einstein condensation in trapped gases. *Rev. Mod. Phys.* 1999, 71, 463–512. [CrossRef]

2. Giorgini, S.; Pitaevskii, L.P.; Stringari, S. Theory of ultracold atomic Fermi gases. *Rev. Mod. Phys.* 2008, 80, 1215–1274. [CrossRef]

3. Castin, Y.; Dum, R. Bose-Einstein Condensates in Time Dependent Traps. *Phys. Rev. Lett.* 1996, 77, 5315–5319. [CrossRef] [PubMed]

4. Kagan, Y.; Surkov, E.L.; Shlyapnikov, G.V. Evolution of a Bose-condensed gas under variations of the confining potential. *Phys. Rev. A* 1996, 54, R1753–R1756. [CrossRef]

5. Castin, Y.; Dum, R. Low-temperature Bose-Einstein condensates in time-dependent traps: Beyond the U(1) symmetry-breaking approach. *Phys. Rev. A* 1998, 57, 3008–3021. [CrossRef]

6. Pitaevskii, L.P.; Rosch, A. Breathing modes and hidden symmetry of trapped atoms in two dimensions. *Phys. Rev. A* 1997, 55, R853–R856. [CrossRef]

7. Sutherland, B. Exact Coherent States of a One-Dimensional Quantum Fluid in a Time-Dependent Trapping Potential. *Phys. Rev. Lett.* 1998, 80, 3678–3681. [CrossRef]

8. Rigol, M.; Muramatsu, A. Fermionization in an Expanding 1D Gas of Hard-Core Bosons. *Phys. Rev. Lett.* 2005, 94, 240403. [CrossRef]
9. Minguzzi, A.; Gangardt, D.M. Exact Coherent States of a Harmonically Confined Tonks-Girardeau Gas. *Phys. Rev. Lett.* 2005, 94, 240404. [CrossRef]

10. Del Campo, A. Fermionization and bosonization of expanding one-dimensional anyonic fluids. *Phys. Rev. A* 2008, 78, 045602. [CrossRef]

11. Castin, Y. Exact scaling transform for a unitary quantum gas in a time dependent harmonic potential. *C. R. Phys.* 2004, 5, 407–410. [CrossRef]

12. Deng, S.; Shi, Z.Y.; Diao, P.; Yu, Q.; Zhai, H.; Qi, R.; Wu, H. Observation of the Efimovian expansion in scale-invariant Fermi gases. *Science* 2016, 353, 371–374. [CrossRef] [PubMed]

13. Masuda, S.; Nakamura, K. Fast-forward of adiabatic dynamics in quantum mechanics. *Proc. R. Soc. A Math. Phys. Eng. Sci.* 2010, 466, 1135–1154. [CrossRef]

14. Del Campo, A. Frictionless quantum quenches in ultracold gases: A quantum-dynamical microscope. *Phys. Rev. A* 2011, 84, 031606. [CrossRef]

15. Schaff, J.F.; Song, X.L.; Capuzzi, P.; Vignolo, P.; Labeyrie, G. Fast optimal transition between two equilibrium states. *New J. Phys.* 2016, 18, 075019. [CrossRef] [PubMed]

16. Choi, S.; Onofrio, R.; Sundaram, B. Optimized sympathetic cooling of atomic mixtures via fast adiabatic strategies. *Phys. Rev. A* 2011, 84, 051601. [CrossRef]

17. Choi, S.; Onofrio, R.; Sundaram, B. Squeezing and robustness of frictionless cooling strategies. *Phys. Rev. A* 2012, 86, 043436. [CrossRef]

18. Beau, M.; Jaramillo, J.; del Campo, A. Scaling-Up Quantum Heat Engines Efficiently via Shortcuts to Adiabaticity. *Entropy* 2016, 18, 168. [CrossRef] [PubMed]
34. Merloti, K.; Dubessy, R.; Longchambon, L.; Olshanii, M.; Perrin, H. Breakdown of scale invariance in a quasi-two-dimensional Bose gas due to the presence of the third dimension. *Phys. Rev. A* 2013, 88, 061603. [CrossRef]

35. Del Campo, A. Exact quantum decay of an interacting many-particle system: The Calogero–Sutherland model. *New J. Phys.* 2016, 18, 015014. [CrossRef]

36. Lohe, M.A. Exact time dependence of solutions to the time-dependent Schrödinger equation. *J. Phys. A Math. Theor.* 2008, 42, 035307. [CrossRef]

37. Mandelstam, L.; Tamm, I. Quantum speed limits: From Heisenberg’s uncertainty principle to optimal quantum control. *J. Phys. USSR* 1945, 9, 249.

38. Shanahan, B.; Chenu, A.; Margolus, N.; del Campo, A. Quantum Speed Limits across the Quantum-to-Classical Transition. *Phys. Rev. Lett.* 2018, 120, 070401. [CrossRef] [PubMed]

39. Okuyama, M.; Ohzeki, M. Quantum Speed Limit is Not Quantum. *Phys. Rev. Lett.* 2018, 120, 070402. [CrossRef] [PubMed]

40. Anandan, J.; Aharonov, Y. Geometry of quantum evolution. *Phys. Rev. Lett.* 1990, 65, 1697–1700. [CrossRef]

41. Demirplak, M.; Rice, S.A. On the consistency, extremal, and global properties of counterdiabatic fields. *J. Chem. Phys.* 2008, 129, 154111. [CrossRef]

42. Campbell, S.; Deffner, S. Trade-Off Between Speed and Cost in Shortcuts to Adiabaticity. *Phys. Rev. Lett.* 2017, 118, 100601. [CrossRef] [PubMed]

43. Carlini, A.; Hosoya, A.; Koike, T.; Okudaira, Y. Time-Optimal Quantum Evolution. *Phys. Rev. Lett.* 2006, 96, 060503. [CrossRef] [PubMed]

44. Takahashi, K. How fast and robust is the quantum adiabatic passage? *J. Phys. A Math. Theor.* 2013, 46, 315304. [CrossRef]

45. Gambardella, P.J. Exact results in quantum many-body systems of interacting particles in many dimensions with SU(1,1) as the dynamical group. *J. Math. Phys.* 1975, 16, 1172–1187. [CrossRef]

46. Girardeau, M. Relationship between Systems of Impenetrable Bosons and Fermions in One Dimension. *J. Math. Phys.* 1960, 1, 516–523. [CrossRef]

47. Haldane, F.D.M. “Fractional statistics” in arbitrary dimensions: A generalization of the Pauli principle. *Phys. Rev. Lett.* 1991, 67, 937–940. [CrossRef] [PubMed]

48. Wu, Y.S. Statistical Distribution for Generalized Ideal Gas of Fractional-Statistics Particles. *Phys. Rev. Lett.* 1994, 73, 922–925. [CrossRef] [PubMed]

49. Murthy, M.V.N.; Shankar, R. Thermodynamics of a One-Dimensional Ideal Gas with Fractional Exclusion Statistics. *Phys. Rev. Lett.* 1994, 73, 3331–3334. [CrossRef]

50. Girardeau, M. Relationship between Systems of Impenetrable Bosons and Fermions in One Dimension. *J. Math. Phys.* 1960, 1, 516–523. [CrossRef] [PubMed]

51. Ares, F.; Gupta, K.S.; de Queiroz, A.R. Orthogonality catastrophe and fractional exclusion statistics. *Phys. Rev. E* 2018, 97, 022133. [CrossRef]

52. Vacek, K.; Okiji, A.; Kawakami, N. Eigenfunctions for SU( nu ) particles with 1/r2interaction in harmonic confinement. *J. Phys. A Math. Gen.* 1994, 27, L201–L206. [CrossRef]

53. Castin, Y.; Werner, F. The Unitary Gas and its Symmetry Properties. In *The BCS-BEC Crossover and the Unitary Fermi Gas*; Zwerger, W., Ed.; Springer: Berlin/Heidelberg, Germany, 2012; pp. 127–191. [CrossRef]

54. Tan, S. Short Range Scaling Laws of Quantum Gases With Contact Interactions. *arXiv* 2004, arXiv:cond-mat/0412764.
61. Werner, F.; Castin, Y. Unitary gas in an isotropic harmonic trap: Symmetry properties and applications. *Phys. Rev. A* **2006**, *74*, 053604. [CrossRef]

62. Rajabpour, M.A.; Sotiriadis, S. Quantum quench of the trap frequency in the harmonic Calogero model. *Phys. Rev. A* **2014**, *89*, 033620. [CrossRef]

63. Deffner, S.; Lutz, E. Nonequilibrium work distribution of a quantum harmonic oscillator. *Phys. Rev. E* **2008**, *77*, 021128. [CrossRef] [PubMed]

64. Campbell, S.; García-March, M.A.; Fogarty, T.; Busch, T. Quenching small quantum gases: Genesis of the orthogonality catastrophe. *Phys. Rev. A* **2014**, *89*, 033620. [CrossRef]

65. Deffner, S.; Lutz, E. Nonequilibrium thermodynamics of harmonically trapped bosons. *New J. Phys.* **2016**, *18*, 103035. [CrossRef]

66. Vicari, E. Particle-number scaling of the quantum work statistics and Loschmidt echo in Fermi gases with time-dependent traps. *Phys. Rev. A* **2019**, *99*, 043603. [CrossRef]

67. Rezek, Y.; Kosloff, R. Irreversible performance of a quantum harmonic heat engine. *New J. Phys.* **2006**, *8*, 83–83. [CrossRef]

68. Salamon, P.; Hoffmann, K.H.; Rezek, Y.; Kosloff, R. Maximum work in minimum time from a conservative quantum system. *Phys. Chem. Chem. Phys.* **2009**, *11*, 1027–1032. [CrossRef]

69. Rezek, Y.; Salamon, P.; Hoffmann, K.H.; Kosloff, R. The quantum refrigerator: The quest for absolute zero. *EPL Europhys. Lett.* **2009**, *85*, 30008. [CrossRef]

70. Hoffmann, K.H.; Salamon, P.; Rezek, Y.; Kosloff, R. Time-optimal controls for frictionless cooling in harmonic traps. *EPL Europhys. Lett.* **2011**, *96*, 60015. [CrossRef]

71. Abah, O.; Roßnagel, J.; Jacob, G.; Deffner, S.; Schmidt-Kaler, F.; Singer, K.; Lutz, E. Single-Ion Heat Engine at Maximum Power. *Phys. Rev. Lett.* **2012**, *109*, 203006. [CrossRef]

72. Kosloff, R.; Rezek, Y. The Quantum Harmonic Otto Cycle. *Entropy* **2017**, *19*, 136. [CrossRef]

73. Lidar, D.A.; Rezakhani, A.T.; Hamma, A. Adiabatic approximation with exponential accuracy for many-body systems and quantum computation. *J. Math. Phys.* **2009**, *50*, 102106. [CrossRef]

74. Kim, S.P.; Kim, W. Construction of exact Ermakov-Pinney solutions and time-dependent quantum oscillators. *J. Korean Phys. Soc.* **2016**, *69*, 1513–1517. [CrossRef]

75. Abah, O.; Lutz, E. Energy efficient quantum machines. *EPL Europhys. Lett.* **2017**, *118*, 40005. [CrossRef]

76. Stefanatos, D.; Ruths, J.; Li, J.S. Frictionless atom cooling in harmonic traps: A time-optimal approach. *Phys. Rev. A* **2010**, *82*, 063422. [CrossRef]

77. Stefanatos, D. Minimum-Time Transitions Between Thermal Equilibrium States of the Quantum Parametric Oscillator. *IEEE Trans. Autom. Control* **2017**, *62*, 4290–4297. [CrossRef]

78. Larocca, M.; Calzetta, E.; Wisniacki, D.A. Exploiting landscape geometry to enhance quantum optimal control. *Phys. Rev. A* **2020**, *101*, 023410. [CrossRef]

79. Fogarty, T.; Deffner, S.; Busch, T.; Campbell, S. Orthogonality Catastrophe as a Consequence of the Quantum Speed Limit. *Phys. Rev. Lett.* **2020**, *124*, 110601. [CrossRef]

© 2020 by the authors. Licensee MDPI, Basel, Switzerland. This article is an open access article distributed under the terms and conditions of the Creative Commons Attribution (CC BY) license (http://creativecommons.org/licenses/by/4.0/).