Mean-field theory of atomic self-organization in optical cavities

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Photons mediate long-range optomechanical forces between atoms in high-finesse resonators, which can induce the formation of ordered spatial patterns. When a transverse laser drives the atoms, the system undergoes a second-order phase transition that separates a uniform spatial density from a Bragg grating maximizing scattering into the cavity and is controlled by the laser intensity. Starting from a Fokker-Planck equation describing the semiclassical dynamics of the *N*-atom distribution function, we systematically develop a mean-field model and analyze its predictions for the equilibrium and out-of-equilibrium dynamics. The validity of the mean-field model is tested by comparison with the numerical simulations of the *N*-body Fokker-Planck equation and by means of a Bogoliubov-Born-Green-Kirkwood-Yvon (BBGKY) hierarchy. The mean-field theory predictions well reproduce several results of the *N*-body Fokker-Planck equation for sufficiently short times and are in good agreement with existing theoretical approaches based on field-theoretical models. The mean field, on the other hand, predicts thermalization time scales which are at least one order of magnitude shorter than the ones predicted by the *N*-body dynamics. We attribute this discrepancy to the fact that the mean-field ansatz discards the effects of the long-range incoherent forces due to cavity losses.

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I. INTRODUCTION

Optically dense atomic ensembles offer a formidable framework to study collective effects induced by atom-photon interactions [1-3]. Correlations are established by multiple-photon scattering [4,5], which can give rise to phenomena such as synchronization [6,7], optomechanical bistability [8,9], and spontaneous spatial ordering [2,3,10–12]. Envisaged applications for these systems range from sensors [13] to quantum-enhanced metrology [14] and quantum simulators [12,15].

Single-mode cavities, furthermore, mediate strong longrange interactions between the atoms [16–18], similar to gravitational and Coulomb potential in two or more dimensions [19]. In view of this analogy, it is relevant to study the dynamics of these systems at and out of equilibrium to test in a laboratory conjectures and predictions, such as ensemble inequivalence and the existence of quasistationary states [16]. The realization in quantum optical setups, like the one sketched in Fig. 1(a), brings additional peculiar features. In fact, these systems are intrinsically lossy, so that nontrivial dynamics can be observed only in the presence of a pump. On the one hand, the conservative potential mediated by the cavity photons shares several analogies with the one of the Hamiltonian Mean Field model [17,19-21], of which several features are well reproduced by a mean-field description [19,20]. On the other hand, cavity losses give rise to damping and diffusion, which are characterized by a spatial structure, thus establishing long-range correlations between the atoms [17,22]. These correlations, in turn, cannot be captured by a mean-field description.

In this work we systematically develop a mean-field model for cold atoms in a standing-wave resonator in the setup illustrated in Fig. 1(a) and test its validity by comparing its predictions with the ones of the Fokker-Planck equation for the *N*-atom distribution [22]. This work completes a series of papers which analyze the equilibrium and out-of-equilibrium dynamics of spatial self-organization of atomic ensembles in a single-mode resonator. Our analysis is based on a semiclassical treatment, specifically on a Fokker-Planck equation (FPE) for the *N*-atom distribution, derived when the atoms are classically polarizable particles and their center-of-mass motion is confined to one dimension [22]. The cavity field, instead, is a full quantum variable. This makes our treatment applicable also in the shot-noise limit [22] and gives access to regimes that are complementary to those based on the model in Ref. [23], where the field is a semiclassical variable.

Our formalism permits us to consistently eliminate the cavity variables from the equations of motion of the atoms and to investigate the properties of the cavity field across the self-organization threshold, where the intracavity field is characterized by large fluctuations. Starting from this model in Ref. [21] we analyzed the stationary state of the *N*-body FPE and showed that (i) this is a thermal state whose temperature is determined by the linewidth of the resonator and (ii) the transition to self-organization is a Landau-type second-order phase transition, as illustrated in Figs. 1(b) and 1(c). In Ref. [21] we also determined the corresponding phase diagram as a function of the physical parameters and predicted the corresponding features in the light emitted by the resonator. In Ref. [24] we investigated the dynamics following sudden quenches across the phase transition and found that the interplay between long-range conservative and dissipative forces gives rise to prethermalization dynamics, where the long-range nature of dissipation plays an essential role.

In this work we derive a mean-field treatment from our N-atom FPE. We then benchmark the limits of validity of the mean-field ansatz by means of numerical simulations using the full N-body FPE and by means of a Bogoliubov-Born-Green-Kirkwood-Yvon (BBGKY) hierarchy. The results we obtain are compared with existing literature on spatial self-organization in single-mode cavities for both the semiclassical treatment [25–28] and the case in which the atomic quantum statistics is assumed to be relevant [15,28–33].

This work is organized as follows. In Sec. II the Fokker-Planck equation at the basis of our analysis is reported, and the corresponding mean-field equation is derived. In



FIG. 1. (a) Atoms in a standing-wave cavity and driven by a transverse laser can spontaneously form ordered patterns when the amplitude of the laser coupling Ω exceeds a threshold value Ω_c , which depends on the rate of photon losses, here due to cavity decay at rate κ . In this regime the system undergoes a second-order phase transition which is characterized by the parameter Θ , indicating spatial ordering of the atoms in Bragg gratings and defined in Eq. (4). Its expectation value in the mean-field description is denoted as Θ_{MF} . (b) and (c) The thermodynamic potential below and above threshold. The bottom panels are schematic pictures of the single-particle density distribution $f_1(x, p)$ in phase space with x in units of the inverse wave number k^{-1} and p in units of the width Δp of the momentum distribution. In (b) the atomic density is uniform; in (c) it is localized at the even or odd sites of the cavity standing wave, $[\cos(kx) = 1 \text{ or }$ -1, respectively]. In this work we derive and discuss a mean-field theory for the dynamics of $f_1(x, p)$.

Sec. III the stationary properties of the mean-field FPE distribution function are analytically determined. The mean-field predictions are compared with the ones of the *N*-body FPE and with further existing theoretical works. In Sec. IV the Vlasov equation, which describes the short-time dynamics of the mean-field FPE, is derived. Its predictions are then determined by means of a stability analysis, and the analytical results are compared with the numerical simulations of the mean-field FPE. Section V reports a critical analysis of the limits of validity of the mean-field treatment. In Sec. VI the conclusions are drawn, while in the Appendix calculations are reported that complement the material presented in Sec. III.

II. DERIVATION OF THE MEAN-FIELD MODEL

In this section we derive the mean-field model starting from the FPE describing the dynamics of an atomic ensemble in the optical potential of a high-finesse resonator of Ref. [22]. The atoms are N and have mass m, and their motion is assumed to be confined along the x axis, which also coincides with the axis of a high-finesse cavity within whose mirrors the atoms are spatially trapped. In the following we denote their canonically conjugated positions and momenta by x_j and p_j (j = 1, ..., N). The atomic dipole strongly couples to one cavity mode and is transversally driven by a laser, as sketched in Fig. 1(a). The parameter regime is such that the atoms coherently scatter photons into the cavity mode, and their external motion is determined by the light forces associated with these processes. The light forces are periodic, and their period is determined by the cavity-mode standing wave, whose spatial mode function is $\cos(kx)$, with k being the cavity-mode wave number.

A. Basic assumptions

Before reporting the FPE which governs the dynamics of the *N*-body distribution function, we summarize the main approximations behind its derivation and the corresponding physical parameters.

One basic assumption of our model is that the only relevant scattering processes are coherent. This regime can be reached when the cavity mode and laser frequencies are tuned far off resonance from the atomic transition [34,35]. We denote by $\Delta_a = \omega_L - \omega_0$ the detuning between laser (ω_L) and atomic (ω_0) frequencies and assume that this is the largest parameter of the problem. It is thus larger than the coupling strengths for the interaction between the dipole and fields. It is also larger than the detuning $\Delta_c = \omega_L - \omega_c$ between laser and cavity-mode frequencies, whose wave numbers are, to good approximation, denoted by the same parameter k. This allows us to eliminate the internal degrees of freedom of the atoms by a perturbative expansion in the lowest order of the small parameter $1/|\Delta_a|$.

The cavity field is treated as a quantum-mechanical variable, and the dynamics can be cast as an optomechanical coupling between atomic motion and cavity field [10,23]. The parameter regime we assume gives rise to a time-scale separation, such that the cavity degrees of freedom evolve on a faster time scale than the motion. This is warranted when the cavity linewidth κ , which determines the relaxation rate of the resonator state, is much larger than the recoil frequency $\omega_r = \hbar k^2/(2m)$, which scales the exchange of mechanical energy between light and atoms. In this limit the cavity field is eliminated from the equations of motion of the atomic external degrees of freedom in a perturbative expansion to first order in the small parameter $1/\kappa$, implementing a procedure first applied in Ref. [36]. The hierarchy of time scales is set by the inequalities $|\Delta_a| \gg \kappa \gg \omega_r$. This is also consistent with a semiclassical treatment since the kinetic energy of the atoms at steady state scales with $\hbar \kappa$, thus warranting that the width Δp of the single-atom momentum distribution is large in comparison to the linear momentum $\hbar k$ carried by each photon [10,17,21].

B. Collective motion of N atoms in a cavity field

The approximations discussed above are at the basis of the theoretical procedure which connects the master equation of atoms in a quantized cavity field with the FPE for the Wigner function $f_N = f_N(x_1, ..., x_N; p_1, ..., p_N; t)$ describing the

positions and momenta of the N atoms at time t. The derivation is detailed in Ref. [22], and the resulting FPE reads

$$\frac{\partial f_N}{\partial t} = -\sum_{i=1}^N \frac{\partial}{\partial x_i} \frac{p_i}{m} f_N + S^2 L[f_N], \qquad (1)$$

where the second summand on the right-hand side (RHS) is due to mechanical effects of the cavity field on the atoms and scales like S^2 . Here, $S = \Omega g / \Delta_a$ is the scattering amplitude between the laser and cavity mode, and it is proportional to the laser strength Ω and to the cavity vacuum Rabi frequency g, which scale the interaction between the dipole and laser and between the dipole and cavity, respectively. Operator $L[f_N]$ takes the form

$$S^{2}L[f_{N}] = \frac{\partial f_{N}}{\partial p_{i}} \frac{\partial V(x_{1}, \dots, x_{N})}{\partial x_{i}}$$
(2a)

$$-S^{2}\sum_{i,j}\frac{\partial}{\partial p_{i}}\Gamma_{0}\sin(kx_{i})\sin(kx_{j})p_{j}f_{N}$$
(2b)

+
$$S^2 \sum_{i,j} \frac{\partial^2}{\partial p_i \partial p_j} D_0 \sin(kx_i) \sin(kx_j) f_N$$
 (2c)

+
$$S^2 \sum_{i,j} \frac{\partial^2}{\partial p_j \partial x_i} \eta_0 \sin(kx_i) \sin(kx_j) f_N$$
. (2d)

Each line on the RHS of Eq. (2) has a physical meaning. The first term describes the dynamics due to the conservative potential

$$V(x_1,\ldots,x_N) = \frac{\hbar\Delta_c}{\kappa^2 + \Delta_c^2} S^2 N^2 \Theta(x_1,\ldots,x_N)^2, \quad (3)$$

where

$$\Theta(x_1,\ldots,x_N) = \frac{1}{N} \sum_{j=1}^N \cos(kx_j), \qquad (4)$$

so that the potential mediates long-range interactions between the atoms. Parameter $\langle |\Theta| \rangle_N$ is the order parameter of self-organization, where $\langle \cdot \rangle_N$ denotes the expectation value taken over the normalized distribution f_N . Specifically, when the atoms form Bragg grating, then $\langle |\Theta| \rangle_N \rightarrow 1$, and the potential depth is maximal. When the atoms are instead uniformly distributed in space, then $\langle |\Theta| \rangle_N \simeq 0$, and the potential vanishes. We note that the Bragg gratings minimize the potential when $\Delta_c < 0$; otherwise, the uniform distribution is energetically favored. We will here denote $\langle |\Theta| \rangle_N$ by magnetization due to the mapping of the self-organization transition to a ferromagnetic model [21].

For later convenience, we define the parameter

$$F_0 = (\hbar k) \frac{2\Delta_c}{\kappa^2 + \Delta_c^2},\tag{5}$$

such that $V = F_0(NS\Theta)^2/(2k)$.

The second term on the RHS [Eq. (2b)] describes a dissipative force and is scaled by the coefficient Γ_0 :

$$\Gamma_0 = \omega_r \frac{8\Delta_c \kappa}{\left(\kappa^2 + \Delta_c^2\right)^2} \,. \tag{6}$$

This term is due to nonadiabatic corrections in the dynamics of the cavity field.

The term in (2c) corresponds to diffusion due to fluctuations of the cavity field associated with losses. The diffusion matrix is the dyadic product of the vector $(\sin(kx_1), \ldots, \sin(kx_N))$ with itself and scales with the coefficient

$$D_0 = (\hbar k)^2 \frac{\kappa}{\kappa^2 + \Delta_c^2}.$$
(7)

Therefore, besides the diffusion due to the diagonal elements, which is a single-particle effect, we also expect that term (2c) establishes long-range correlations.

The last line, (2d), contains cross derivatives and scales with the coefficient

$$\eta_0 = 2\hbar\omega_r \frac{\kappa^2 - \Delta_c^2}{\left(\kappa^2 + \Delta_c^2\right)^2},\tag{8}$$

whose sign depends on whether the ratio $|\Delta_c/\kappa|$ is smaller or larger than unity, while it vanishes for $|\Delta_c/\kappa| = 1$. An analogous term has also been reported in the semiclassical description of cold atoms in optical lattices [36], where it has been neglected under the assumption of uniform spatial densities. Such an assumption cannot be applied in the selforganized regime; nevertheless, we will show that this term can be consistently discarded in the thermodynamic limit we apply, which warrants Kac's scaling [19].

C. Mean-field ansatz

To derive a mean-field FPE we assume that the Wigner function is factorized into single-particle distribution functions according to the prescription

$$f_N(x_1, \dots, x_N; p_1, \dots, p_N; t) = \prod_{i=1}^N f_1(x_i, p_i; t),$$
 (9)

where $f_1(x_i, p_i; t)$ denotes the distribution for particle *i* at time *t* and is thus defined on the phase space of this particle. We use then Eq. (9) in the FPE (1) and integrate out all particles' variables but one. In this way we derive the mean-field FPE, which reads

$$\frac{\partial f_1}{\partial t} = -\frac{\partial}{\partial x} \frac{p}{m} f_1 + S^2 \mathfrak{L}[f_1]$$
(10)

and has same structure as the FPE in Eq. (1). Operator \mathfrak{L} describes, as *L*, the mechanical effects of light. However, it is now a nonlinear operator of f_1 and takes the form

$$\mathfrak{L}[f_1] = -\frac{\partial}{\partial p} F_0\{\cos(kx) + (N-1)\Theta_{\rm MF}[f_1]\}\sin(kx)f_1$$
(11a)

$$-\frac{\partial}{\partial p}\Gamma_0\{\sin(kx)p + (N-1)\Xi_{\rm MF}[f_1]\}\sin(kx)f_1$$

(11b)

$$+\frac{\partial^2}{\partial^2 p} D_0 \sin^2(kx) f_1 \tag{11c}$$

$$+\frac{\partial^2}{\partial p \partial x}\eta_0 \sin^2(kx)f_1, \qquad (11d)$$

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where we have introduced the functionals

$$\Theta_{\rm MF}[f_1] = \frac{1}{\lambda} \int_0^\lambda dx \int_{-\infty}^\infty dp \cos(kx) f_1, \qquad (12)$$

$$\Xi_{\rm MF}[f_1] = \frac{1}{\lambda} \int_0^\lambda dx \int_{-\infty}^\infty dp \sin(kx) p f_1.$$
(13)

The mean-field order parameter Θ_{MF} is the expectation value $\langle \cos(kx) \rangle$, where $\langle \cdot \rangle$ indicates the average taken over the single-particle distribution function $f_1(x, p)$. The terms on the RHS in (11a) and (11b) have a different origin but a similar structure, which can be recognized by analyzing the form of the two summands within the respective inner brackets. The first summand in each line describes the interaction of the atom with itself, mediated by the cavity field. The second summand in each line emerges from the interaction between the atom and all other N - 1 atoms.

We further notice that the term in (11a) can be cast in terms of a conservative force originating from the potential

$$V_{1}[f_{1}](x) = \frac{F_{0}}{2k} S^{2}[\cos^{2}(kx) + 2(N-1)\Theta_{\rm MF}[f_{1}]\cos(kx)] + \frac{\Gamma_{0}}{k} (N-1)S^{2}\Xi_{\rm MF}[f_{1}]\cos(kx)$$
(14)

and contains a term whose corresponding term in Eq. (1) has dissipative nature [see (2b)]. Using this result, we can rewrite Eq. (11) in the compact form

$$\mathfrak{L}[f_1] = \frac{\partial V_1}{\partial x} \frac{\partial f_1}{\partial p} - \frac{\partial}{\partial p} \left(\Gamma_0 p - \frac{\partial}{\partial p} D_0 - \frac{\partial}{\partial x} \eta_0 \right) \sin^2(kx) f_1,$$

which allows us to simply read out the physical meaning of the other terms; they are, in fact, the diagonal component of friction, diffusion, and the cross-derivative term in Eq. (1).

III. STATIONARY STATE OF THE MEAN-FIELD EQUATION

The stationary properties of the mean-field distribution are analyzed by means of the single-particle distribution $f_{st}(x, p)$ that solves Eq. (10) with

$$\partial_t f_{\rm st}(x,p) = 0. \tag{15}$$

In the following we determine $f_{st}(x,p)$ and then analyze its predictions for relevant physical quantities.

A. Derivation of the steady-state solution

In order to solve Eq. (15) we consider the ansatz

$$f_{\rm st}(x,p) = f_0 \exp[a(x) + b(p)],$$

where a(x) and b(p) are functions which depend only on position and momentum, respectively, and f_0 is the normalization constant. Using this ansatz in Eq. (10), we obtain differential equations for a(x) and b(p), whose solutions read $b(p) = -\beta p^2/(2m)$ and

$$a(x) = (Y/2 - 1)\ln[1 + Z\sin^2(kx)] - (N - 1)Y\Theta_{\rm MF}[f_{\rm st}]$$
$$\times \sqrt{\frac{Z}{1+Z}} \operatorname{arctanh}\left(\sqrt{\frac{Z}{1+Z}}\cos(kx)\right), \tag{16}$$

with $Y = F_0/(k\eta_0)$, $Z = \beta \eta_0 S^2$, and

$$\beta = -\frac{\Gamma_0 m}{D_0} = \frac{-4\Delta_c}{\hbar \left(\kappa^2 + \Delta_c^2\right)}.$$
(17)

Therefore,

$$f_{\rm st}(x,p) = \mathcal{F}(\cos kx) \exp\left(-\beta \frac{p^2}{2m}\right),\tag{18}$$

with $\mathcal{F}(\cos kx) = f_0 \exp[a(x)]$. Equation (18) describes a thermal distribution provided that $\Delta_c < 0$: In this limit parameter β , Eq. (17), plays the role of an inverse temperature at steady state. This temperature coincides with the value found by solving the steady state of the *N*-body FPE, Eq. (1), as shown in Refs. [17,21].

We note that the function $\mathcal{F}(\cos kx)$ depends on $\Theta_{\text{MF}}[f_{\text{st}}]$, which leads to the fixed-point equation

$$\Theta_{\rm MF} \equiv \langle \cos(kx) \rangle = \sqrt{\frac{2\pi m}{\beta}} \frac{1}{\lambda} \int_0^\lambda dx \cos(kx) \mathcal{F}(\cos kx).$$
(19)

Its solution is, in general, not transparent, but it gets simpler in an appropriately defined thermodynamic limit. This consists of scaling the coupling strength $g \sim 1/\sqrt{N}$ as the number of atoms is increased, leading to the scaling relation $S \propto 1/\sqrt{N}$ [37,38]. In this limit function a(x), Eq. (16), can be cast into the form

$$a(x) = 2\frac{\bar{n}}{\bar{n}_c}\Theta_{\rm MF}\cos(kx),\tag{20}$$

with

$$\bar{n} = \frac{NS^2}{\kappa^2 + \Delta_c^2} \tag{21}$$

and

$$\bar{n}_c = \frac{\kappa^2 + \Delta_c^2}{4\Delta_c^2}.$$
(22)

This leads to a compact form of the stationary distribution in the mean-field limit:

$$f_{\rm st}(x,p) = f_0 \exp\left[-\beta \left(\frac{p^2}{2m} + \hbar \Delta_c \bar{n} \Theta_{\rm MF} \cos(kx)\right)\right], \quad (23)$$

with

$$f_0^{-1} = \sqrt{\frac{2m\pi}{\beta}} I_0 \left(2\frac{\bar{n}}{\bar{n}_c} \Theta_{\rm MF} \right),$$

where I_j is the modified Bessel function of *j*th order [39].

We thus see that in the thermodynamic limit the effect of the cross derivatives vanishes. For finite N, parameter η_0 is small but finite, and in the stationary state it gives rise to a correction to the effective potential term, as seen in Eq. (16).

B. Stationary properties in the thermodynamic limit

The mean-field distribution, Eq. (23), allows one to analytically determine several properties of the steady state. First, functional Θ_{MF} in the exponent has to be determined



FIG. 2. Plot of $q(2\bar{n}\zeta/\bar{n}_c)$, Eq. (25), as a function of ζ and for different values of \bar{n} . The intersection points with the curve $y = \zeta$ (dashed line) give the solutions of Eq. (24). Stable points are at the crossing where $q' < \bar{n}_c/(2\bar{n})$ and are the equilibrium values of the order parameter $\Theta_{\rm MF}$. Inset: The resulting stable solution $\bar{\Theta} \ge 0$ as a function of \bar{n} (in units of \bar{n}_c).

self-consistently. Using Eq. (20) in Eq. (19) gives the relation

$$\Theta_{\rm MF} = q \left(2 \frac{\bar{n}}{\bar{n}_c} \Theta_{\rm MF} \right), \tag{24}$$

where q is a function of the form

$$q\left(2\frac{\bar{n}}{\bar{n}_c}\zeta\right) = \frac{I_1\left(2\frac{\bar{n}_c}{\bar{n}_c}\zeta\right)}{I_0\left(2\frac{\bar{n}}{\bar{n}_c}\zeta\right)} \tag{25}$$

and is plotted in Fig. 2 for values of \bar{n} below, at, and above \bar{n}_c . The solutions of Eq. (24) are the crossing between the curve $y = \zeta$ and $y = q(2\bar{n}\zeta/\bar{n}_c)$ [see Eq. (25)]. For $\bar{n} < \bar{n}_c$ this equation allows for one solution, corresponding to $\Theta_{MF} = 0$. For $\bar{n} > \bar{n}_c$, there are three solutions, two of which are stable and one of which is unstable. The stable solutions give $\Theta_{MF} = \pm \bar{\Theta}$, with $0 \leq \bar{\Theta} < 1$, and correspond to the self-organized state. Close to, but above, the critical point the value $\bar{\Theta}$ can be analytically determined and reads

$$\bar{\Theta} = \sqrt{2(\bar{n}/\bar{n}_c - 1)}.$$
(26)

The value $\bar{n} = \bar{n}_c$, with \bar{n}_c defined in Eq. (22), hence determines a critical point at which the transition to selforganization occurs and that is controlled by the detuning from the cavity field and the cavity loss rate for the thermodynamic limit we chose. The results we have obtained so far for the stationary mean-field distribution are in full agreement with the ones found for the stationary distribution of Eq. (1) (see Ref. [21]). The stationary mean-field distribution in Eq. (23) corresponds to the one that is found from the stationary *N*-particle distribution after integrating out the other N - 1 position and momentum variables and then taking the thermodynamic limit. The equation for the order parameter, Eq. (24), agrees with the one obtained for the *N*-particle case and obtained by means of a saddle-point approximation. This agreement is found also for the critical value of Eq. (22) and for the temperature of Eq. (17). Hence, the mean-field model predicts the same phase diagram as the *N*-body FPE.

It is also instructive to consider the value of the bunching parameter \mathcal{B} as a function of \bar{n} . This is defined as

$$\mathcal{B} = \langle \cos^2(kx) \rangle \tag{27}$$

and gives a measure of localization of the particles at the minima of the mechanical potential [21,25]. Using Eq. (20), we obtain

$$\mathcal{B} = \begin{cases} 1/2, & \bar{n} \le \bar{n}_c, \\ 1 - \bar{n}_c/(2\bar{n}), & \bar{n} > \bar{n}_c, \end{cases}$$
(28)

in the stationary state. Therefore, below threshold the atoms are uniformly distributed, while above threshold they increasingly localize at the minima of the Bragg potential. In particular, when the atoms are tightly bound at the minima, the abovethreshold expression in Eq. (28) delivers the amplitude of the fluctuations, namely,

$$k^2 \langle x^2 \rangle \approx \frac{\bar{n}_c}{2\bar{n}},$$
 (29)

showing that these are inversely proportional to the laser intensity.

C. Comparison with existing literature

The results obtained so far by means of the mean-field model show a remarkable agreement with the predictions of the stationary solution of the N-particle FPE, Eq. (1). It is further worthwhile to compare the results derived here with the results obtained in the literature by means of different approaches.

We first discuss Ref. [25], where, among other studies, a mean-field approach is developed based on plausible conjectures. Here, the mean-field potential is calculated, and the threshold of self-organization is determined by (i) assuming that the stationary state is thermal, with temperature given by the linewidth of the cavity, and (ii) performing a stability analysis of the uniform density distribution. By means of this study a threshold value for self-organization is identified, which agrees with the prediction in Eq. (22), which becomes evident after defining the threshold amplitude S_c such that

$$\frac{NS_c^2}{\Delta_c^2 + \kappa^2} \equiv \bar{n}_c.$$

In particular, the quantity η^* in [25] is in our notations $S_c \Delta_a/g$ calculated for the case $\Delta_c = -\kappa$.

The stationary state of self-organization was first derived in Refs. [26,27] by means of a FPE as a function of the atomic and field variables. This description assumes that the field fluctuations are small and thus cannot reliably reproduce the field correlation functions below and at threshold. It predicts, nevertheless, that the atoms' steady state is thermal and its temperature coincides with the inverse of Eq. (17), apart from corrections of the order ω_r/κ that are systematically neglected in our approach because they are of higher order. It further predicts the same behavior of the order parameter as in Eq. (26) above, but close to, the threshold.

It is also interesting to compare our results with a series of other theoretical studies which focus on self-organization of ultracold atomic ensembles in cavities but discard retardation effects: In these works only the conservative part of the cavity potential is considered, while the temperature at steady state is due to the coupling to an external heat bath [15,28–33]. Even though the conditions seem quite different from our case, remarkable agreement is found in the appropriate limits. References [15,29] analyze the self-organization transition of an ultracold gas of bosonic atoms and derive the mapping to the Dicke model. Here, the recoil energy plays a role analogous to the temperature, and the threshold which is derived agrees with the threshold in Eq. (22) after setting

$$NS_c^2 = \frac{1}{\beta} \frac{\kappa^2 + \Delta_c^2}{-\Delta_c},\tag{30}$$

with $\beta = 4/\hbar\omega_r$. By means of this prescription, the threshold also agrees with the one calculated in Ref. [28]. Furthermore, it also coincides with the one evaluated in Ref. [33] when using the Boltzmann distribution for the atoms statistics.

Another quantity which has been determined in these works is the photon flux, which corresponds to the intracavity photon number in our treatment. In Refs. [28–31] the photon flux scales as $1/|\bar{n} - \bar{n}_c|$ below threshold, while at threshold it diverges as \sqrt{N} . These predictions are in perfect agreement with the results we find when taking the stationary distribution of Eq. (1) [see Eqs. (A4) and (A5)]. In particular, the intracavity photon number at threshold, Eq. (A5), coincides with the one calculated in Ref. [28] after substituting in their equation $\omega_z = (\omega_0^2 + \kappa^2)/\omega_0$ for the temperature, with $\omega_0 = -\Delta_c$. The result for the intensity-intensity correlations at zero time delay and below threshold, Eq. (A11), further agrees with the result derived in Refs. [30,31].

IV. MEAN-FIELD DYNAMICS

We now study the dynamics predicted by the mean-field FPE. We focus on the Vlasov equation, which we derive from Eq. (10) by taking the thermodynamic limit, according to our prescription. The Vlasov equation for our problem reads

$$\frac{\partial f_1}{\partial t} + \frac{p}{m} \frac{\partial f_1}{\partial x_1} - \frac{\partial V_0[f_1](x)}{\partial x} \frac{\partial f_1}{\partial p} = 0, \qquad (31)$$

with

$$V_0[f_1](x) = 2\hbar\Delta_c\bar{n}\cos(kx)\Theta_{\rm MF}[f_1] - \frac{\hbar^2k}{m}\bar{n}\beta\kappa\cos(kx)\Xi_{\rm MF}[f_1], \qquad (32)$$

and it corresponds to the potential in Eq. (14) after neglecting the self-reaction term, which is of order 1/N. Therefore, the validity of the predictions we will extract is limited to sufficiently short time scales for which the corrections can be discarded. We will quantify this statement in the next section.

A. Preliminary considerations: Energy conservation

We first analyze whether Eq. (31) warrants energy conservation. We consider a class of functions for which $\Xi_{MF}[f_1] = 0$. This includes the stationary solution of Eq. (23). For these solutions, the energy of one particle takes the form

$$\epsilon(t) = \frac{\langle p^2 \rangle}{2m} + \hbar \Delta_c \bar{n} \Theta_{\rm MF}^2.$$
(33)

In order to determine $\dot{\epsilon}(t)$ we thus calculate $\dot{\Theta}_{\rm MF}$ and $\langle p^2 \rangle$. This gives

$$\dot{\Theta}_{\rm MF} = -\frac{k}{m} \Xi_{\rm MF},$$
$$\frac{\langle \dot{p}^2 \rangle}{2m} = 2 \frac{\hbar}{m} \bar{n} (k \Delta_c \Theta_{\rm MF} - \omega_r \kappa \beta \Xi_{\rm MF}) \Xi_{\rm MF}$$

and therefore, we get for the derivative of the energy

$$\dot{\epsilon} = -2\frac{\hbar}{m}\bar{n}\omega_r\kappa\beta\Xi_{\rm MF}^2.$$

These derivatives hence vanish when $\Xi_{MF} = 0$, and thus, for the class of distribution fulfilling this condition, energy, with the potential term given in Eq. (33), is conserved. Fluctuations, on the other hand, can give rise to finite values of Ξ_{MF} . The purpose of the next section is to analyze the stability and shorttime dynamics of solutions of the Vlasov equation, Eq. (31), after quenches of the laser parameters.

B. Stability analysis of spatially homogeneous distributions

We now analyze the short-time dynamics described by Eq. (31), assuming that at t = 0 the distribution is thermal and has uniform spatial density; thus, $f_1(x, p, 0) = f_1(p, 0)$ and $\Theta_{\text{MF}}|_{t=0^-} = 0$, with

$$f_1(p,0) = \left(\frac{2m\pi}{\beta_0}\right)^{-\frac{1}{2}} \exp\left(-\beta_0 \frac{p^2}{2m}\right),$$
 (34)

where β_0 is the inverse temperature. This distribution is a stable solution of the Vlasov equation after setting $\bar{n} = 0$. At t = 0 the laser strength is quenched above threshold, so that parameter \bar{n} takes a finite value larger than \bar{n}_c . We then let evolve the distribution of Eq. (34) by taking this value \bar{n} in Eq. (31). Figure 3 shows the results of the numerical integration of Eq. (31) for different values of \bar{n} . We analyze these results, keeping in mind that they are strictly valid for short times since the Vlasov equation discards effects, such as diffusion, which are crucial in determining the stationary state. In Fig. 3(a) the order parameter evolves from zero to a finite value, about which it oscillates. This value is smaller than the one predicted by the stationary solution of the mean-field FPE. It is reached after an initial dynamics characterized by an exponential increase, whose slope is steeper the larger \bar{n} is. Figures 3(b) and 3(c) display the corresponding evolution of the quantities Ξ_{MF}^2 [see Eq. (13)]. This quantity emerges from the retardation effects of the dynamics; it is thus a signature of memory effects and mathematically corresponds to the buildup correlations between momentum and position that cannot be factorized. The initial distribution, Eq. (34), is chosen so that $\Xi_{MF} = 0$, and we observe that the dynamics give rise to a buildup of a finite value of Ξ^2_{MF} , with an exponential increase that leads to a maximum where the curve for Θ_{MF} reaches the plateau. Then, it oscillates like Θ_{MF} (one can well understand the behavior of these oscillations observing that Ξ_{MF} is proportional to the time derivative of Θ_{MF}) and is exponentially damped to zero. In the initial phase, the exponential growth of Ξ_{MF}^2 increases with \bar{n} ; as in the second phase of the dynamics, where Θ_{MF} oscillates about a finite mean value, the amplitude of the oscillations of Ξ_{MF}^2 is also larger the larger \bar{n} is.



FIG. 3. Time evolution of (a) the order parameter Θ_{MF} , Eq. (12), and (b) and (c) parameter Ξ_{MF}^2 , Eq. (13), calculated by numerical integration of the Vlasov equation (31) for different values of \bar{n} and for $\Delta_c = -\kappa$. The initial distribution is given in Eq. (34) with $\beta_0 = 2/(\hbar\kappa)$.

We now analyze the initial exponential increase, which is in the regime where the Vlasov equation is a reliable approximation of the full dynamics, as we also verified in Ref. [24]. In order to do so, we use a standard procedure, which is also detailed in Refs. [19,40]. For short times t after the quench, we write the distribution as

$$f_1(x, p, t) = f_1(p, 0) + \delta f_1(x, p, t), \tag{35}$$

where δf_1 describes small fluctuations which can be due to the finite size of the system and thus scale with $1/\sqrt{N}$. Using Eq. (35) in the Vlasov equation (31) and neglecting the terms of order 1/N, we obtain the linearized Vlasov equation

$$\frac{\partial \delta f_1}{\partial t} + \frac{p}{m} \frac{\partial \delta f_1}{\partial x} - \frac{\partial \delta V}{\partial x} \frac{\partial f_1(p,0)}{\partial p} = 0, \qquad (36)$$

where $\delta V = V[\delta f_1(x, p, t)]$ and we dropped the argument of function δf_1 . We look for solutions of Eq. (36) by means of the ansatz of Fourier waves with frequency ω and wave number k:

$$\delta f_1 = g_1(p)e^{i(\omega t - kx)} + g_{-1}(p)e^{i(\omega t + kx)}, \qquad (37)$$

$$\delta V = 2A\cos(kx)e^{i\omega t},\tag{38}$$

where $A \propto 1/\sqrt{N}$ is some constant and the amplitudes $g_1(p)$ and $g_{-1}(p)$ are sole functions of the momentum p. The dispersion relation $\omega = \omega(k)$ can be derived after using Eqs. (37) and (38) in the linearized Vlasov equation (36). By equating the coefficients of exp(ikx) and exp(-ikx) we get expressions for the functions $g_1(p)$ and $g_{-1}(p)$. With those expressions one finds the dispersion relation by using the definition $\delta V = V[\delta f_1]$ and Eqs. (32) and (38):

1

$$0 = 1 + \left(\hbar\Delta_c + i\frac{\hbar\kappa}{2}\hbar\omega\beta\right)\bar{n}\frac{1}{2}$$
$$\times \int_{-\infty}^{\infty} dp \left(\frac{-k}{\frac{pk}{m} + \omega} + \frac{-k}{\frac{pk}{m} - \omega}\right)\partial_p f_1(p,0). \quad (39)$$

This relation holds for any initial distribution that describes a uniform spatial density. We now use the Gaussian distribution in Eq. (34) and obtain

$$0 = 1 + \left(\hbar\Delta_c + i\frac{\hbar\kappa}{2}\hbar\omega\beta\right)\bar{n}\beta_0$$

$$\times \left[1 - \bar{a}\exp(-\bar{a}^2)\left(i\sqrt{\pi} - 2\int_0^{\bar{a}}du\exp(u^2)\right)\right], (40)$$

where we defined $\bar{a} = \sqrt{\beta_0/(2m)}(m\omega/k)$. We then introduce $\bar{b} = i\bar{a}$ and

 $\gamma = i\omega$

and cast Eq. (40) in the form

$$0 = 1 + \left(\hbar\Delta_c + \frac{\hbar\kappa}{2}\hbar\gamma\beta\right)\bar{n}\beta_0$$

$$\times \left[1 - \bar{b}\exp(\bar{b}^2)\left(\sqrt{\pi} - \int_{-\bar{b}}^{\bar{b}} du\exp(-u^2)\right)\right], \quad (41)$$

where $\bar{b} \propto \gamma$. It can be shown that parameter γ , which solves Eq. (41), is a real number. Therefore, ω is an imaginary number. In particular, if $\gamma < 0$, both Eqs. (37) and (38) describe fluctuations which are exponentially damped, and therefore, $f_1(x, p, t)$ will tend to the initial distribution, which is stable. If, instead, the solution of Eq. (40) gives $\gamma > 0$, the initial distribution is unstable against fluctuations. The value $\gamma = 0$ separates the two regimes. After setting $\gamma = 0$ in Eq. (41), we thus get the critical condition

$$1 = -\hbar \Delta_c \bar{n} \beta_0, \tag{42}$$

which connects Δ_c , \bar{n} , and the initial temperature $1/\beta_0$, which is an external parameter. If β_0 coincides with the value in Eq. (17), then Eq. (42) corresponds to the same relation as in Eq. (22), which defines the critical value of \bar{n} for selforganization. For the values of the parameters for which $\gamma > 0$, the uniform distribution is unstable and tends to form a grating at the wave vector k of the resonator with exponential increase, giving rise to a violent relaxation. Parameter γ gives the rate at which the amplitude of this density modulation grows.

Figure 4 compares the value of γ extracted by fitting the exponential increase of Θ_{MF} in the first phase of the dynamics of Fig. 3 and for different values of \bar{n} with the one determined by Eq. (41), showing very good agreement. In particular, we note that in the limit $|\Delta_c| \gg |\gamma|$, Eq. (41) can be reduced to the form [41]

$$\gamma = \omega_0 (1 - p\chi) \frac{\ln\left(\frac{\chi}{1.135}\right) - \ln(1 - p\chi)}{1.4(1 - p\chi) + \hbar\kappa\beta\omega_0/(2|\Delta_c|)}, \quad (43)$$

with $\chi = \hbar |\Delta_c| \bar{n} \beta_0 = (\bar{n}/\bar{n}_c)(\beta_0/\beta), \, \omega_0 = \sqrt{2\omega_r/(\hbar\beta_0)}$, and p = 27/227.



FIG. 4. Slope γ of the initial increase of Θ_{MF} . The dots are extracted by fitting the curve obtained from the numerical simulations in Fig. 3, the dashed line is the value predicted by Eq. (43), which agrees well with Eq. (41) (solid line). For these parameters the threshold for the Vlasov stability, Eq. (42), reads $1 = \bar{n}/\bar{n}_c$.

V. VALIDITY OF THE MEAN-FIELD ANSATZ

The mean-field treatment is based on the assumption that the distribution function for the N particles can be approximated by the product of the single-particle distribution. This ansatz thus discards interparticle correlations which emerge from the photon-mediated interactions: the factorized ansatz is very different from the form of the distributions one obtains from the full N-particle FPE [17,21]. Nevertheless, the assumption still captures essential features of the shorttime dynamics of distributions, which initially have the form of Eq. (9). We will follow the procedure illustrated in Refs. [19,42] and study the validity of the mean-field ansatz within a BBGKY hierarchy, which we derive from the N-particle FPE, Eq. (1). We will particularly focus on the dynamics of two-particle correlations and determine the characteristic time scale of their dynamics.

For convenience, we introduce the vectors $\mathbf{x} = (x_1, \dots, x_N)^T$ and $\mathbf{p} = (p_1, \dots, p_N)^T$, and define $f_N(\mathbf{x}; \mathbf{p}; t) \equiv f_N(x_1, \dots, x_N; p_1, \dots, p_N; t)$.

A. BBGKY hierarchy of the photon-mediated Fokker-Planck equation

For the derivation of the BBGKY hierarchy we assume that the energy of the system is finite. This corresponds to assuming that the following limit holds,

$$\lim_{\mathbf{p}|\to\infty} f_N(\mathbf{x};\mathbf{p};t) = 0, \tag{44}$$

where $|\mathbf{p}| = \sqrt{\sum_{i=1}^{N} p_i^2}$, and that expectation values of all moments exist. Furthermore, f_N is periodic with wavelength λ in every x_i , which implies

$$f_N(\mathbf{x} + \lambda \mathbf{z}; \mathbf{p}; t) = f_N(\mathbf{x}; \mathbf{p}; t)$$
(45)

for every $\mathbf{z} \in \mathbb{Z}^N$. The distribution function f_N is invariant under particle exchange, which we can express by means of the permutation matrix **P**, such that

$$f_N(\mathbf{Px};\mathbf{Pp};t) = f_N(\mathbf{x};\mathbf{p};t), \qquad (46)$$

where each row and column of \mathbf{P} contain only one entry different from zero and equal to 1.

In order to derive the BBGKY hierarchy of the FPE in Eq. (1) we first define the *l*-particle distribution function:

$$f_l = \int_0^\lambda \frac{dx_{l+1}}{\lambda} \int_{-\infty}^\infty dp_{l+1} \cdots \int_0^\lambda \frac{dx_N}{\lambda} \int_{-\infty}^\infty dp_N f_N, \quad (47)$$

where f_l inherits the three properties in Eqs. (44), (45), and (46) from f_N . Index *l* takes the value l = 1, ..., N, such that for l =1 the distribution f_l is the single-particle phase-space function, and for l = N it describes the *N*-particle state. The evolution of f_l is found from Eq. (1) after integrating out the other N - lparticle variables and can be cast in the form

$$\frac{\partial f_l}{\partial t} = \sum_{j=1}^{l} \left(\mathcal{L}_j^{(l)} f_l + \mathcal{G}_j^{(l)} [f_{l+1}] \right), \tag{48}$$

where the first operator on the RHS solely depends on the variables of the *l* particles and reads

$$\mathcal{L}_{j}^{(l)}f_{l} = -\frac{\partial}{\partial x_{j}}\frac{p_{j}}{m}f_{l}$$

$$-S^{2}\frac{\partial}{\partial p_{j}}\sum_{i=1}^{l}[F_{0}\cos(kx_{i}) + \Gamma_{0}\sin(kx_{i})p_{i}]\sin(kx_{j})f_{l}$$

$$+S^{2}\frac{\partial}{\partial p_{j}}\sum_{i=1}^{l}\left(D_{0}\frac{\partial}{\partial p_{i}} + \eta_{0}\frac{\partial}{\partial x_{i}}\right)\sin(kx_{i})\sin(kx_{j})f_{l}.$$
(49)

The second operator, instead, depends nonlinearly on the (l + 1)-particle distribution function. This term vanishes when l = N, while for l < N it describes the dynamics of correlations, which are established by the interparticle potential. It reads

$$\mathcal{G}_{j}^{(l)}[f_{l+1}] = -S^{2}(N-l)\frac{\partial}{\partial p_{j}}\sin(kx_{j})$$
$$\times (F_{0}\Theta_{l}[f_{l+1}] + \Gamma_{0}\Xi_{l}[f_{l+1}]), \qquad (50)$$

where

$$\Theta_{l}[f_{l+1}] = \int_{0}^{\lambda} \frac{dx_{l+1}}{\lambda} \int_{-\infty}^{\infty} dp_{l+1} \cos(kx_{l+1}) f_{l+1}, \qquad (51)$$

$$\Xi_{l}[f_{l+1}] = \int_{0}^{\lambda} \frac{dx_{l+1}}{\lambda} \int_{-\infty}^{\infty} dp_{l+1} \sin(kx_{l+1}) p_{l+1} f_{l+1}, \quad (52)$$

while $\Theta_0[f_1] = \Theta_{MF}$ and $\Xi_0[f_1] = \Xi_{MF}$. Note that when the factorization ansatz of Eq. (9) holds, $\Theta_1[f_2] = \Theta_{MF}f_1$ and $\Xi_1[f_2] = \Xi_{MF}f_1$. A closed set of equations for f_l can thus be strictly obtained for l = N, giving Eq. (1), or for S = 0, hence in the absence of the cavity field.

B. The Lenard-Balescu equation

For l = 2 we can generally decompose the distribution function into two terms:

$$f_2(x_1, x_2, p_1, p_2) = f_1(x_1, p_1)f_1(x_2, p_2) + g_2(x_1, x_2, p_1, p_2),$$
(53)

where the first term on the RHS is the mean-field term and the second term describes all corrections beyond mean field. When at t = 0 the distribution function is factorized in a form like Eq. (9), the dynamics beyond mean field will tend to build correlations which are described by g_2 . We obtain the mean-field FPE, Eq. (10), by performing the approximation $\mathcal{G}_1^{(1)}[f_2] \rightarrow \mathcal{G}_1^{(0)}[f_1]f_1$. In the following we analyze the regime in which this approximation is justified by studying the equation describing the evolution of the function g_2 under some approximation, which permits us to truncate the BBGKY hierarchy to second order. This equation is known in the literature as the Lenard-Balescu equation [19], and it will allow us to identify a time scale where the mean-field treatment provides reliable predictions. In order to derive the Lenard-Balescu equation we first consider the distribution function for l = 3. Using the same type of decomposition as in Eq. (53), this can be written as

$$f_{3}(x_{1}, x_{2}, x_{3}, p_{1}, p_{2}, p_{3})$$

$$= f_{1}(x_{1}, p_{1})f_{1}(x_{2}, p_{2})f_{1}(x_{3}, p_{3})$$

$$+ \sum_{i,j,k=1}^{3} |\epsilon_{ijk}|f_{1}(x_{i}, p_{i})g_{2}(x_{j}, p_{j}, x_{k}, p_{k})$$

$$+ g_{3}(x_{1}, x_{2}, x_{3}, p_{1}, p_{2}, p_{3}),$$

where ϵ_{ijk} is the Levi-Civita tensor and g_3 describes all threebody correlations which cannot be written as a function of f_1 and/or f_2 . We assume now that g_3 is of higher order (from the treatment below we will see that $g_3 \propto 1/N^2$) and drop g_3 in the equation describing the dynamics of f_2 , Eq. (48). By means of this assumption we obtain two coupled equations for f_1 and f_2 , which can then be cast into the Lenard-Balescu equations for f_1 and g_2 using Eq. (53) and read

$$\frac{\partial f_{1}}{\partial t} = \mathcal{L}^{(1)} f_{1} + \mathcal{G}^{(1)}[f_{1}]f_{1} + \mathcal{G}^{(1)}[g_{2}],$$

$$\frac{\partial g_{2}}{\partial t} = -\frac{\partial}{\partial x_{1}} \frac{p_{1}}{m} g_{2} - \frac{\partial}{\partial x_{2}} \frac{p_{2}}{m} g_{2} - S^{2} \sum_{j=1}^{2} \sum_{i \neq j} \frac{\partial}{\partial p_{j}} F_{0} \sin(kx_{j}) \{\cos(kx_{i}) - \Theta_{\mathrm{MF}}[f_{1}]\} f_{1} f_{1}$$

$$-S^{2} \sum_{j=1}^{2} \sum_{i \neq j} \frac{\partial}{\partial p_{j}} \Gamma_{0} \sin(kx_{j}) \{\sin(kx_{i})p_{i} - \Xi_{\mathrm{MF}}[f_{1}]\} f_{1} f_{1} + S^{2} \sum_{j=1}^{2} \sum_{i \neq j} \frac{\partial}{\partial p_{j}} \sin(kx_{j}) \left(D_{0} \frac{\partial}{\partial p_{i}} + \eta_{0} \frac{\partial}{\partial x_{i}} \right) \sin(kx_{i}) f_{1} f_{1}$$

$$-NS^{2} F_{0} \sum_{j=1}^{2} \sum_{i \neq j} \frac{\partial}{\partial p_{j}} \sin(kx_{j}) \{\Theta_{1}[g_{2}]_{i} f_{1}(x_{j}, p_{j}) + \Theta_{\mathrm{MF}}[f_{1}]g_{2}\} - NS^{2} \Gamma_{0} \sum_{j=1}^{2} \sum_{i \neq j} \frac{\partial}{\partial p_{j}} \sin(kx_{j}) \{\Xi_{1}[g_{2}]_{i} f_{1}(x_{j}, p_{j}) + \Xi_{\mathrm{MF}}[f_{1}]g_{2}\},$$
(54a)

where we specified the arguments when necessary and introduced the notation $\Theta_1[g_2]_i$ and $\Xi_1[g_2]_i$ to indicate that these are functions of (x_i, p_i) .

The validity of the mean-field FPE, Eq. (10), relies on whether one can discard the term $\mathcal{G}^{(1)}[g_2]$ on the RHS of Eq. (54a). Let us recall the thermodynamic limit for which $S^2 \sim 1/N$. If we now assume that g_2 is of order 1/N with respect to f_1 , then the term $\mathcal{G}^{(1)}[g_2]$ is of order 1/N with respect to $\mathcal{G}^{(1)}[f_1]f_1$. A detailed analysis of Eq. (54b) shows that, if $g_2 \sim 1/N$ at t = 0, this scaling is preserved by the dynamics. In fact, (i) the first two terms on the RHS of Eq. (54b) give a scaling with 1/N because they are proportional to g_2 , while all other quantities are independent of N; (ii) the last term in the first line and the terms in the second line are all proportional to $S^2 \sim 1/N$, and (iii) the last two lines scale with $NS^2g_2 \sim$ 1/N. Therefore, for sufficiently short times the contribution of g_2 to the dynamics in the mean-field equation can be neglected.

We note that in Eq. (54a) the term $\mathcal{L}^{(1)}f_1$ also has components which scale with 1/N. If one consistently neglects all terms scaling with 1/N, then Eq. (54a) reduces to the

Vlasov equation (31) and therefore also neglects the diffusion processes leading to equilibrium. Figure 5 illustrates the order of magnitude of the corrections to the Vlasov and Lenard-Balescu equations, as well as the types of correlations that these describe.



Lenard-Balescu equation

FIG. 5. Illustration of the order of magnitude of the corrections of the Vlasov and of the Lenard-Balescu equations and of which type of correlations they include.



FIG. 6. Time evolution of the squared order parameter evaluated by numerically simulating the mean-field FPE, Eq. (10) (dash-dotted lines), and the N-particle FPE of Eq. (1) (solid lines). The curves correspond to different particle numbers N = 20 (top two brown curves), N = 50 (middle two green curves), and N = 200 (bottom two blue curves) and are calculated taking $\Delta_c = -\kappa$ and $\bar{n} = 2\bar{n}_c$. The number T of trajectories taken for the *N*-body FPE is T = 1000for N = 20, T = 500 for N = 50, and T = 100 for N = 200 (see Refs. [21,22] for details on the simulations). The horizontal dotted line indicates the asymptotic value of the squared order parameter. The inset shows the curves of the onset with the time axis rescaled by N. Note that the initial distribution of the full N-body FPE is the one which statistically corresponds to a spatially uniform distribution with the same temperature as the asymptotic one. Therefore, the value of $\langle \Theta^2 \rangle_N$ at t = 0 does not vanish due to finite-size effects. In order to compare these dynamics with the mean-field FPE, we have taken into account these finite-size effects in the initial mean-field distribution given by $\tilde{f}_0(x,p) = [1 + \delta_N \cos(kx)] f_1(p,0)$, where f_1 is given in Eq. (34) and δ_N is a spatial modulation depending on N.

C. Mean-field versus full N-atom dynamics

In order to complete our analysis of the limits of validity of the mean-field ansatz, we now compare its predictions with the ones obtained by numerical simulations of the *N*-particle FPE of Eq. (1). The latter are performed by means of stochastic differential equations (see Refs. [21,22] for details). We focus now on the evolution of the expectation value of Θ^2 , which explicitly depends on two-particle correlations and scales the strength of the conservative many-body potential. We recall the definition $\langle \cdot \rangle_N$ in order to indicate the mean value of an *N*particle observable taken over the *N*-particle distribution f_N .

Figure 6 compares the *N*-particle description where the evolution of f_N is governed by FPE (1) (solid line) and the mean-field description, where $f_N(\mathbf{x}; \mathbf{p}; t) =$ $f_1(x_1, p_1; t) \cdots f_1(x_N, p_N; t)$ and the evolution of f_1 is governed by the mean-field FPE (10) (dash-dotted line). The curves are plotted as a function of time and for different particle numbers, N = 20,50,200, where the parameter S^2 has been rescaled according to our thermodynamic limit to warrant a threshold \bar{n}_c which is independent of *N*. The parameters have been fixed so that, initially, the distribution is spatially uniform, while the momentum distribution is a Gaussian whose width coincides with the asymptotic temperature of the dynamics, Eq. (17). The strength of the field is such that $\bar{n} = 2\bar{n}_c$; therefore, the asymptotic spatial distribution is a Bragg grating with $|\Theta_{\rm MF}| \sim 0.83$. The dynamics we observe is the one which leads to the formation of the Bragg gratings starting from a uniform spatial distribution and exhibit three stages, which have been extensively discussed in Ref. [24]: a violent relaxation, a prethermalized phase, and a slow approach to equilibrium. The solid lines are simulations of the full FPE, and the dash-dotted lines are the corresponding mean-field prediction, which indeed qualitatively reproduces the three-stage dynamics.

The violent relaxation is a stage of the dynamics where there is good agreement between the mean-field and N-body FPEs. This is the short-time regime where the Vlasov equation, Eq. (31), is valid, and the behavior of the N-body FPE is reproduced by the one observed by numerically integrating the Vlasov equation [see Fig. 3(a)]. This has also been verified in Ref. [24]. The prethermalized regime is also predicted by the Vlasov equation [see Fig. 3(a)]. The mean-field FPE, however, provides a more accurate description and qualitatively reproduces the N-body FPE. Nevertheless, a clear difference between mean-field and N-body dynamics is found at the onset of the prethermalized stage: In fact, the oscillations are damped at a faster rate in the N-body FPE. Apart from this difference, there is a qualitative agreement between the mean-field and N-body FPEs also for this stage.

While both mean-field and *N*-body FPEs agree in the asymptotic value, we observe a striking difference between the two results in the relaxation to equilibrium after prethermalization. This is the stage where the role of dissipation and diffusion becomes relevant, as shown in Ref. [24] by comparing this behavior with the one where the dynamics is due to only the Hamiltonian term. In particular, the relaxation time scale predicted by the full simulation is about one order of magnitude larger than the corresponding mean-field prediction. This becomes even more evident by plotting the curves rescaling the time axis with N, which is visible in the inset. The curves of the mean-field FPE collapse to one curve, whereas the ones of the *N*-body FPE collapse to a significantly different curve.

Let us now summarize these results. First, the short-time behavior of the fluctuations of the order parameter are well described by the mean-field equation, in particular by the Vlasov equation. This is well understood in terms of the typical contributions to the dynamics: For short times the dominant contributions are indeed the terms of Eq. (31), and interparticle correlations are small, as we argued in the previous section. Discrepancies are due to finite-size effects. The prethermalized regime, moreover, exhibits good agreement between mean-field and full dynamics. This regime is dominated by the Hamiltonian dynamics, and the results show that Hamiltonian dynamics with long-range interactions is well reproduced by the meanfield description. Big deviations instead appear for long times, where the mean-field ansatz is expected to fail, and at the time scales dominated by relaxation to the stationary state.

VI. CONCLUSIONS

In this work we have systematically developed a mean-field description of the self-organization dynamics of atoms in a high-finesse cavity. The predictions of the mean-field model have been explored at equilibrium and out of equilibrium; its limits of validity have been tested by comparing them with those of the *N*-body FPE. We have found that the mean-field equation provides an excellent description of the dynamics when the latter is prevailingly Hamiltonian. It further describes the equilibrium properties of single-particle observables, including the asymptotic temperature and the order parameter. It fails, however, to reproduce the long-time out-of-equilibrium dynamics.

Despite these differences, this analysis shows that from the mean-field model one can analytically extract several predictions of the system dynamics. It is, indeed, remarkable that several predictions reproduce in the corresponding limits the ones obtained by means of other theoretical treatments, some of which start from a fully quantum mechanical treatment of the atoms. This, on the one hand, leads us to conjecture that quantum fluctuations play a marginal role in determining the steady-state properties of the cavity field. It further urges one to develop a full quantum kinetic theory, analogous to the full *N*-body semiclassical theory, which will overcome all limitations of simplifying theoretical assumptions performed so far. Only such a model, in fact, can give full access to the dynamical interplay between matter waves and cavity photons.

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APPENDIX: CAVITY-FIELD CORRELATION FUNCTION AT STEADY STATE

Experimentally accessible quantities are the correlation functions of the field at the cavity output, which allows one to monitor the atoms' state and is proportional to the intracavity field. In our formalism, the intracavity field is closely connected to the atomic state by the relation $E_{\text{cav}} \propto \sqrt{N\bar{n}}\Theta$; therefore, the correlation functions of the cavity field are proportional to the correlation functions of the magnetization Θ [21,22]. In the following we determine the autocorrelation function of the magnetization, which can be detected by means of the first-order correlation function of the field and the fourth-moment of the magnetization $\langle \Theta^4 \rangle_N$. As we showed in Ref. [21], in fact, $\langle \Theta^4 \rangle_N$ delivers the value of the intensity-intensity correlation of the field at zero-time delay and at zero order in the retardation effects.

1. Field intensity across the transition

We first determine the intracavity photon number n_{cav} at steady state for \bar{n} below, at, and above threshold. For this purpose we use the relation [21,22]

$$n_{\rm cav} = N\bar{n} \langle \Theta^2 \rangle_N,\tag{A1}$$

which, by introducing $\alpha = \bar{n}/\bar{n}_c$, can be cast in the form (see also the Sec. A 3)

$$n_{\rm cav} = \frac{1}{2}\bar{n}_c + \bar{n}\frac{\partial}{\partial\alpha}\mathcal{G}(\alpha), \tag{A2}$$

where

$$\mathcal{G}(\alpha) = \ln\left(\int_{-\infty}^{\infty} dy \exp\left\{-N\{\alpha y^2 - \ln[I_0(2\alpha y)]\}\right\}\right).$$
(A3)

We then analyze the prediction of this expression close to threshold for $\bar{n} \sim \bar{n}_c$ and thus $\alpha \sim 1$. For this purpose we expand the exponent of $\mathcal{G}(\alpha)$ about the value y = 0 and consider the behavior of n_{cav} for $\alpha \to 1^-$ and hence for $\bar{n} < \bar{n}_c$ but sufficiently close to the transition point that the truncation of the expansion is valid. In this limit we find

$$n_{\rm cav} \approx \frac{\bar{n}_c^2/2}{\bar{n}_c - \bar{n}},\tag{A4}$$

where the details of the derivation are reported in Sec. A 3. The value at the transition point is reported at leading order in N and reads (see Sec. A 3)

$$n_{\rm cav} \approx 2\sqrt{N}\bar{n}_c \frac{\Gamma\left[\frac{3}{4}\right]}{\Gamma\left[\frac{1}{4}\right]},\tag{A5}$$

where $\Gamma[x]$ denotes the gamma function [39].

The value of the intracavity photon number above threshold is found after observing that the exponent of function $\mathcal{G}(\alpha)$ has two minima that are given by the nonvanishing solutions of the fixed-point equation (24), which we denote by $\Theta_{MF} = \pm \overline{\Theta}$, with $\overline{\Theta}$ given in Eq. (26). Therefore, it holds that

$$n_{\rm cav} = N\bar{n}\Theta^2 \approx 2N(\bar{n}-\bar{n}_c),$$

sufficiently close to the critical point. In particular, the mean number of photons increases linearly with \bar{n} . We analyze now some properties of the first-order correlation function of the intracavity field, $g^{(1)}(\tau) = \lim_{t\to\infty} \text{Re}\langle E_{\text{cav}}(t+\tau)E_{\text{cav}}(t)\rangle_N$. This function has been extensively studied in Ref. [21] by numerically solving the *N*-particle FPE. Here, we will use the mean-field ansatz in order to better understand the two sidebands of its Fourier transform, at which it exhibits maxima above threshold. For this purpose we first notice that the correlation function is proportional to the autocorrelation function $C(\tau)$ of the magnetization by the relation $g^{(1)}(\tau) =$ $N\bar{n}C(\tau)$ and

$$C(\tau) = \lim_{t \to \infty} \langle \Theta(t) \Theta(t+\tau) \rangle_N.$$
 (A6)

We want to derive $C(\tau)$ in the mean field, and hence, the mean value now has to be taken over the factorized distribution as in Eq. (9) with the stationary mean-field distribution given in Eq. (23). We calculate $C(\tau)$ by solving the equations of the mathematical pendulum

$$\dot{x} = \frac{p}{m},$$

$$\dot{p} = 2\hbar k \Delta_c \bar{n} \bar{\Theta} \sin(kx),$$
(A7)

with initial conditions $x(0) = x_0$ and $p(0) = p_0$. The value $\overline{\Theta}$ is here the positive stable solution of Eq. (24). In the limit of small oscillations, these equations describe harmonic motion



FIG. 7. Contour plot of the spectrum of the autocorrelation function $\tilde{S}(\omega)$ as a function of \bar{n} and of the frequency (in units of κ) evaluated from the numerical data of $\Theta(x_1, \ldots, x_N)$, Eq. (4), by integrating the *N*-particle FPE, Eq. (1), for 100 trajectories of N = 50atoms, $\Delta_c = -\kappa$, and evolution time $t_{\text{tot}} = 10^4 \kappa^{-1}$ (see Ref. [21]). The lines are analytical estimates of the spectrum maximum for $\bar{n} > \bar{n}_c$. The dashed line corresponds to the frequency of the corresponding harmonic oscillator in Eq. (A8). The solid line is at the frequency extracted by solving Eq. (A7) for a mathematical pendulum and in good agreement with the peak position of the numerically evaluated spectra.

at the frequency

$$\omega_0 = \sqrt{-4\omega_r \Delta_c \bar{n}\bar{\Theta}}.$$
 (A8)

The mean frequency, however, is the result of the possible trajectories of the mathematical pendulum weighted by the probability density function $f_{st}(x_0, p_0)$. For $x_0 \neq 0$ and $p_0 \neq 0$ the oscillation period turns out to be larger than $2\pi/\omega_0$, and this prediction fits quite well the maximum found numerically by integrating the coupled equations of *N* particles, as shown in Fig. 7.

2. Intensity-intensity correlations at zero time delay

The intensity-intensity correlation function at zero time delay $g^{(2)}(0)$ provides a direct measurement of the fourth moment of the magnetization when retardation effects are sufficiently small [21]:

$$g^{(2)}(0) = \langle \Theta^4 \rangle_N / \langle \Theta^2 \rangle_N^2. \tag{A9}$$

Above threshold $\langle \Theta^n \rangle_N = \overline{\Theta}^n + O(1/N)$, with $\overline{\Theta}$ being the solution of Eq. (24). Therefore, for $\overline{n} > \overline{n}_c$ we obtain

$$g^{(2)}(0)_{\bar{n}>\bar{n}_c} = 1,$$
 (A10)

which corresponds to coherent light and is valid at leading order, with an error that scales with 1/N. In the mean field for the factorized distribution, Eq. (9), we get

$$\langle \Theta^2 \rangle_N = \frac{1}{N} \mathcal{B} + \frac{N-1}{N} \bar{\Theta}^2$$

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and

$$\begin{split} \langle \Theta^4 \rangle_N = & \frac{N(N-1)(N-2)(N-3)}{N^4} \bar{\Theta}^4 \\ &+ \frac{6N(N-1)(N-2)}{N^4} \bar{\Theta}^2 \mathcal{B} + \frac{3N(N-1)}{N^4} \mathcal{B}^2 \\ &+ \frac{4N(N-1)}{N^4} \bar{\Theta} \langle \cos^3(x) \rangle + \frac{N}{N^4} \langle \cos^4(x) \rangle. \end{split}$$

Notice that above threshold for $\overline{\Theta} \neq 0$ we can again write $\langle \Theta^4 \rangle_N = \overline{\Theta}^4 + O(1/N)$. Hence, we get the same value for $g^{(2)}(0) = 1$ (above threshold) in the thermodynamic limit $N \to \infty$. Below threshold, in Sec. A 3 we show that the expression takes the value

$$g^{(2)}(0)_{\bar{n}<\bar{n}_c}=3,\tag{A11}$$

which corresponds to super-Poissonian light. Corrections scale with 1/N. This also holds for the calculation with the factorized ansatz. Below threshold we get

$$\langle \Theta^2 \rangle_N = \frac{1}{N} \mathcal{B}$$

and

$$\langle \Theta^4 \rangle_N = \frac{3}{N^2} \mathcal{B}^2 + O\left(\frac{1}{N}\right)$$

and therefore the same value of $g^{(2)}(0) = 3$ (below threshold) as for the *N*-particle description. Finally, at threshold we obtain

$$g^{(2)}(0)_{\bar{n}=\bar{n}_c} \approx \frac{1}{4} \left(\frac{\Gamma\left[\frac{1}{4}\right]}{\Gamma\left[\frac{3}{4}\right]} \right)^2, \tag{A12}$$

with corrections scaling with $1/\sqrt{N}$, thus giving a slower convergence than the one found for the values above and below threshold. We want to mention here that the meanfield description cannot reproduce the value in Eq. (A12). Figure 8 displays the mean-field predictions for $g^{(2)}(0)$ at the thermodynamic limit and as a function of \bar{n} . These curves are compared with the mean-field calculation at finite *N* and with the corresponding one of the *N*-particle FPE. Even though the mean-field curve at finite *N* is tendentially closer to the thermodynamic limit than the *N*-particle FPE prediction, they both converge to the values of Eqs. (A10), (A12), and (A11), depending on whether $\bar{n} < , = , > \bar{n}_c$ for $N \to \infty$.

3. Useful relations

In order to demonstrate Eq. (A2) we first consider the relation

$$\int_{-\infty}^{\infty} dy \exp\left[-\alpha N\left(y - \frac{1}{N}\sum_{i=1}^{N}\cos(kx_i)\right)^2\right] = \sqrt{\frac{\pi}{\alpha N}}$$

and cast it into the form

$$\int_{-\infty}^{\infty} dy e^{-\alpha N y^2} \exp\left(2\alpha N y \frac{1}{N} \sum_{i=1}^{N} \cos(kx_i)\right)$$
$$= \sqrt{\frac{\pi}{\alpha N}} \exp[\alpha N \Theta(\mathbf{x})^2].$$



FIG. 8. Intensity-intensity correlation function at zero time delay $g^{(2)}(0)$, Eq. (A9), as a function of \bar{n} for (a) N = 50 atoms and (b) N = 200 atoms. The solid blue lines are the curves evaluated using in Eq. (A9) the mean-field steady state (23). The dashed lines are calculated for the corresponding full *N*-particle distribution given in [17]. The horizontal black solid lines are the values at the thermodynamic limit given at $\bar{n} < \bar{n}_c$ by Eq. (A11) and at $\bar{n} > \bar{n}_c$ by Eq. (A10). The point at $\bar{n} = \bar{n}_c$ is at the value of Eq. (A12). The discrepancy between the mean-field curve and the full *N*-particle predictions decreases as $N \to \infty$, where they both converge to the value given by the thermodynamic limit.

From these relations, it follows

$$\ln\left\{\frac{1}{\lambda^{N}}\int d\mathbf{x} \exp[\alpha N\Theta(\mathbf{x})^{2}]\right\}$$
$$=\frac{1}{2}\ln\left(\frac{N}{\pi}\alpha\right)$$
$$+\ln\left(\int_{-\infty}^{\infty} dy \exp\left\{-N\left\{\alpha y^{2}-\ln\left[I_{0}(2\alpha y)\right]\right\}\right\}\right).$$

We use it for evaluating expression (A1) and obtain

$$n_{\text{cav}} = \bar{n} \frac{\partial}{\partial \alpha} \ln \left\{ \frac{1}{\lambda^{N}} \int d\mathbf{x} \exp[\alpha N \Theta(\mathbf{x})^{2}] \right\}$$
$$= \bar{n} \left(\frac{1}{2\alpha} + \frac{\frac{\partial}{\partial \alpha} \int_{-\infty}^{\infty} dy \exp\left\{-N\{\alpha y^{2} - \ln[I_{0}(2\alpha y)]\}\right\}}{\int_{-\infty}^{\infty} dy \exp\left\{-N\{\alpha y^{2} - \ln[I_{0}(2\alpha y)]\}\right\}} \right),$$
(A13)

which leads to Eq. (A2) by using definition (A3).

In order to determine the intracavity photon number close to threshold, we expand the exponent of Eq. (A3) about y = 0 to fourth order:

$$\alpha y^{2} - \ln[I_{0}(2\alpha y)] = \alpha(1-\alpha)y^{2} + \frac{\alpha^{4}}{4}y^{4} + O(y^{6}).$$

For $\bar{n} < \bar{n}_c$, the coefficient of the quadratic term is positive, and we thus discard the fourth-order term. Expression (A2) takes the form

$$\begin{split} n_{\text{cav}} &\approx \bar{n} \left(\frac{1}{2\alpha} + \frac{\frac{\partial}{\partial \alpha} \int_{-\infty}^{\infty} dy \exp\left[-N\alpha(1-\alpha)y^2 \right]}{\int_{-\infty}^{\infty} dy \exp\left[-N\alpha(1-\alpha)y^2 \right]} \right) \\ &= \bar{n} \left(\frac{1}{2\alpha} + \frac{2\alpha - 1}{2\alpha(1-\alpha)} \right) = \frac{\bar{n}}{2(1-\alpha)}. \end{split}$$

Using the explicit value of α ,

ł

$$n_{\rm cav} = \frac{\bar{n}\bar{n}_c}{2(\bar{n}_c - \bar{n})} \approx \frac{\bar{n}_c^2/2}{\bar{n}_c - \bar{n}},$$
 (A14)

which thus gives Eq. (A4).

At the transition point $\bar{n} = \bar{n}_c$ the integral in Eq. (A2) diverges in the limit $N \to \infty$. We determine its value for finite N and keep the leading order. Moreover, since the coefficient of the quadratic term in the expansion in y vanishes, we include the fourth order and evaluate the integral at $\alpha = 1$, obtaining

$$\begin{split} n_{\rm cav} &\approx \bar{n}_c \left(\frac{1}{2} + \frac{\int_{-\infty}^{\infty} dy (Ny^2 - Ny^4) \exp\left[-\frac{N}{4} y^4 \right]}{\int_{-\infty}^{\infty} dy \exp\left[-\frac{N}{4} y^4 \right]} \right) \\ &\approx \bar{n}_c \frac{2\sqrt{N}\Gamma\left[\frac{3}{4} \right]}{\Gamma\left[\frac{1}{4} \right]}, \end{split}$$

which is the expression in Eq. (A5).

To calculate $g^{(2)}(0)$ below and at threshold we notice that

$$N^{2} \langle \Theta^{4} \rangle_{N} - N^{2} \langle \Theta^{2} \rangle_{N}^{2} = \frac{\partial^{2}}{\partial \alpha^{2}} \ln \left\{ \frac{1}{\lambda^{N}} \int d\mathbf{x} \exp[\alpha N \Theta(\mathbf{x})^{2}] \right\}$$
$$= N \frac{\partial}{\partial \alpha} \langle \Theta^{2} \rangle_{N}$$

holds. Below threshold for $\alpha < 1$ we calculated in leading order that

$$\frac{\partial}{\partial \alpha} \frac{1}{2(1-\alpha)} = \frac{1}{2(1-\alpha)^2},$$

which then delivers the expression

$$g^{(2)}(0)_{\bar{n}<\bar{n}_c} = \frac{\frac{1}{2(1-\alpha)^2} + \frac{1}{4(1-\alpha)^2}}{\frac{1}{4(1-\alpha)^2}} = 3$$

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and thus Eq. (A11). In order to calculate the value at threshold we use

$$N^{2} \langle \Theta^{4} \rangle_{N} - N^{2} \langle \Theta^{2} \rangle_{N}^{2} \approx N - 4N \left(\frac{\Gamma\left[\frac{3}{4}\right]}{\Gamma\left[\frac{1}{4}\right]} \right)^{2}, \qquad (A15)$$

which is valid in leading order and which gives Eq. (A12).

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