Development and regression of a large fluctuation

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We study the evolution leading to (or regressing from) a large fluctuation in a statistical mechanical system. We introduce and study analytically a simple model of many identically and independently distributed microscopic variables n_m (m = 1, M) evolving by means of a master equation. We show that the process producing a nontypical fluctuation with a value of $N = \sum_{m=1}^{M} n_m$ well above the average $\langle N \rangle$ is slow. Such process is characterized by the power-law growth of the largest possible observable value of N at a given time t. We find similar features also for the reverse process of the regression from a rare state with $N \gg \langle N \rangle$ to a typical one with $N \simeq \langle N \rangle$.

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I. INTRODUCTION

The occurrence of fluctuations is at the heart of most physical phenomena [1]. Typically, in an extended system made of a large number M of microscopic constituents, like those usually considered in equilibrium thermodynamics, a collective variable N (like the particle number or the energy) evolves as to stay most of the time close to its average value $\langle N \rangle$. Large deviations are rare and become progressively less frequent as N moves away from the average. For this reason they are neglected in many practical applications. However, in some cases they can have important consequences. This happens, for instance, when their occurrence leads the system to an absorbing state, namely, a configuration that cannot be escaped [2]. Examples include the extinction of a species, the failure of a device, or the bankruptcy of a company. The latter indeed was the first problem for which the rigorous results of large deviation theory were applied [3]. In addition, large deviations play a prominent role in many nonequilibrium phenomena, e.g., in the decay of metastable states [4].

Configurations corresponding to a large fluctuation are usually very different from those typically observed when $N \simeq \langle N \rangle$ (to ease the notation we use the same symbol for the stochastic variable and its possible outcomes), because the system explores a seldom visited region of phase space that may have peculiar properties. Hence, the question arises of how the representative point moves to reach such lowprobability sectors, or, in other words, what are the properties of the dynamical process producing a large deviation. This issue is not only an important and largely unexplored topic in large deviation theory, but might also represent a first step towards the detection and control of fluctuations, with important applications concerning the predictability of catastrophic events. Likewise, the reverse process, whereby the typical behavior is recovered after a rare event, has also theoretical and practical interest.

In this paper we study such problems in a simple but sufficiently general model where $N = \sum_{m=1}^{M} n_m$ is the sum of a large number M of independent variables $n_m = 0, 1, ...$ identically distributed with probability $p(n_m)$. The creation of a fluctuation is studied by evolving the initial distribution $p(n,0) = p(n_m = n, t = 0)$ of the microvariables n_m in a typical state of the system with $N = \langle N \rangle$, until a deviation with $N \neq \langle N \rangle$ is observed. Solving the master equation yields the evolution of the probability P(N, M, t) that the collective variable takes a given value N. This quantity provides a detailed description of the whole fluctuation spectrum of N, characterizes the event whereby the fluctuation is built, and identifies its relevant properties. In the same way, one can study the disruption of a large deviation by studying the evolution from an initial condition with $N \neq \langle N \rangle$.

In the present article we choose the master equation governing the dynamics of p(n,t) such as to have the stationary solution

$$p_{st}(n) \propto (n+1)^{-k}.$$
 (1)

Systems with fat tail distributions analogous to the one considered here are found in natural sciences, social sciences, and economics. Among many examples we can mention the magnitude of earthquakes [5], the spreading of forest fires [6], rain events [7], size of cities [8], wealth distribution [9], price returns of stock indices [10], and degree distribution of networks [11].

Our choice of p_{st} not only is motivated by its ubiquitous character, but also stems from general considerations regarding the actual probability to observe large fluctuations. As we discuss below, such probability is particularly large for the model under consideration.

We have already mentioned that in extended systems the large deviations of a collective variable N are generally strongly suppressed. Indeed, the probability P(N,M) usually obeys [12] the following *large deviation principle* (LDP):

$$\lim_{M \to \infty} \ln P(N, M) = -MR(\rho), \tag{2}$$

where $\rho = \frac{N}{M}$ and *R* is the *rate function*. For simplicity, in the above equation and in the following we omit the dependence on *t*. Equation (2) implies that P(N, M) is always exponentially small (in *M*) except for the values of the *density* ρ for which $R(\rho) = 0$. Notice that at least one such configuration is bound to exist in order to preserve normalization of probability as $M \to \infty$. The simplest case is when there is a single value of ρ yielding $R(\rho) = 0$. This value trivially coincides with the average $\langle \rho \rangle = \frac{\langle N \rangle}{M}$ for large *M*. In this scenario, the outcome of a measurement is almost always close to $\langle \rho \rangle$, whereas sizable fluctuations are extremely rare and can be observed only in systems of mesoscopic scale, namely, with *M* not too large. The same mathematical structure applies to

all those cases where *N* is formed by the addition of many microscopic contributions. This is nicely illustrated by the much studied [13] problem of the fluctuations of a *charge* $N = \int_0^{\tau} j(t) dt$, *j* being a current, flowing in a certain time interval τ through a system.

Despite all of the above, there are important cases where large fluctuations are not exponentially suppressed as in Eq. (2), because the LDP breaks down for some range of values of *N*. In this situation, $R(\rho)$ can vanish not only for one (or some isolated) specific value(s) of ρ , but in a whole interval $\rho \in \mathcal{I}$:

$$R(\rho) \equiv 0, \quad \text{for} \quad \rho \in \mathcal{I}, \tag{3}$$

In this case Eq. (2) must more properly be rewritten as

$$\ln P(N,M) \simeq -M[R(\rho) + \mathcal{R}(M,\rho)], \quad \text{for large } M. \quad (4)$$

When the LDP holds, *R* is finite and the second term on the right-hand side of the equation above is a correction to the leading behavior for finite $M: \lim_{M\to\infty} \frac{\mathcal{R}(M,\rho)}{R(\rho)} = 0$. Then, in the large-*M* limit, Eq. (2) is meaningful, since fluctuations are fully described by *R*. However, if Eq. (3) holds, then \mathcal{R} becomes the relevant term and Eq. (2) is useless. This implies that fluctuations are reduced more softly in \mathfrak{I} than where Eq. (2) holds, and, depending on the system, significant deviations may have a good chance to form and be detected.

Situations where Eq. (3) holds are observed in a variety of systems. A first notable example is represented by magnetic materials where, below the critical temperature, the probability distribution of the (fluctuating) spontaneous magnetization Nexhibits [14] a structure like the one in Eq. (4), where M is the number of spins. In this case R vanishes in the whole region \mathfrak{I} such that $\rho \in [-\rho_0, \rho_0]$, where ρ_0 is the absolute value of the (average) spontaneous magnetization (per spin). Another example is the case of a Brownian walker on a line, with hopping rates retaining memory of the previous history [15]. The probability of moving a certain distance N after M steps takes a form like the one of Eq. (4). For instance, for a particular choice of the memory term, Rvanishes in the entire region $\rho > 0$, i.e., when the particle, starting from the origin, moves to the right. Other examples include fluctuations of driven Maxwell-Lorentz particles [16], quantum quenches [17], disordered systems [18], and many others [12,19].

As already anticipated, the model that we will study in this article is conceived in such a way that the stationary probability $P_{st}(N,M)$ obeys the crucial LDP-breaking property [Eq. (3)]. Indeed, it is well known [20] that assuming the distribution Eq. (1) with k > 2, it follows that $-\lim_{M\to\infty} M^{-1} \ln P_{st}(N,M) = R_{st}(\rho) = 0$ in the entire range $\mathcal{I} \equiv \{\rho \mid \rho > \langle \rho \rangle\}$. Correspondingly, the LDP [Eq. (2)] does not apply in that region (notice that the same phenomenon does not occur for $k \leq 2$; see the discussion in Sec. III). As a consequence, the condition for the observability [in the sense discussed above, after Eq. (3)] of large deviations is met in the interval $\rho \in \mathcal{I}$. The detailed theoretical study of the formation and regression of a large deviation that we will carry out in the present article for an analytically tractable model could therefore pave the way to the experimental investigation of such processes. Moreover, the model studied here could

represent a simple paradigm for a class of systems, like those mentioned above, where the condition (3) is satisfied. For example, in the previously discussed magnetic context, the problem at hand would correspond to study the spontaneous process whereby a typical configuration with $\rho = \rho_0$ evolves to a probabilistically unfavored one with $\rho < \rho_0$ due to thermal fluctuations, and the regression to the initial state.

Probabilistic setups, similar in spirit to the one introduced and studied in the present work, have been already considered in models of simplicial quantum gravity [21,22], non-equilibrium-driven systems [23], Lévy walks [24], and other systems [25] (see Sec. II for more details). However, to the best of our knowledge, the process whereby fluctuations form and regress has never been previously investigated.

Due to the analytical tractability of the model, we can derive several significant results from its solution. First, the nature of the process associated with the production and regression of fluctuations is radically different if it occurs in the region $\mathcal{I} \equiv \{\rho \mid \rho \leq \langle \rho \rangle\}$ where the LDP holds, or in \mathfrak{I} , where it is violated. In the former case the evolution is fast and relatively simple, with an exponential convergence towards the stationary form. In the latter, it displays a slow nontrivial evolution. This happens because the mechanism whereby P(N,M,t) approaches $P_{st}(N,M)$ is effective only up to a typical finite fluctuation scale $N \sim v(t)$, while larger values of N are left untouched. The characteristic value v(t) increases slowly in an algebraic way. This leads to an everlasting aging phenomenon, which closely resembles the dynamics of systems crossing a phase transition [26]. The probability P(N, M, t) attains the stationary value for increasingly large values of N. However, at each time there will always exist a sufficiently large value of N beyond which stationarity is not reached.

Second, considering the growth of spontaneous fluctuations, the mechanism whereby LDP breaks down, starting from an initial configuration that satisfies it, shows a nontrivial interplay between M and t. Specifically, while violations of LDP are enhanced at large times, as expected, they are reduced by increasing M. This shows that, when studying the large time behavior of a fluctuating system, attention should be paid to the order of the limits $\lim_{M\to\infty}$ and $\lim_{t\to\infty}$, which again reminds us of the physics of phase transitions.

This paper is organized as follows: In Sec. II we introduce the statistical model and set the notations. In Sec. III we discuss the properties of the model in the large-time limit when stationarity is reached. The breakdown of LDP is discussed and related to a *condensation* phenomenon. In Sec. IV the formation (Sec. IV A 1) and suppression (Sec. IV B) of fluctuations are studied by solving analytically the master equation for the microscopic probabilities p(n,t) and inferring the time evolution of the global probability P(N,M,t). The phenomenon of *partial condensation* is also discussed. Finally, in Sec. V we briefly summarize our results and draw our conclusions.

II. THE STATISTICAL MODEL

We consider *M* independent random variables that take integer values $n_m = 0, 1, 2, ...$ (m = 1, M) subject to a probability distribution $p(n_m)$, which in general depends on some parameters among which, possibly, the time *t*.

The closely related problem where n_m are continuous variables behaves very similarly. In the following, in order to make the idea more concrete, we will speak of *M* boxes containing a total number

$$N = \sum_{m=1}^{M} n_m \tag{5}$$

of *particles*, with an average value $\langle N \rangle = \sum_{m=1}^{M} \langle n_m \rangle$, where $\langle n_m \rangle = \sum_n n p(n)$.

Particles can be exchanged with the external environment. Assuming that the dynamics amounts to elementary moves where a single entity can be added to or removed from a specific box, the probability p obeys the following master equation:

$$\frac{dp(n,t)}{dt} = -[w^{+}(n) + w^{-}(n)]p(n,t) + w^{-}(n+1,t)p(n+1,t) + w^{+}(n-1,t)p(n-1,t),$$
(6)

where $w^+(n)$ is the transition rate to increase the number of particles, i.e., $n \rightarrow n + 1$, and $w^-(n)$ the one to decrease it, $n \rightarrow n - 1$. Here and in the following we denote by p(n,t) the probability, making the time dependence explicit. The master equation (6) is completely general for systems of discrete variables where *N* is not conserved as, e.g., spin models (Ising, Potts, Clock, etc.) [27]. We consider the following transition rates:

$$w^{+}(n) = (n+2)^{-k}$$

$$w^{-}(n) = n^{-k}(1-\delta_{n,0}).$$
(7)

The Kronecker δ function guarantees that particles cannot be extracted from an empty box. The form (7) is such that the evolution of a large cluster of *n* particles located in a single box is much slower than in the less populated ones, similarly to what happens in certain models of irreversible aging processes [28]. Notice that the transition rates (7) obey detailed balance:

$$\frac{w^+(n-1)}{w^-(n)} = \frac{p_{st}(n)}{p_{st}(n-1)}, \quad \forall n > 0,$$
(8)

where

$$p_{st}(n) = \zeta^{-1}(k)(n+1)^{-k}$$
(9)

and the normalization factor $\zeta(k) = \sum_{n=0}^{\infty} (n+1)^{-k}$ is the Riemann ζ function. Let us stress that the more general choice for a transition rate obeying the detailed balance condition (8) is $w^+(n) = (n+2)^{-k}g(n+1)$, $w^-(n) = n^{-k}(1-\delta_{n,0})g(n)$, where g is an arbitrary function. Here we make the simplest choice $g(n) \equiv 1$.

The form of p_{st} above provides an average occupation

$$\langle \rho \rangle = \langle n \rangle = \frac{\zeta(k-1)}{\zeta(k)} - 1,$$
 (10)

a result that will be useful in the following. Notice also that $\langle \rho \rangle$ exists only for k > 2, and, similarly, there is a finite variance only for k > 3.

With the w^{\pm} of Eq. (7) the master equation (6) reads

$$\frac{dp(n,t)}{dt} = -[(n+2)^{-k} + n^{-k}]p(n,t) + (n+1)^{-k}[p(n+1,t) + p(n-1,t)], \quad \forall n > 0$$

$$\frac{dp(0,t)}{dt} = -2^{-k}p(0,t) + p(1,t).$$
(11)

The probability to have a total number N of particles at time t is

$$P(N,M,t) = \sum_{n_1,n_2,...,n_M} p(n_1,t)p(n_2,t)\cdots p(n_M,t)\,\delta_{\mathcal{N},N}$$

= $\frac{1}{2\pi i} \oint dz \, e^{M[\ln Q(z,t) - \rho \ln z]},$ (12)

where $\mathcal{N} = \sum_{m=1}^{M} n_m$, we have used the representation $\delta_{\mathcal{N},N} = \frac{1}{2\pi i} \oint dz \, z^{-(N-\mathcal{N}+1)}$ and

$$Q(z,t) = \sum_{n} p(n,t)z^{n}$$
(13)

with $\rho = \frac{N+1}{M} \simeq \frac{N}{M}$ is the particle density. The relation

$$P(N,M,t) = \sum_{n=0}^{N} \pi(n,N,M,t)$$
(14)

with

$$\pi(n, N, M, t) = P(N - n, M - 1, t) p(n, t)$$
(15)

is easily proved; see, for instance, Ref. [22]. $\pi(n, N, M, t)$ is the conditional probability that, at time *t*, there are *n* particles in the *M*th box, given that a total number *N* is found in all the boxes. Since the random variables are identically distributed, the same probability applies to a generic box, not only to the *M*th. The recursion (14) allows one to determine the probability distribution of *M* variables from the one for M - 1. Specifically, once *p* is known from the solution of the evolution equations (11), Eqs. (14) and (15) can be used with the *boundary condition* P(N, M = 1, t) = p(n = N, t) to obtain *P*, step by step, for larger and larger values of *M*.

It should be stressed that Eq. (14) makes the exact determination of *P* feasible also for reasonably large values of *N* and *M*. Indeed, the computational complexity using this formula is only polynomial, whereas there is an exponential

number of redundant operations involved in the determination of P by using the first line of Eq. (12). This point is discussed in more detail in Appendix A.

Besides Eq. (14), which is always exact, for large M one can alternatively determine P by evaluating the integral in Eq. (12) by the method of steepest descent

$$P(N,M,t) \simeq e^{-MR(\rho,t)},\tag{16}$$

where

$$R(\rho, t) = -\ln Q[z^*(\rho, t)] + \rho \ln z^*(\rho, t)$$
(17)

is the rate function and z^* is the value of z for which the exponential argument in Eq. (12) is maximum. This is provided by the following saddle-point equation:

$$z^* \frac{Q'(z^*,t)}{Q(z^*,t)} = \rho.$$
 (18)

As we will see soon, however, a straightforward saddle-point evaluation of the integral in Eq. (12) is not always doable.

In this paper the model introduced insofar is studied to understand the basic mechanisms governing the occurrence of fluctuations and the mathematical structure behind. As already pointed out in the introduction, its formulation is similar to other, physically inspired and intensively studied models of statistical mechanics.

To begin with, collections of independent identically distributed random variables obeying Eq. (9) have been introduced as a simple description of quantum gravity [21,22]. This same model is also sometimes referred to as *urn* model, or *balls and boxes* model. In this approach N is an external control parameter. This means that, at variance with our analysis, fluctuations of this quantity are forbidden by construction. What is usually studied in that context are, instead, the properties of the stationary state as the control parameters k and N are varied. The nonequilibrium dynamics following an abrupt change of k (playing the role of an inverse temperature) has also been considered in [29]. The evolution of the model in that case, however, is ruled by a N-conserving stochastic equation different from Eq. (11).

Another class of related problems are descriptions of non-equilibrium-driven systems with particles hopping on a lattice, like the zero range process [23]. In these systems the probability at stationarity is factorized into single-site distributions that, for particular choices of the hopping rates, can take the form (9) [23]. N is a conserved quantity also in these cases. Furthermore, the properties of independent random variables distributed according to Eq. (9) have been discussed in relation to a wealth of different physical situations, like in the notable case of Lévy walks [24] and in other models [25].

III. STATIONARY STATE

Equation (11) has the stationary solution (9). The properties of the probability $P_{st}(N,M)$ in the stationary state have been studied elsewhere [20]. Here we briefly mention some basic results that will be needed in the following. We first derive them in a somewhat simplified framework that provides physical hints to the mathematically more refined exposition that will be presented in the following section.

A. Simplified framework

A relatively simple description of the properties of the stationary state can be obtained by considering the large-*k* behavior. In this limit $\langle \rho \rangle \rightarrow 0$. In fact, given the form of the microscopic probabilities p_{st} in Eq. (9), as *k* grows the chance of a nonvanishing outcome *n* becomes progressively smaller. This result can be easily derived from the exact expression (10). Then, for large *k*, deviations with $\langle \rho \rangle \ll \rho \ll 1$ are possible. We will focus our analysis in this range of densities.

Since the variables n_m are identically distributed there is an obvious symmetry among the boxes. Then, if such a symmetry is not spontaneously broken, the representative configurations of the stationary state are expected to have the *N* balls fairly distributed among all the boxes. In this case, given that $\rho \ll 1$, most of them will be empty, and a comparatively smaller number, of order *N*, will contain one ball. Given that *k* is large, the chance for a single box to host more than one particle is very small and will be neglected in the following. The probability to have states with fairly distributed particles is

$$P_{\text{sym}}(N,M) = p_{st}(1)^N \cdot p_{st}(0)^{M-N} \cdot \Omega(N,M)$$
$$= \zeta(k)^{-M} \cdot 2^{-kN} \cdot \Omega(N,M), \qquad (19)$$

where $\Omega(N, M) = {M \choose N}$ is the number of ways to choose the *N* occupied sites out of the total *M* and we have used Eq. (9).

We show now that, for large M, the probability P_{sym} of this symmetric state can be negligible compared to that of a *condensed* one, where the symmetry among the boxes is broken and a macroscopic number $n_c \propto M$ of particles is accumulated in one of them. The probability P_{cond} of such a state is

$$P_{\text{cond}}(N,M) = M \cdot p_{st}(n_c) \cdot P_{\text{sym}}(N - n_c, M - 1).$$
(20)

Here the first term, $M \cdot p_{st}(n_c)$, represents the probability to place n_c particles in the condensing box (the factor M in front accounts for the M ways to choose it). The last term, P_{sym} , is the probability [given by Eq. (19)] associated with the remaining $N - n_c$, which are uniformly spread among the remaining M - 1 boxes.

At large *M*, using the Stirling approximation for Ω and introducing the density $\overline{\rho}$ of balls in noncondensing boxes through $n_c = (\rho - \overline{\rho})M$ one finds

$$\frac{P_{\text{sym}}(N,M)}{P_{\text{cond}}(N,M)} \simeq e^{-Mk\{(\rho-\overline{\rho})\ln 2 - k^{-1}[s(\rho) - s(\overline{\rho})]\}}, \qquad (21)$$

where $s(x) = -x \ln x - (1 - x) \ln(1 - x)$ and subdominant terms have been dropped. Equation (21) shows that the formation of the condensed state is surely favored for $\rho > \overline{\rho}$, because for large *k* the first term in the argument of the exponential, namely, the positive quantity $(\rho - \overline{\rho}) \ln 2$, prevails over the second. Similarly, the formation of the condensed phase is unfavored for $\rho < \overline{\rho}$.

The discussion presented insofar is valid for large k. However, the basic results apply also to the small-k regime, provided that k > 2. This will be shown with the somewhat more refined calculation sketched in Sec. III B. We will also identify $\overline{\rho}$ with the average value $\langle \rho \rangle$, given in Eq. (10), and establish that condensation always occur, when k > 2, for $\rho > \langle \rho \rangle$.

B. Some mathematical refinements

We now show how the results obtained with the simple approach of the preceding section are confirmed by a more accurate treatment of the model equations where the condition of large k is released. At stationarity and for large M one has

$$Q(z) = \zeta^{-1}(k) \sum_{n=0}^{\infty} (n+1)^{-k} z^n = \zeta^{-1}(k) \frac{Li_k(z)}{z}, \quad (22)$$

where $Li_k(z)$ is the polylogarithm (Jonquière's function), and therefore

$$zQ'(z) \simeq \zeta^{-1}(k) \frac{Li_{k-1}(z) - Li_k(z)}{z}.$$
 (23)

The saddle-point condition (18) then reads

$$\frac{Li_{k-1}(z^*)}{Li_k(z^*)} = \rho + 1.$$
(24)

For $k \leq 2$ this equation always admits a solution and this corresponds to the fact that condensation does not occur. In the following we will concentrate on the sector with k > 2. In this case Eq. (24) has solution only in the region \mathcal{I} with $\rho \leq \langle \rho \rangle$ [21,22,25,29], where $\langle \rho \rangle$ is given in Eq. (10). In this range Eq. (16) holds with

$$R_{st}(\rho) = -\ln\left\{\zeta(k)^{-1} \frac{Li_k[z^*(\rho)]}{z^*(\rho)}\right\} + \rho \ln z^*(\rho).$$
(25)

In the complementary sector \mathfrak{X} , that is for $\rho > \langle \rho \rangle$, a straightforward saddle-point approach is not available and the phenomenon of condensation occurs, namely, a macroscopic number $n_c \simeq N - \langle N \rangle$ of particles, those that cannot be accommodated in the *normal* state, is accumulated in a single box, as discussed in Sec. III A. In this case $P_{st}(N,M)$ is, for large N, determined by the probability $Mp_{st}(N - \langle N \rangle)$ that a single box contains such a huge amount of particles. Comparing with Eq. (20), this shows that $\overline{\rho} = \langle \rho \rangle$. In summary, one finds the following behavior:

$$P_{st}(N,M) \sim \begin{cases} e^{-MR_{st}(\rho)} & \text{for } \rho \leqslant \langle \rho \rangle \\ Mp_{st}(N-\langle N \rangle) & \text{for } \rho \gg \langle \rho \rangle, \end{cases}$$
(26)

with R_{st} given in Eq. (25).

The above expressions, covering the regions $\rho \leq \langle \rho \rangle$ and $\rho \gg \langle \rho \rangle$ only, can be derived analytically in the large-*M* limit. The exact expression for P_{st} , which is valid for any *M* and ρ , can instead only be obtained by iteration of the recursion (14), and is plotted in the lower half of Fig. 1 (continuous green curve) considering the case with k = 3 and M = 100. The curve has a maximum around $\langle N \rangle \simeq 36.84$. We also compare the exact solution with the power-law behavior of Eq. (26) (dashed violet line), finding perfect agreement for $N \gg \langle N \rangle$.

Equation (26) shows that fluctuations with $\rho \leq \langle \rho \rangle$ behave *normally*, in the sense that a large-deviation principle with rate function R_{st} is obeyed. On the other hand, fluctuations with $\rho > \langle \rho \rangle$ are peculiar since P_{st} is not exponentially suppressed in M and relatively large fluctuations are possible. Indeed, rewriting the second line of Eq. (26) in terms of the density as

$$P_{st}(N,M) \sim M^{1-k} \zeta^{-1}(k) (\rho - \langle \rho \rangle)^{-k} \quad \text{for} \quad \rho \gg \langle \rho \rangle,$$
(27)



FIG. 1. In the upper panel the quantity $\frac{\pi_{st}(n,N,M)}{\pi_{st}(0,N,M)}$ of Eq. (28) is plotted against n + 1 in a double logarithmic scale, for M = 100and different values of N (see key). The dotted green line is the small-*n* behavior $(n + 1)^{-k}$. In the lower panel the exact probability $P_{st}(N,M)$ with k = 3 and M = 100, obtained by iteration of Eq. (14) using the form (9) in Eq. (15), is plotted against N with double logarithmic scales. The dashed violet curve is the behavior $(N - \langle N \rangle)^{-3}$ of Eq. (26).

one sees that, upon increasing M, a fluctuation with $\rho \gg \langle \rho \rangle$ is reduced only as M^{1-k} . This gives a much better chance to detect large deviations with respect to a case where the LDP holds.

Condensation can be understood by considering the conditional probability $\pi_{st}(n, N, M)$ [Eq. (15) at stationarity], which is shown in the upper part of Fig. 1. Here π_{st} is plotted for k = 3 and M = 100 and is normalized by $\pi_{st}(n = 0, N, M)$ in order to better compare curves for different values of N. In this figure π_{st} is obtained from Eq. (15) by evaluating P_{st} by means of the recurrence (14).

Let us discuss the properties of π_{st} . Given its meaning, which is expressed below Eq. (15), it is clear that $\pi_{st}(n > N, N, M) = 0$ in all cases, as can be seen in the upper half of Fig. 1. Furthermore, by plugging Eq. (26) into Eq. (15) one has

$$\pi_{st}(n,N,M) \sim \begin{cases} e^{-MR_{st}(\rho-\frac{n}{M})}(n+1)^{-k} & \text{for } N-n \leqslant \langle N \rangle \\ M^{1-k} \left(\rho-\langle \rho \rangle - \frac{n}{M}\right)^{-k}(n+1)^{-k} & \text{for } N-n \gg \langle N \rangle \end{cases}$$
(28)

where we have replaced M - 1 with M and neglected $\frac{1}{M}$ for large M.

This equation shows that, for fixed *N* and *M*, $\lim_{n\to 0} \pi_{st}(n, N, M) \propto (n + 1)^{-k}$ for any value of *N*. However, what makes the big difference between the normal and the condensed case is the behavior of π_{st} at large $n \leq N$. Indeed, for $N \leq \langle N \rangle$ a simple study of Eq. (28) found in Appendix B) shows that $\pi_{st}(n, N, M)$ is a monotonically decreasing function of *n*. Hence, large values of *n* are associated with a very small probability π_{st} , and this implies that condensation, namely, a large fraction of particles in a single box, is probabilistically negligible.

This can be checked in the upper panel of Fig. 1 The cases with $N \leq \langle N \rangle$ discussed above are represented by the curves with N = 10 and N = 20, because using Eq. (10) with k = 3 one finds $\langle N \rangle \simeq 36.84$. These curves decay monotonically, as expected.

Conversely, when condensation occurs, there is an extensive number n_c of particles in a box, meaning that π_{st} must be non-negligible also for values of n as large as $n = n_c$. Indeed, for $N > \langle N \rangle$ (curves with $N \ge 50$), $\pi_{st}(n, N, M)$ develops a pronounced maximum [30], as can be seen in Fig. 1 (upper panel). The location of the peak (see Appendix B) is in

$$n = n_c(N) = \eta \left(N - \langle N \rangle \right), \tag{29}$$

where $\eta \leq 1$ is a function weakly dependent on N and M such that $\eta \rightarrow 1$ as N or M become large (the dependence on N can be checked by inspection of Fig. 1).

The phenomenon of condensation of fluctuations is not restricted to the present model, but it has been observed in a variety of different systems [17,31,32], not only related to physics. Despite different in principle from the usual condensation *on average* occurring in the prototypical example of a boson gas [33] and in many other systems [29,34], these two kinds of condensation are related, as explained in Ref. [31], and a strong mathematical similarity exists.

IV. DYNAMICS

In the stationary state, for large M, a typical observation of the system will give a value N_{obs} of N very close to $\langle N \rangle$. However, if one waits enough, starting from this initial state a fluctuation with $N \neq \langle N \rangle$ will develop spontaneously. The aim of this paper is to describe the properties of the dynamical process associated with the formation of such large deviations. This is particularly interesting for a fluctuation with $N > \langle N \rangle$, since in this case a macroscopic number n_c of particles must pile up in a single box whose occupancy was initially very small, and this might be a slow and complex phenomenon. Furthermore, once such a fluctuation sets on, it must regress and this implies once again the dislocation of a large number n_c of particles. These two processes will be studied in the following sections. In order to do that, we start at t = 0 with the typical form of the single-variable probabilities in a system where the value $N = N_{obs}$ is observed [35]. Recalling the meaning of π [Eq. (15) and discussion below], this reads

$$p(n,t=0) = \frac{\pi_{st}(n, N_{\text{obs}}, M)}{P_{st}(N_{\text{obs}}, M)}.$$
(30)

For instance, we will consider in the following section the case where $N_{obs} = \langle N \rangle$ is the most likely value of N in order to study how a large deviation forms. Then we will consider the evolution of the p's by means of Eq. (11), and, from the knowledge of p(n,t) at all times, we will derive the form of P(N,M,t) and of $\pi(n,N,M,t)$ using Eqs. (14) and (15) and/or Eqs. (16) and (17).



FIG. 2. The single-variable probability p(n,t) with k = 3 is plotted against n + 1 with double logarithmic scales for different times (see key) exponentially spaced. The dotted green line is the asymptotic form (9). In the lower inset data collapse is tested by plotting $\frac{p(n,t)}{p_{st}(n)}$ against $\frac{n+1}{\tilde{v}(t)}$. The quantity $\tilde{v}(t)$ is plotted against *t* in the upper inset, in a double logarithmic plot. The dashed magenta line is the expected long-time behavior (33).

A. Creation of a fluctuation

1. Evolution of the single-variable probabilities p(n,t)

In this case we take $N_{obs} = \langle N \rangle$. The dependence on *n* of the initial condition (30) is contained in the function $\pi_{st}(n, N_{obs} = \langle N \rangle, M)$. As discussed regarding Eq. (28), this quantity behaves as $p(n,0) \simeq (n+1)^{-k}$ for small *n* and goes rapidly to zero for $n \simeq \langle N \rangle$. This can be seen in the upper panel of Fig. 1 (curve with $N = \langle N \rangle$) or in Fig. 2 (leftmost black curve, for t = 0, corresponding to the initial condition (30) with $N_{obs} = \langle N \rangle$). Hence we can write $p(n,0) \simeq p_{st}(n) f[(n + 1)/\nu]$, where $\nu \simeq \langle N \rangle$ and *f* has the properties

$$f(x) \simeq \begin{cases} 1 & \text{for } x \ll 1\\ 0 & \text{for } x \gg 1 \end{cases}$$
(31)

For long times one must approach the stationary condition (9). Therefore, for large t, we search for a *scaling* solution of Eq. (11) of the form

$$p(n,t) = \sigma(t) p_{st}(n) f\left[\frac{n+1}{\nu(t)}\right],$$
(32)

where f(x) has the properties (31), v is an increasing function of t, and σ is a weakly time-dependent normalization such that $\lim_{t\to\infty} \sigma(t) = 1$. Plugging this ansatz into the first line of Eq. (11) and performing the calculations as detailed in Appendix C one can determine the exact form of the scaling function f(x) and of the growth law of v,

$$\nu(t) = bt^{\frac{1}{k+2}},\tag{33}$$

where b is a constant.

The behavior of p(n,t) as a function of *n* at different times, obtained by numerical integration of Eq. (11), is shown in Fig. 2. In the lower inset of this figure we illustrate the data collapse obtained by plotting $\frac{p(n,t)}{p_{st}(n)}$ against $\frac{n+1}{\tilde{v}}$, according to the scaling (32) [recalling that $\sigma(t) \simeq 1$]. The



FIG. 3. The probability P(N, M, t) with k = 3 is plotted against N with double logarithmic scales for different times (see key), exponentially spaced. The dotted green line is the exact asymptotic form given by iteration of Eq. (14), using the form (9) in Eq. (15) (it is the same curve as the green one plotted in the lower panel of Fig. 1). The dashed violet curve is the behavior $(N - \langle N \rangle)^{-3}$ of Eq. (26). In the inset P(N, M, t) is plotted against ρ for $t = 4.85 \times 10^8$ and different choices of M (M = 10, 20, 50, 100, 200, 500 starting from the curve on the top) on a double logarithmic scale. The dashed violet curve is the behavior $(N - \langle N \rangle)^{-3}$.

function $\tilde{v}(t)$ [eventually to be identified with v(t)] has been obtained looking for the best superposition of the curves at different times. As shown in the upper inset, \tilde{v} satisfies the behavior Eq. (33) asymptotically. The collapse of the curves displayed in the lower inset shows some correction at short times, progressively improving with increasing *t*. Furthermore, data fall on a master curve that is almost indistinguishable from the exact form of the scaling function f(x) given in Eq. (C6) of Appendix C, when the undetermined parameter *a* appearing in that expression is appropriately tuned. This confirms the validity of our solution based on the scaling ansatz.

2. Evolution of the collective probability P(N, M, t)

Once the form of the microscopic probabilities p is found at all times, one can obtain the exact time evolution of the probability P of the collective quantity N by inserting p(n,t)in the recurrence equation (14). The outcome of this procedure is shown in Fig. 3, where P(N, M, t) is plotted against N and different curves correspond to different times (see caption). One observes that in the region $N < \langle N \rangle$, where condensation does not occur, the asymptotic form P_{st} (dotted green line) is reached already at an early stage. On the contrary, in the condensing part with $N > \langle N \rangle$ the recovery towards P_{st} is slow and proceeds gradually from small to large values of Nas time goes on. In this way, at any time, no matter how long, there exists a region of sufficiently large values of N where stationarity is not yet reached.

This behavior can be understood analytically. The analysis is presented in Appendix D. The main outcome of this study is that, at any time t, the dynamical probability P(N,M,t) catches up with its stationary value $P_{st}(N, M)$ for densities

$$\rho - \langle \rho \rangle \lesssim \frac{\nu}{M(k-1)}$$
(34)

larger then the average value. Recalling the discussion in Sec. III, this implies that the LDP is violated. Violation of the LDP in the dynamics occurs due to a mechanism analogous to the one operating at stationarity. Indeed, also in the dynamical case, the steepest descent evaluation of the integral in Eq. (12) cannot be done straightforwardly. We also show that, for $\rho > \langle \rho \rangle$ outside the range (34), the validity of the LDP is restored, because the integral in Eq. (12) does admit a saddle-point evaluation. Clearly, this is trivially true in any case also for small densities $\rho \leq \langle \rho \rangle$.

Notice that the interval (34) shrinks to zero as M increases. Hence, for any finite value of v, e.g., at any time, the validity of the steepest descent solution is recovered by considering a number M of boxes sufficiently large. However, if $v = \infty$ (namely, in the stationary state), the saddle-point evaluation fails for any $\rho > \langle \rho \rangle$ and condensation occurs. Let us remark that the above analysis implies that $\lim_{M\to\infty} \lim_{t\to\infty} P(N, M, t) \neq \lim_{t\to\infty} \lim_{M\to\infty} P(N, M, t)$.

This interplay between N (or equivalently ρ) and M is shown in the inset of Fig. 3, where P is plotted against ρ for $t = 4.85 \times 10^8$ (corresponding to the third to last curve in the main picture), for different values of M. Here it is clearly seen that, according to Eq. (34), the region of ρ where P coincides with P_{st} shrinks as M is increased until, at M = 500it is practically absent, meaning that the LDP is recovered basically everywhere. The dashed violet line illustrates the large-N behavior of Eq. (26).

It is clear that in the range of Eq. (34) something akin to condensation occurs, although its mathematical definition is less sharp than in the stationary state, since we cannot let $M \to \infty$ because the interval (34) would shrink to zero. This is supported by the observation that for ranges of ρ increasing with time P(N, M, t) becomes basically indistinguishable from $P_{st}(N,M)$ (see Fig. 3). This implies that in such ranges condensation occurs as at stationarity. For larger values of ρ , however, P departs from P_{st} , signaling that condensation is absent. Since LDP is recovered, P decays exponentially fast in M, like in Eq. (2), as opposed to the much softer algebraic decrease, expressed by Eq. (27). This explains why P(N, M, t)drops off faster than $P_{st}(N,M)$ (violet dashed line).

In order to understand the differences between the dynamical and the stationary state, it is useful to consider the conditional probability $\pi(n, N, M, t)$ of Eq. (15). We have evaluated this quantity at time $t = 4.85 \times 10^8$, which coincides with the time at which we have plotted the third to last indigo curve in Figs. 2 and 3. From Fig. 2 one can infer that $\nu(t) \simeq 120$ at this particular time. Indeed, one sees that p(n,t)(indigo curve) is practically identical to p_{st} (dotted green line) up to this cutoff value of n, above which p decreases much more rapidly then p_{st} .

We have plotted $\pi(n, N, M, t)$ in Fig. 4 (normalized by $\pi(n = 0, N, M, t)$). It is useful to contrast this probability with $\pi_{st}(n, N, M)$ plotted in the upper half of Fig. 1. Figure 4 shows that π behaves similarly to π_{st} for $N \leq v$: Upon increasing N a peak is developed around a value $n = n_c(N)$ [given in Eq. (29)] growing with N. However, while for π_{st} this



FIG. 4. The quantity $\frac{\pi(n,N,M,t)}{\pi(0,N,M,t)}$ for $t = 4.85 \times 10^8$ of Eq. (28) is plotted against n + 1 in a double logarithmic scale, for M = 100 and different values of N (see key). The dotted green line is the small-*n* behavior $(n + 1)^{-k}$.

continues to be true for any value of N, no matter how large, the position of the relative maximum of π saturates around $n \simeq v$. This means that not all the $N - \langle N \rangle$ particles exceeding the average condense, but only a quantity of order v. This *partial condensation* is obviously related to the fact that, for n > v, the microscopic probability p(n,t) rapidly vanishes and the probability to condense more than v balls is negligible.

In conclusion, at a given time t the probability P(N,M,t) has reached the stationary form $P_{st}(N,M)$ only up to a value of ρ given in Eq. (34), while for larger values it is strongly suppressed. Correspondingly, in the range (34) a condensation phenomenon similar to the one observed at stationarity is observed, with a number $n_c(N)$ in Eq. (29) of particles populating a single box. For larger values of N, outside the interval (34), only an incomplete condensation occurs and a reduced number $n_c(v) < n_c(N)$ (with respect to what occurs at stationarity) of particles is accumulated. Notice that the approach of P to P_{st} is a slow, everlasting process, since it is regulated by the power-law growth (33) of v.

B. Regression of a fluctuation

1. Evolution of the single-variable probabilities p(n,t)

In order to study the process of the regression of a large fluctuation we take $N_{obs} \gg \langle N \rangle$. According to Eq. (28) and the following discussion, the *n*-dependence of the initial condition (30) is given by the quantity $\pi_{st}(n, N_{obs}, M)$. This behaves like $p(n,0) \simeq (n+1)^{-k}$ for small *n* and there is a peak at large *n*, centered around the value $n_c(N_{obs})$ of Eq. (29), as shown in the upper part of Fig. 1. The initial condition is represented in Fig. 5 (curve for t = 0, black). We can express these features with the form

$$p(n,0) \simeq p_{st}(n) f\left[\frac{(n+1)}{\nu}\right] + \nu^{-\alpha} g\left[\frac{(n+1)}{\nu}\right], \quad (35)$$

where $\nu \simeq n_c(N_{\text{obs}})$, *f* has the properties (31), and $\nu^{-\alpha}$ and *g* are an amplitude and a function describing the behavior of the condensate (the peak), respectively.



FIG. 5. The single-variable probability p(n,t) with k = 3 is plotted against n + 1 with double logarithmic scales for different times (see key), exponentially spaced. The dotted green line is the asymptotic form (9). In the upper inset data collapse is tested by plotting $q(t)[p(n,t) - p_{st}(n)]$ against $x = \frac{n+1}{\tilde{v}(t)}$. The dashed indigo line is the behavior x^4 . In the lower inset the quantities $\tilde{v}(t)$ (below) and q(t) (above) are plotted against t in a double logarithmic plot. The dashed magenta line (below) is the expected behavior (33), and the dashed turquoise line (above) is the behavior $\sim t^{3/2}$.

200 165

n+1

10

Proceeding as in Sec. IV A 1, for large t we search for a *scaling* solution of the form

$$p(n,t) = \sigma(t)p_{st}(n)f\left[\frac{n+1}{\nu(t)}\right] + \nu^{-\alpha}(t)g\left[\frac{n+1}{\nu(t)}\right], \quad (36)$$

where ν , σ , and f have the same meaning as in Sec. IV A 1, α is a dynamical exponent and the scaling function g has the following limiting behaviors:

$$\lim_{x \to 0} g(x) = 0$$

$$\lim_{x \to \infty} g(x) = 0.$$
(37)

Inserting the form (36) into the first line of Eq. (11) and proceeding like in Appendix C one has the time dependence (33) of v together with the form (C6) of f(x). The expression for g(x) can also be determined. This is detailed in Appendix E.

The behavior of p(n,t) as a function of n at different times, obtained by numerical integration of Eq. (11), is shown in Fig. 5. According to Eq. (36) a superposition of curves at different times should be obtained by plotting $\nu^{\alpha} \{p(n,t) - \sigma(t)p_{st}(n)f[\frac{n+1}{\nu}]\}$ against $\frac{n+1}{\nu}$. Given that $\sigma(t) \simeq 1$ for long times and $f(x) \simeq 1$ for $x \lesssim 1$, data collapse can be checked by plotting $\nu^{\alpha} \{p(n,t) - p_{st}(n) = 1\}$ as well.

Since the value of α cannot be obtained from the above calculation, we evaluate it from the numerical data as follows: The second term in Eq. (36) describes the peak of p observed in Fig. 5. The amplitude $v^{-\alpha}$ of such contribution at the peak position $n_c(t)$ can then be estimated by measuring the difference $q^{-1}(t) = p(n_c,t) - \sigma(t)p_{st}(n_c)f[\frac{n_c+1}{v(t)}]$. For large times it is $\sigma(t) \simeq 1$, $n_c(t) \simeq v(t)$, and the first term in Eq. (36) behaves as $p_{st}(n)$ up to $n \simeq v(t) \simeq n_c(t)$ (given the form of the function f; see inset of Fig. 2). Therefore the quantity $q^{-1}(t)$ can be approximately simplified to $q^{-1}(t) \simeq p(n_c,t) - p_{st}(n_c)$.

100

1000

In the upper inset of Fig. 5 we plot $q(t)\{p(n,t) - p_{st}(n)\}$ against $x = \frac{n+1}{\tilde{v}(t)}$, where $\tilde{v}(t)$ is defined like in Sec. IV A 1. As pointed out above, data collapse of the curves at different times is expected in this plot in the region of the peak. In this figure, in fact, an excellent superposition is found for long times. Notice that, using the small-z behavior $L_n^{\lambda}(z) \simeq \text{const}$ of the Laguerre polynomials entering the form of g [Eq. (E2) in Appendix E], one has $g(x) \sim x^{k+1}$ for small x, which is indeed very well observed in the upper inset of Fig. 5. The behavior of v(t) and $q(t) \sim v(t)^{\alpha} \sim t^{\frac{\alpha}{k+2}}$ as time changes is shown in the lower inset of the figure. This plot confirms the growth law (33) of v and indicates a value of α consistent with $\alpha = 15/2$ for the case considered here with k = 3.

2. Evolution of the collective probability P(N, M, t)

Equation (36) shows that the single-variable probability is roughly the one at stationarity with a cutoff at $n + 1 \simeq v(t)$ and an extra contribution (the second term on the right-hand side) concentrated around $n + 1 \simeq v(t)$. The latter, which represents the condensed fraction, is clearly visible as a *bump* in Fig. 5. Given this form of p(n,t) it is easy to show that the collective probability P exhibits a series of maxima located in $N = N_{\ell}(t)$ $(\ell = 0, 1, 2, ...)$ given by

$$N_{\ell}(t) = \langle N \rangle + \ell \nu(t). \tag{38}$$

Indeed, there is an obvious maximum of the probability in the average value, $N_0 = \langle N \rangle$. Furthermore, since the *p*'s have a large support around ν , the situation where, besides the $\langle N \rangle$ particles distributed among the boxes, there are ν marbles stored in a single box (or, in general, in ℓ boxes) is also largely probable, thus giving a series of maxima located like in Eq. (38).

This structure of *P* with many relative maxima is shown in Fig. 6. Also in this case *P* has been computed by inserting the time-dependent form of the *p*'s in Eq. (14). Clearly, as time goes on and ν increases the relative *strength* of the condensed term decreases (as $\nu^{-\alpha}$) and the maxima are gradually smeared out. The location of the maxima (38) can be checked by



FIG. 6. The probability P(N,M,t) with k = 3 is plotted against N (with double logarithmic scales) for different times (see key), exponentially spaced. The dotted green line is the asymptotic form (26). In the inset the same quantity is plotted against $\frac{N-\langle N \rangle}{\nu(t)}$.

plotting *P* against $\frac{N-\langle N \rangle}{\nu(t)}$, since on this axis the maxima are placed on the integer values $\ell = 0, 1, 2, \ldots$. This is very neatly observed in the inset of Fig. (6). From the discussion above it is clear that the maxima with $\ell > 0$ are due to the presence of the second term on the right-hand side of Eq. (36). In the region $N < N_1(t)$, since the effect of this term is negligible, one recovers the asymptotic behavior $P(N,M,t) \simeq P_{st}(N,M)$ of Eq. (26), as can be seen in Fig. 6.

By computing $\pi(n, N, M, t)$ one finds a behavior analogous to the one discussed in Sec. IV A 2, signaling that also in this case full condensation can occur only for sufficiently small values of N, whereas only a partial one is possible for larger N. This is due to the same mechanism already discussed in Sec. IV A 2, namely, to the fact that the microscopic probabilities p are negligibly small for $n > \nu$. Again, the approach of P to the stationary form is an everlasting slow evolution.

V. SUMMARY AND CONCLUSIONS

We have investigated the kinetics leading to the formation or to the resorption of a large fluctuation of a collective variable N in a statistical system. We have considered a simple model where N is the sum of a relatively large number *M* of stochastic (micro-) variables n_m (m = 1, ..., M) identically and independently distributed. We speak of nballs stored in a box with probability p(n), to make the idea more concrete. The evolution equation of the p's is chosen as to have a stationary solution p_{st} with fat tails, namely $p_{st}(n) \propto (n+1)^{-k}$ (k > 2). It is known that this form induces a probability $P_{st}(N,M)$ to observe a given value of N at stationarity that does not obey the LDP for $N > \langle N \rangle$, due to a condensation phenomenon. This feature implies that large fluctuations in this range have a better chance to be observed in a system with finite but large M with respect to a case where the LDP holds. This possibly makes some of our results amenable to numerical and/or experimental verification.

We have considered the evolution of the model starting from (1) a typical situation where the most probable value $\langle N \rangle$ is observed and (2) a case where a measurement with unlikely outcome $N > \langle N \rangle$ just happened. In (1) we follow the dynamical process whereby large fluctuations, which are not present in the *average* initial state, form. Conversely, in (2) we consider a realization where a rare initial state is disrupted upon approaching stationarity. Both these cases can be solved analytically. We have worked out the solutions in detail and reported them in the appendices.

We have shown that the convergence to stationarity is, in any case, a slow, everlasting process akin to those observed in aging systems. This happens because the evolution is slaved by the power-law growth of a characteristic value N = v(t) of N separating a region N < v(t), where the stationary behavior has been attained, from one with N > v(t), where reminiscence of the initial condition is retained. During such evolution one observes a condensation phenomenon with a twofold character: For $N \leq v(t)$ the phenomenon is indistinguishable from the one observed at stationarity, with the huge number $n_c(N)$ (29) of particles stored in a single box. For N > v(t), instead, condensation is incomplete: a relevant number of balls is still accumulated, but this number only equals $n_c(v)$, which is smaller than the value expected when full condensation occurs. As time passes and v(t) diverges, full condensation is gradually recovered on increasing values of N.

Related to that, the interval I where the LDP breaks down has a nontrivial time dependence. We stress again that in this sector large deviations may occur more easily. To assess I is therefore of practical interest in nanoscale applications where fluctuations play an important role. In the stationary state, full condensation happens, and this gives $\mathcal{I} \equiv \{\rho | \rho > \langle \rho \rangle\}.$ However, at finite times, since condensation is incomplete, the LDP is spoiled only in the region given by Eq. (34). This means that, at any time, for a finite M the LDP is restored for sufficiently large values of ρ , at variance with what happens at stationarity. The region (34) expands as time elapses and v(t) increases. In this way the violation of the LDP progressively extends towards the whole sector with $\rho > \langle \rho \rangle$, and the stationary properties of P are recovered. Equation (34) shows that the size of \mathcal{I} can be tuned not only by changing t, but also by acting on M. Since t and M enter in the combination $\nu(t)/M$, this implies that $t \to \infty$ and $M \to \infty$ are noncommuting limits.

The simple probabilistic setup discussed in this paper is suited to describe at an elementary level the dynamics of fluctuations in a variety of systems ranging from physics to chemistry, biology, and the social sciences. It has the advantage of being amenable to analytic investigation. Some of the general features displayed here rely only on general aspects of probability, such as the violation of the LDP. Therefore, they are expected to be observed with similar characteristics in a class of problems wider than the one considered in this paper, e.g., with nonidentically and/or nonindependently distributed microscopic variables. This makes the issue a rather broad and general research topic worth of further investigations. Finally, we remark that collective variables defined differently from the sum N, e.g., energy or heat fluxes in solvable models of statistical mechanics, have been shown [31] to display a different condensation phenomenon, without violation of the LDP. The question then arises of how the dynamics of fluctuations in these systems compares with the case considered in the present paper. This investigation will be the subject of future work.

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APPENDIX A: COMPUTATIONAL COMPLEXITY OF EQS. (14) AND (12)

Suppose we want to determine *P* by means of Eq. (14). Let us consider the procedure at a certain step when *P* is known for a number m - 1 of boxes and the task is to determine it for *m* of them. Equation (14) says that, in order to find P(N,m,t), we must preliminarily know $\pi(n,N,m,t)$ for any value of *n*. It is obvious that in order to find $\pi(n,N,m,t)$ (for any *n*) by means of Eq. (15) the value of P(j,m - 1,t) must be previously known for any j = 0, 1, ..., N. This means that at any step $\pi(n,i,m - 1,t)$ must be determined through Eq. (15) for any n = 0, 1, ..., N and i = 0, 1, ..., N. This requires $(N + 1)^2$ computations. Once $\pi(n,i,m,t)$ is known in this way, we can get P(j,m,t) through Eq. (14) with further $(N + 1)^2$ computations, since we have to sum up N + 1 terms, and this operation must be repeated for any j = 0, 1, ..., N. Then, for any step (namely, going from m - 1 to m) a number of order $2(N + 1)^2$ of elementary computations is needed. Since the recurrence must be repeated up to m = M, a total number $2(N + 1)^2M$ of such calculations is needed. This must be compared with the exponentially large number $(N + 1)^M$ of operations involved in the determination of P by using the first line of Eq. (12).

APPENDIX B: PROPERTIES OF π_{st}

For $N \leq \langle N \rangle$ the first line of Eq. (28) necessarily applies, which shows that not only $(n + 1)^{-k}$ decreases upon increasing *n* but also $e^{-MR_{st}(\rho-\frac{n}{M})} = P_{st}(N-n,M)$. Referring to Fig. 1 (lower part), this can be understood as follows: In the first line of Eq. (28), $P_{st}(x, M)$ is evaluated for x = N - n. Recalling that the condition $N - n \leq \langle N \rangle$ applies, this value of x is located on the left of the maximum of P_{st} located in $x = \langle N \rangle$ (or, at most, on the maximum itself). Consequently, raising nmoves the argument x of $P_{st}(x, M)$ further and further away on the left of the maximum (which amounts to descend towards the left along the green curve of the lower panel of Fig. 1). This makes $P_{st}(x, M)$ decrease monotonically. The quantity $e^{-MR_{st}(\rho-\frac{n}{M})}$ in the first line of Eq. (28) behaves similarly. Then large values of *n* are associated with a very small probability π_{st} , and this implies that condensation, namely, a large fraction of particles in a single box, is probabilistically negligible.

Conversely, for $N \gg \langle N \rangle$ and values of *n* such that the lower row of Eq. (28) applies [i.e., for $N - n \gg \langle N \rangle$], while increasing *n* the term $(n + 1)^{-k}$ decreases, the factor $(\rho - \langle \rho \rangle - \frac{n}{M})^{-k}$ increases. The effect of this is the development of a pronounced maximum [30] in $\pi_{st}(n, N, M)$, as can be seen in Fig. 1 (upper panel) for $N > \langle N \rangle$. There is therefore a relatively high probability of having a macroscopic, namely, of order *N*, number n_c of particles condensed in a single box.

Using the second line of Eq. (28) the location of the maximum is at

$$n = n_c(N) = \frac{1}{2} \left(N - \langle N \rangle \right), \tag{B1}$$

where e = 1/2 is a constant. Notice that the simple calculation presented above to determine n_c is not exact, since the second line of Eq. (28) is accurate only for $N - n \gg \langle N \rangle$ while n_c is located outside this range. It can be shown, however, that the result (B1) is basically correct, since the true behavior, expressed by Eq. (29), differs only by the value of the prefactor.

APPENDIX C: CREATION OF A FLUCTUATION: SOLUTION OF THE EQUATION FOR THE EVOLUTION OF THE *p*'S

Inserting Eq. (32) into the first line of Eq. (11) one arrives at

$$v(t)^{k-1} \frac{dv(t)}{dt} x^{k+1} \left[\frac{v \frac{d\sigma}{dt}}{\sigma \frac{dv}{dt}} \frac{f(x)}{x} - f'(x) \right]$$
$$= -\left[\left(1 + \frac{1}{\nu x} \right)^{-k} + \left(1 - \frac{1}{\nu x} \right)^{-k} \right] f(x)$$
$$+ \left[\left(1 + \frac{1}{\nu x} \right)^{-k} f\left(x + \frac{1}{\nu} \right) \right]$$

$$+\left(1-\frac{1}{\nu x}\right)^{-k}f\left(x-\frac{1}{\nu}\right)\Big],\tag{C1}$$

where $x = (n + 1)/\nu$ like before, and f' = df/dx. Now we make the ansatz that the quantity $\frac{\nu \frac{da}{dt}}{\sigma \frac{d\nu}{dt}}$ vanishes in the long-time limit. This will be checked for consistency at the end of the calculation. In the same limit, when ν is large, we can expand the terms $(1 \pm \frac{1}{\nu x})^{-k} \simeq 1 \mp \frac{k}{\nu x} + \frac{k(k+1)}{2\nu^2 x^2}$ and $f(x \pm \frac{1}{\nu}) \simeq f(x) \pm \frac{1}{\nu} f'(x) + \frac{1}{2\nu^2} f''(x)$ to second order in the small quantity $1/(\nu x)$, and retaining the leading terms one obtains

$$-\nu(t)^{k+1}\frac{d\nu(t)}{dt}x^{k+1}f'(x) = f''(x) - \frac{2k}{x}f'(x).$$
 (C2)

Regarding Eq. (C2) in the variables x, t, since the right-hand side does not depend on t one must have, on the left-hand side, $v^{k+1}dv/dt = a$, where a > 0 is a constant. Hence

$$\nu(t) = bt^{\frac{1}{k+2}},\tag{C3}$$

where $b = [a(k+2)]^{\frac{1}{k+2}}$. Equation (C2) then reads

$$f''(x) + \left(ax^{k+1} - \frac{2k}{x}\right)f'(x) = 0,$$
 (C4)

which, with the limiting behaviors (31), has the solution

$$f'(x) = -cx^{2k}e^{-\frac{a}{k+2}x^{k+2}},$$
 (C5)

where c > 0 is a constant. Integrating once again one arrives at

$$f(x) = \frac{\Gamma\left(-\frac{3}{k+2}, \frac{ax^{k+2}}{k+2}\right)}{\Gamma\left(-\frac{3}{k+2}\right)} - \frac{u\left(1 + \frac{k+2}{k^2+k-2}x^{k+2}\right)}{x^3}e^{-\frac{a}{k+2}x^{k+2}},$$
(C6)

where $\Gamma(\alpha)$ and $\Gamma(\alpha, y)$ are the Γ and the incomplete Γ function, and $u = \frac{\left[\frac{(k+2)^{k+5}}{a^3}\right]^{\frac{1}{k+2}}}{3\Gamma(-\frac{3}{k+2})}$. This form depends on the single parameter *a*, which is difficult to determine since our solution is exact only asymptotically. The quantity $\sigma(t)$ in Eq. (32) can be easily obtained from the normalization of the probability as

$$\sigma(t) - 1 \propto \nu(t)^{1-k}.$$
 (C7)

Finally, from this equation and Eq. (C3) it is easy to verify the ansatz made after Eq. (C1), namely, that $\frac{v \frac{d\sigma}{dt}}{\sigma \frac{dv}{dt}} \to 0$ for $t \to \infty$.

APPENDIX D: BEHAVIOR OF THE COLLECTIVE PROBABILITY P

Equation (32) shows that the form of the single-variable probability p is basically the one at stationarity with a cutoff at $n + 1 \simeq v(t)$. We simplify the discussion about the evolution of P by assuming that such cutoff is sharp. This amounts to approximate the actual behavior of f given by Eq. (C6) with the schematic form $f(x) = 1 - \theta(x - 1)$. Reparametrizing time in terms of v by means of Eq. (33), let us study the lefthand side, $S(z,v) = z \frac{Q'(z,v)}{Q(z,v)}$ of the saddle-point equation (18), for arbitrary z. Since now

$$Q(z,v) = \sum_{n=0}^{v} p_{st}(n) z^n = \zeta^{-1}(k) \sum_{n=0}^{v} (n+1)^{-k} z^n, \quad (D1)$$

one has

$$S(z,\nu) = \frac{\sum_{n=0}^{\nu} n(n+1)^{-k} z^n}{\sum_{n=0}^{\nu} (n+1)^{-k} z^n}.$$
 (D2)

Notice that, in the two equations above, the notation should be specified since *n* must run up to an integer number, e.g., the closest to v. However, this would not change the discussion below, and we prefer to keep the simple notation of Eqs. (D1) and (D2). As a function of *z*, *S* rises steeply from S(z = 0, v) = 0 to an asymptotic value (for large *z*) that can be easily determined by retaining only the dominant term with n = v in the sums defining *S* in Eq. (D2):

$$\lim_{z \to \infty} S(z, \nu) = \nu.$$
 (D3)

This asymptotic value is assumed for $z \gtrsim z_r$, where z_r can be evaluated as follows. Let us consider the numerator on the right-hand side of Eq. (D2). For $z \gtrsim 1$, as a function of *n*, the terms $n(n + 1)^{-k} z^n$ decrease down to a value n_r given by the largest solution of the following equation:

$$\ln z = \frac{k}{n_r + 1} - \frac{1}{n_r}.$$
 (D4)

For $n > n_r$, the argument of the sum in the numerator of Eq. (D2) very rapidly diverge, because of the term z^n . A similar analysis for the denominator shows that it behaves similarly, but with a slightly larger value of n_r , that we denote by $n_{r,D}$, given by $\ln z = \frac{k}{n_{r,D}+1}$. Starting from $n_r = \infty$ when z = 1, n_r decreases upon raising z. Notice that, if z is too close to unity, one has $n_r > v$, meaning that such value is not contained in the sums defining S in Eq. (D2). Therefore a critical value z_r exists,

$$\ln z_r = \frac{k}{\nu + 1} - \frac{1}{\nu},$$
 (D5)

such that, for $z > z_r$ this term starts to be contained in the sums in Eq. (D2), and, beyond that value, these sums rapidly diverge. A similar analysis carried out for the denominator of Eq. (D2) leads to a value, denoted by $z_{r,D}$, given by $\ln z_{r,D} = \frac{k}{\nu+1}$. Since $z_r < z_{r,D}$, $S(z,\nu)$ is a very steep function for values of z close to z_r , and then flattens when z also crosses $z_{r,D}$ and the denominator diverges as well. As Eq. (D5) shows, $z_r \ge 1$ for large ν , which allows one to expand $\ln z_r$ around $z_r = 1$ on the left-hand side of Eq. (D5) and to set $\nu + 1 \simeq \nu$ on the right-hand side, thus arriving at

$$z_r = 1 + \frac{k-1}{\nu}.$$
 (D6)

This means that when the solution z^* of Eq. (18) moves the short distance from $z^* = 1$ to the nearby value $z^* = z_r$ given by Eq. (D6), ρ widely varies from $\rho = \langle \rho \rangle$ [according to Eq. (10)] to a value $\rho \sim \nu$ since, because of Eq. (D3), the left-hand-side *S* of the saddle-point equation (18) rapidly converges to this value as soon as $z \gtrsim z_r$.

Let us now consider the argument $-M\mathcal{R}(z,\rho,\nu)$ of the exponential defining *P* on the right-hand side of Eq. (12). Given that the largest contribution to the integral comes from $z \simeq z^*$ and, for $\langle \rho \rangle \leq \rho \lesssim \nu, z^*$ is close to $z^* = 1$, in this range

of ρ one can write

$$M\mathcal{R}(z,\rho,\nu) \simeq -M\left[\frac{d\ln Q}{dz}\bigg|_{z=1}(z-1) - \rho(z-1)\right]$$
$$= M(\rho - \langle \rho \rangle)(z-1), \tag{D7}$$

where we have expanded the argument of the exponential in Eq. (12) to first order in z - 1 and we have used $\ln Q(1) = 0$ [after Eq. (13) and normalization of the probabilities p] and $\frac{d \ln Q}{dz}|_{z=1} = \langle \rho \rangle$ [from Eqs. (18) and (10)].

Roughly speaking, for a given value of M, a saddle-point evaluation of the integral defining P in Eq. (12) is accurate if the positive quantity of Eq. (D7), as a function of z, has a pronounced minimum at $z = z^*$. This in turn means that, whatever the value of $M\mathcal{R}$ at $z = z^*$ is, it must become much larger than its value for other choices of z. However, we know that, in the region $\rho \ge \langle \rho \rangle$ where condensation is possible, z^* ranges at most up to $z^* = z_r$. This implies that a saddle-point solution cannot be invoked if $M(\rho - \langle \rho \rangle)(z_r - 1)$ is not a large number, namely, if $\rho - \langle \rho \rangle < [M(z - 1)]^{-1}$. Using Eq. (D6), this means that the steepest descent evaluation breaks down for all the densities

$$ho - \langle \rho \rangle \lesssim rac{
u}{M(k-1)}$$
 (D8)

larger but sufficiently near to the average value. Its validity is restored only for larger values of ρ (besides, clearly, for small densities $\rho \leq \langle \rho \rangle$).

APPENDIX E: REGRESSION OF A FLUCTUATION: SOLUTION OF THE EQUATION FOR THE EVOLUTION OF THE *p*'S

Inserting the form (36) into the first line of Eq. (11) and proceeding like in Appendix C one has the time dependence (33) of ν and the form (C6) of f(x), whereas for g(x) one arrives at

$$g''(x) + ax^{k+1}g'(x) + \left[a\alpha x^k - \frac{k(k+1)}{x^2}\right]g(x) = 0, \quad (E1)$$

where $a = v^{k+1} dv/dt > 0$ is the same constant introduced in Appendix C [below Eq. (C2)]. With the boundary conditions (37) the solution is

$$g(x) = d x^{k+1} L_{-\frac{\alpha+k+1}{k+2}}^{\frac{2k+1}{k+2}} \left(-\frac{a}{k+2} x^{k+2} \right),$$
(E2)

where *d* is a constant and $L_m^{\lambda}(z)$ are the generalized Laguerre polynomials.

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- [35] Notice that our initial condition p(n,t = 0) is the typical form of the *p*'s when the fluctuating variable *N* takes the value $N = N_{obs}$, but having the microscopic probabilities set to p(n,t = 0)does not necessarily imply that the outcome of a measurement of *N* is $N = N_{obs}$, since obviously fixing p_{st} we do not constrain the value of *N*.