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Statistical mechanics of a discrete nonlinear system

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Statistical mechanics of the discrete nonlinear Schrödinger equation is studied by means of analytical and numerical techniques. The lower bound of the Hamiltonian permits the construction of standard Gibbsian equilibrium measures for positive temperatures. Beyond the line of \( T = \infty \), we identify a phase transition, through a discontinuity in the partition function. The phase transition is demonstrated to manifest itself in the creation of breather-like localized excitations. Interrelation between the statistical mechanics and the nonlinear dynamics of the system is explored numerically in both regimes.

The pioneering studies of Fermi, Pasta and Ulam (FPU) showed that energy exchange between coupled systems may be suppressed in the presence of nonlinearity; instead a type of behavior that severely contrasts equipartition among the linear modes is observed. The question of whether equipartition of excitation energy always appears is a contemporary issue in various fields of physics. Many manifestations of nonequilibrium and non-equipartition phenomena equivalent to the dynamical behavior of systems with few degrees of freedom contrasting statistical mechanics expectations have been observed. Some of these phenomena, and therefore the absence of immediate equipartition expressed in terms of self-trapping of energy, play an important role for optical storage patterns in nonlinear fibers, condensed matter physics, and biophysics.

A particularity of discrete nonlinear systems is their ability to sustain strong localization of energy. This is accomplished via intrinsic localized modes (breathers) which are modes that remain stable for extremely long times. So far it is a largely unaddressed problem how to handle and describe these excitations in a statistical mechanics framework although it has been argued that breathers may act as virtual bottlenecks delaying the thermalization process.

In this work, we develop a statistical understanding of the dynamics, including the breathers, in a discrete nonlinear Schrödinger (DNLS) equation. The DNLS equation plays a significant role in several branches nonlinear physics as a simple physical model because it may approximate many of the above mentioned nonlinear systems. We study analytically and numerically the thermalization of the lattice for \( T \geq 0 \). We identify the regime in phase space wherein regular statistical mechanics considerations apply, and hence, thermalization is observed numerically and explored analytically using regular, grand-canonical, Gibbsian equilibrium measures. However, the nonlinear dynamics of the problem renders permissible the realization of regimes of phase space which would formally correspond to “negative temperatures” in the sense of statistical mechanics. The novel feature of these states is that the energy tends to be localized in certain lattice sites forming breather-like excitations. Returning to statistical mechanics, such realizations, which would formally correspond to negative temperatures, are not possible (since the Hamiltonian is unbounded from above, as is seen by a simple scaling argument similar to the continuum case) unless one refines the grand-canonical Gibbsian measure to correct for that. This correction will necessitate a discontinuity in the partition function signaling a phase transition that we identify, numerically, with the appearance of breather modes.

In order to explore and illustrate the scenario described above we consider the one-dimensional DNLS equation in the form

\[
 i\dot{\psi}_m + (\psi_{m+1} + \psi_{m-1}) + \nu |\psi_m|^2 \psi_m = 0 , \tag{1}
\]

where the overdot denotes time derivative, \( m \) is a site index, and \( \nu \) is a tunable coefficient to the nonlinear term. Equation (1) is the equation of motion, \( \dot{\psi}_m = -\frac{\partial H}{\partial \psi_m} \), where \( H \) is the Hamiltonian function given by

\[
 H = \sum_m (\psi_m^* \psi_{m+1} + \psi_m \psi_{m+1}^*) + \sum_m \frac{\nu}{2} |\psi_m|^4 ,
\]

for which \( i\psi_m^* , \psi_m \) form canonically conjugate pairs of variables. In addition to the conserved energy \( H \), the quantity \( A = \sum_m |\psi_m|^2 \), is also conserved by the dynamics of Eq. (1), and serves as the norm of the system.

In order to study the statistical mechanics of the system, we calculate the classical grand-canonical partition function \( Z \). We first apply the canonical transformation...
\[ \psi_m = \sqrt{A_m} \exp(i \phi_m), \]
leading to
\[ \mathcal{H} = \sum_m 2 \sqrt{A_m A_{m+1}} \cos(\phi_m - \phi_{m+1}) + \frac{\nu}{2} \sum_m A_m^2. \]

The partition function then becomes,
\[ Z = \int_0^\infty \prod_m dA_m \exp[-\beta(\mathcal{H} + \mu A)], \quad (2) \]
where the multiplier \( \mu \) is introduced in analogy with a chemical potential to ensure conservation of \( A \). Straightforward integration over the phase variable \( \phi_m \) reduces the symmetrized partition function to,
\[ Z = (2\pi)^N \int_0^\infty \prod_m dA_m I_0(2\beta \sqrt{A_m A_{m+1}}) \times \exp \left[ -\beta \sum_m \left( \frac{\nu}{4} (A_m^2 + A_{m+1}^2) + \frac{\mu}{2} (A_m + A_{m+1}) \right) \right]. \]

This integral can be evaluated exactly in the thermodynamic limit of a large system \( (N \to \infty) \) using the eigenfunctions and eigenvalues of the transfer integral operator \( \mathcal{S} \),
\[ \int_0^\infty dA_m \kappa(A_m, A_{m+1}) y(A_m) = \lambda y(A_{m+1}), \]
where the kernel \( \kappa \) is
\[ \kappa(x, z) = I_0(2\beta \sqrt{xz}) \times \exp \left[ -\beta \left( \frac{\nu}{4} (x^2 + z^2) + \frac{\mu}{2} (x + z) \right) \right]. \quad (3) \]

Similar calculations have been performed for the statistical mechanics of the \( \phi^4 \) field \( \mathcal{S} \), and for models of DNA denaturation \( \mathcal{S} \). One obtains \( Z \approx (2\pi \lambda_0)^N \), as \( N \to \infty \) where \( \lambda_0 \) is the largest eigenvalue of the operator. From this expression the usual thermodynamic quantities such as the free energy, \( F \), or specific heat can be calculated. More importantly, for our purpose we can obtain the averaged energy density, \( h = \langle \mathcal{H} \rangle / N \), and the average excitation norm, \( a = \langle A \rangle / N \) as
\[ a = \frac{1}{\beta \lambda_0} \frac{\partial \lambda_0}{\partial \mu} \quad h = -\frac{1}{\lambda_0} \frac{\partial \lambda_0}{\partial \beta} - \mu a. \]

The average excitation norm \( a \) can also be calculated as \( a = \frac{1}{2} \int_0^\infty \prod_m dA_m A_m \exp[-\beta(\mathcal{H} + \mu A)] \), where the integral again can be calculated using the transfer technique \( \mathcal{S} \) and yields \( a = \int_0^\infty y_0^2(A) A dA \), where \( y_0 \) is the normalized eigenfunction corresponding to the largest eigenvalue \( \lambda_0 \) of the kernel \( \kappa \) (Eq. \( \mathcal{S} \)). This shows that
\[ p(A) = y_0^2(A) \]
is the probability distribution function (PDF) for the amplitudes \( A \).

The problem is now reduced to finding the largest eigenvalue \( \lambda_0 \) and the corresponding eigenfunction \( y_0 \) of the transfer operator, Eq. \( \mathcal{S} \). This we do numerically. However, two limits \( (\beta \to \infty \text{ and } \beta \to 0) \) are also amenable to analytical treatment.

First, notice that the Hamiltonian is bounded from below and one can observe that this minimum is realized by a plane wave, \( \psi_m = \sqrt{A} \exp(im\pi) \), whose energy density is \( h = -2a + \frac{\nu}{4} a^2 \). This relation defines zero temperature, or the \( \beta = \infty \) line.

The high temperature limit is also tractable. When \( \beta \ll 1 \) the modified Bessel function in the transfer operator can be approximated, to leading order, by unity (this amounts to neglecting the coupling term in the Hamiltonian). This allows us to reduce the remaining eigenvalue problem to the approximate solution valid for thermalized independent units,
\[ y_0(A) = \frac{1}{\sqrt{\lambda_0}} \exp \left[ -\frac{\beta}{4} (\nu A^2 + 2\mu A) \right]. \]

Using this approximation we can, enforcing the constraint \( \beta \mu = \gamma \) (where \( \gamma \) remains finite as we take the limits \( \beta \to 0 \) and \( \mu \to \infty \)), obtain \( h = \nu / \gamma^2 \) and \( a = 1 / \gamma \). Thus, we get \( h = va^2 \) at \( \beta = 0 \).

FIG. 1. Parameter space \((a, h)\), where the shaded area is inaccessible. The thick lines represents the \( \beta = \infty \) and \( \beta = 0 \) lines and thus bound the Gibbsian regime. The dashed line represents the \( h = 2a + \frac{\nu}{4} a^2 \) line along which the reported numerical simulations are performed (pointed by the symbols).
density \( a \) above the \( T = 0 \) line in an infinite system.

A statistical treatment of the remaining domain of parameter space can be accomplished introducing formally negative temperatures. But the partition function \( \beta, \mu' \) is not suited for that purpose since the constraint expressed in the grand-canonical form fails to bound the Hamiltonian from above. In all the alternative approaches of the study of negative temperatures we will have to consider a finite system of size \( N \). As suggested in \cite{4,10} we can still consider the grand-canonical ensemble using the modified partition function \( Z'(\beta, \mu') = \int \exp(-\beta(\mathcal{H} + \mu'A^2)) \prod_m d\psi_m d\psi^*_m \); but this introduces long range coupling and \( \mu' \) will have to be of order \( 1/N \). Now \( \beta \) can be negative since \( \mathcal{H} + \mu'A^2 \) can be seen to be bounded from above when \( \mu' < -\nu/2N \).

The important consequence of this explicit modification of the measure, is a jump discontinuity in the partition function, that in turn signals a phase transition. More explicitly, if one starts in a positive \( T \), thermalizable (in the Gibbsian sense) state in phase space with \( \beta, \mu \), respectively. Thus, the number of microstates \( Z(\beta, \mu) \) can be negative since \( \mathcal{H} + \mu'A^2 \) is really achieved and whether we can observe different kinds of initial conditions (with same energy and norm densities) produced the same results. In conclusion, the system reaches an equilibrium state which is perfectly recovered by means of the transfer operator method. Moreover it can be checked on Fig. 2 that the curvature of \( \log p(A) \) (i.e., \( -\beta \)) tends to zero when \( h = \nu a^2 \). (The cut-off at high amplitudes is due to finite size effects). In this domain of parameter space, high amplitude excitations are highly improbable and can be considered as simple fluctuations; as shown on Fig. 2 large amplitude fluctuations have been recorded but were checked in the evolution pattern to disappear rapidly.

The scenario is very different when the energy and norm densities are above the \( \beta = 0 \) line. We can observe a rapid creation of breather excitations due to the modulational instability accompanied by thermalization of the rest of the lattice. Once created, these localized excitations remain mostly pinned and because the internal frequency increases with amplitude their coupling with the small amplitude radiation is very small. This introduces a new time scale in the thermalization process necessitating simplectic integration for as long as \( 10^6 - 10^7 \) time units in order to reach a stationary PDF. This can also be qualitatively justified by the effective long range interactions, introduced in the modified partition function, which will produce stronger memory effects as one

\[
\text{FIG. 2. Distribution of } A = |\psi|^2 \text{ for three cases under (and on) the transition line. The solid lines show the results of simulations and the symbols are given by the transfer operator. Curves are vertically shifted to facilitate visualization.}
\]
observes regimes in phase space which are further away from the transition line (since the long range interaction will be stronger).

Typical distribution functions of the amplitudes are shown in Fig. 3. The presence of high amplitude excitations is directly seen here (more straightforwardly we observed standing breathers in the spatial pattern). The dotted line represents the PDF in the case where the initial condition is chosen at random, using a larger system size: we check that the initial condition seems unimportant, but the system size does influence the amplitude of the highest breathers. The cut-off value in the very large system limit, as well as the persistence of a bump in the PDF, is still an open question, since we have no prediction for the PDF above the $\beta = 0$ line. However the positive curvature of the PDF at small amplitudes clearly indicates that the system evolves in a regime of negative temperature and the appearance of the phase transformation is signified in the dynamics by the appearance of these strongly localized persistent breathers.

The actual dynamics in the negative temperature regime is studied more closely in Ref. [11].

Finally, we can draw an interesting parallel with what has been known in plasma physics and hydrodynamics for several years [12], where the appearance of localized structures (of vortices in that case) is also related to a description in terms of negative temperatures.

In conclusion, studying the DNLS system, we have been able to quantify and explain, through analytical calculations supported by numerical computations, the behavior in different regimes of the $(h,a)$ phase space. We have been able to link the regime of thermalization to the regime where regular statistical mechanics is applicable in the Gibbsian sense. Further, we have traced the explanation of the appearance of localized modes in different regimes in phase space to the need for a modified measure to ensure normalizability, which therefore necessitates a phase transition leading to these localized modes. Our numerical simulations strongly support this theoretical picture, illuminating this novel quantitative connection between nonlinear dynamics and statistical mechanics.

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[5] Negative temperatures appear as an artifact of applying the Hamiltonian, which in this system is not a true measure of the energy, in the statistical formalism.


[7] In the numerical examples we take $\nu = 1$, since a rescaling of $\psi_m$ allows restriction to this case.


