

Collective quantum enhancement in critical quantum sensing

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Critical systems represent a valuable resource in quantum sensing and metrology. Critical quantum sensing (CQS) protocols can be realized using finite-component phase transitions, where criticality is not due to the thermodynamic limit but rather to the rescaling of the system parameters. In particular, the second-order phase transitions of parametric Kerr resonators are of high experimental relevance, as they can be implemented and controlled with various quantum technologies currently available. Here, we show that collective quantum advantage can be achieved with a multipartite critical quantum sensor based on a parametrically coupled Kerr resonators chain in the weak-nonlinearity limit. We derive analytical solutions for the low-energy spectrum of this unconventional quantum many-body system, which is composed of *locally* critical elements. We then assess the performance of an adiabatic CQS protocol, comparing the coupled-resonator chain with an equivalent ensemble of independent critical sensors. We evaluate the scaling of the quantum Fisher information with respect to fundamental resources, and find that the critical chain achieves a quadratic enhancement in the number of resonators. Beyond the advantage found in the case of zero Kerr, we find that there is a collective enhancement even in the scenario of finite Kerr nonlinearity.

I. INTRODUCTION

Critical quantum sensing (CQS) is by now an established approach, based on the exploitation of quantum properties spontaneously developed in proximity of phase transitions. In the context of quantum metrology, the performance of parameter estimation protocols [1] is assessed based on the precision that can be achieved using a limited amount of fundamental resources, such as the size of the probe system and/or the protocol duration time. By using non-classical properties such as entanglement and superposition, it is possible to overcome the performance of any sensing strategy based on purely classical resources [2]. Numerous theoretical studies have shown that a quantum-enhanced sensing precision can be achieved exploiting static [3–13] or dynamical [14–17] properties of many-body systems in proximity of the critical point. Despite critical quantum systems becoming divergently slow as they approach the phase transition, the optimal precision scaling can, in principle, be saturated with respect to both probe size and time [18]. Furthermore, even if prior information on the parameter to be estimated is required to operate the sensor close to the critical point, CQS protocols work well also in a global-sensing scenario when adaptive strategies are applied [19, 20]. First implementations with Rydberg atoms [21], nuclear magnetic resonance [22] and superconducting quantum circuits [23, 24] demonstrate the experimental feasibility of the CQS approach.

Recently, it has been shown [25] that CQS protocols can also be conceived using finite-component phase transitions (FCPTs), where the thermodynamic limit is replaced with a rescaling of the system parameters [26, 27]. This class of phase transitions emerges in quantum resonators with atomic [26, 28–31] or Kerr-like [27, 32–34] nonlinearities, and it is of high relevance for CQS

for two main reasons: (i) It provides a tractable theoretical framework to analyze the fundamental precision bounds and to design optimal parameter estimation protocols. For instance, FCPTs have been used to demonstrate the intrinsic constant-factor advantage of dynamical over statical CQS protocols [35, 36] and the presence of apparent super-Heisenberg scalings when focusing on a single resource [37–39]. Continuous-measurement schemes [40, 41] have been designed to efficiently retrieve information in the dissipative case, where the optimal precision bounds are in general not achievable. In the presence of a thermal bath, CQS outperforms passive strategies when preparation and measurement times are not negligible [42]. (ii) As demonstrated by recent experiments, FCPTs can be *controllably* implemented using atomic [43], polaritonic [44, 45] and circuit-QED [46–50] platforms. Accordingly, CQS protocols can be implemented with small-scale devices, without requiring the implementation and control of many-body quantum systems. CQS protocols based on FCPTs have been designed for quantum resonators [23, 51–54], single trapped-ions [55], optomechanical [56, 57] or magnomechanical [58] devices, spin impurities [59, 60] and Rabi-like systems [61–64]. Finally, physical devices undergoing FCPTs can be interconnected to form one- or two-dimensional arrays of critical systems, thereby extending CQS approaches to complex extended systems.

Here, we theoretically show that collective quantum enhancement in CQS can be achieved using an array of coupled critical quantum sensors. In particular, we consider a model composed of a chain of nonlinear quantum resonators coupled via a parametric nearest-neighbor interaction. We focus on the weak-nonlinearity limit, where each resonator of the array can become locally critical and undergo an FCPT. We assess the metrological performance of a static CQS protocol by analyti-

cally evaluating the quantum Fisher information (QFI) over the ground state manifold, under Gaussian approximation. In proximity to the critical point, which is the optimal working regime, we find closed-form analytical solutions for the asymptotic scaling of the QFI with respect to fundamental resources, i.e., the number of resonators, the number of excitations and the protocol-duration time. This analysis allows us to benchmark the estimation precision achievable with the coupled-resonator chain against that of an equal number of independent critical sensors. We find that the coupled chain yields a quadratic enhancement of the QFI scaling with respect to the number of resonators, compared to the linear scaling observed in an uncoupled array. Beyond the scaling advantage, we develop a perturbation theory to estimate the saturation point due to local Kerr nonlinearities, and show that the resonator chain presents a practically relevant advantage for finite values of physical parameters.

This article is organized as follows. In Sec. II, we introduce the adiabatic CQS protocol, the physical model of a single-mode critical quantum sensor, and we summarize the main results obtained for the coupled-resonator chain. In Sec. III, we discuss the analytical solutions for the local critical models, i.e., single- and two-resonator systems, which are propaedeutic to analyze the collective case. In Sec. IV, we first present the full analytical solution for the critical resonator chain. Then, we provide approximated closed-form solutions for two relevant parameter regimes: the asymptotic scaling near the critical point and the limit of an infinite number of resonators, dubbed continuous limit. In Sec. V, we exploit the analytical solutions to compare the sensing performance of the critical resonator chain with that of an ensemble of uncoupled critical sensors, examining both the asymptotic scaling and the saturation for a finite nonlinearity. Finally, we discuss possible experimental implementations and future perspectives.

II. SUMMARY OF RESULTS

A. Critical quantum sensing: adiabatic protocol

Here, we briefly introduce the CQS protocol and the methods used to assess parameter-estimation performance. Let us consider as a sensor a system whose Hamiltonian depends on an unknown parameter ω to be estimated, while all other parameters are known, and such that one control parameter ϵ can be tuned in real-time. A static CQS protocol can be defined as a three-step algorithm: (i) the system is initialized in its ground state in a trivial region of the parameter space, say $\epsilon = 0$; (ii) the control parameter ϵ is adiabatically tuned close to the critical value ϵ_c , slowly enough so that the system never leaves its instantaneous ground state; (iii) a measurement is performed on the final state, and the outcome is used to infer information about the value of ω . The working

principle behind adiabatic CQS protocols is the exploitation of the high susceptibility of the system ground state with respect to small variations of the parameter to be estimated.

The ultimate precision achievable in parameter estimation is given by the quantum Cramér-Rao bound [2]. The mean squared error is lower-bounded by $\Delta\omega^2 \geq (M\mathcal{I})^{-1}$, where M is the number of independent measurements and \mathcal{I} is the QFI. For our adiabatic protocol, the QFI must be evaluated [as in Eq. (8)] on the manifold defined by the system ground state as the parameter ω is varied. The QFI represents the maximum amount of information that can be extracted on average for a given repetition of the protocol, assuming that the optimal observable is measured after the adiabatic sweep. The performance of a parameter-estimation protocol can be assessed by analyzing the scaling of the QFI with the fundamental resources used during the protocol, e.g., the size of the probe system N . For example, in Ramsey interferometry, the QFI for phase estimation can grow at most linearly as $\mathcal{I} \sim N$ when the spins are in a separable state, while it can achieve the quadratic scaling $\mathcal{I} \sim N^2$ when the spins are entangled [65]. This quadratic enhancement is dubbed Heisenberg scaling and is achievable only using nonclassical resources such as entanglement or squeezing.

The size of the probe system is not the only relevant resource. When dealing with critical systems the protocol time must unavoidably be taken into account as a fundamental resource. When the critical point is approached, the energy gap between the ground and first-excited state tends to zero, and so the time T required to perform an adiabatic sweep unavoidably diverges. Analyzing the scaling of the QFI only with respect to the number of probe systems can then lead to an apparent super-Heisenberg scaling. Once time is taken into account [18, 25], under very general assumptions, the optimal Heisenberg scaling is given by $\mathcal{I} \sim T^2 N^2$. However, this is strictly valid only for closed quantum systems. When decoherence and dissipation are included, stricter asymptotic bounds apply [66]. Here, we limit ourselves to the unitary case, as our focus is on the collective enhancement achievable with quantum-correlated probes, and our results can be generalized to the dissipative case.

B. Critical parametric quantum sensor

Let us now introduce our physical model of a critical quantum sensor, which is a quantum nonlinear resonator with two-photon pumping terms, described by the Hamiltonian

$$H = \omega a_1^\dagger a_1 + \frac{\epsilon}{2} (a_1^2 + a_1^{\dagger 2}) + \chi a_1^\dagger a_1^\dagger a_1 a_1. \quad (1)$$

Here and throughout the manuscript, we set ($\hbar = 1$). This system undergoes an FCPT in the limit $\chi \rightarrow 0$, when ϵ approaches the critical point $\epsilon_c = \omega$. This model is of direct experimental relevance [48], and it effectively describes the critical scaling of fully-connected

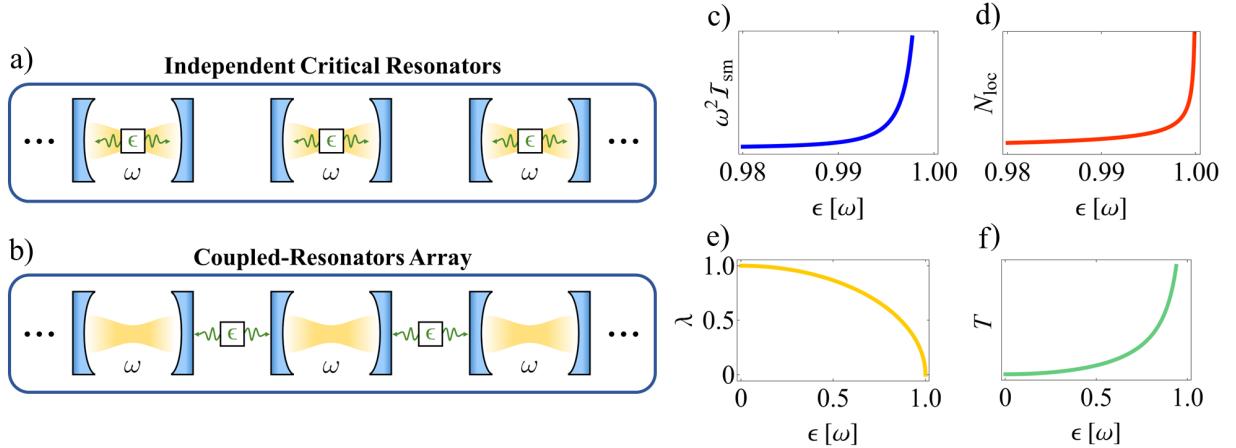


FIG. 1. We illustrate **a)** an ensemble of independent critical resonators and **b)** a chain of parametrically-coupled critical resonators. A single critical sensor is described by the Hamiltonian (1), while the coupled-resonator chain is modeled by the Hamiltonian (2). Focusing on a single independent critical sensor, we plot **c)** the QFI \mathcal{I}_{sm} and **d)** the local number of photons N_{loc} as functions of ϵ . As the squeezing parameter ϵ approaches the critical point $\epsilon_c = \omega$, both \mathcal{I}_{sm} and N_{loc} diverge. Furthermore, we show **e)** the energy gap λ and **f)** the estimate of the protocol duration time T as a function of ϵ . When $\epsilon \rightarrow \epsilon_c$, λ tends to zero, indicating that the energy gap between the ground and first-excited state closes. Consequently, the time required to perform the adiabatic sweep diverges. To obtain the plot for T we set $\eta = 1$.

models [36], a broad class of quantum-optical critical systems. The CQS protocol consists of initializing the system in the vacuum state, and adiabatically tuning ϵ to a value close to ϵ_c . To work in this critical regime, we assume to have good prior information on the parameter to be estimated, namely $\omega = \omega_0 + \delta\omega$, where ω_0 is known and $\delta\omega \ll \omega_0$ is an unknown frequency shift to be estimated. For global quantum sensing, i.e., when prior information is not available, efficient adaptive strategies can be implemented [19, 20].

As shown in Sec. III, under Gaussian approximation ($\chi = 0$), we can obtain analytical results. Indeed, we first compute the ground state manifold, and then evaluate the QFI for the estimation of ω . Consequently, we provide an upper-bound on the estimation precision with respect to the involved resources, i.e., the number of photons $N_{\text{loc}} = \langle a_1^\dagger a_1 \rangle$ and the protocol duration time T . Here, the subscript stands for *local* and will be meaningful in the multipartite case. By applying time-dependent perturbation theory, it can be shown [25] that the time required to perform the adiabatic sweep is of the order of $T \sim (\eta\lambda)^{-1}$, where λ is the energy gap between the ground and first-excited state, and $\eta < 1$ is a parameter which controls the adiabaticity of the protocol. As shown in Fig. 1, when approaching the critical point the QFI and the photon number diverge, while the energy gap between the ground and the first-excited state vanishes, as expected. For a single-mode critical sensor, the QFI scales as $\mathcal{I}_{\text{sm}} \sim 2\eta^2 T^2 N_{\text{loc}}^2$. Here, the subscript “sm” stands for “single-mode”. We stress that in the limit $\chi \rightarrow 0$, this result can be analytically derived. In the case of finite nonlinearity, this scaling is valid until the photon number becomes large enough to break the Gaussian approximation and reach saturation. As shown

in Appendix C, perturbation theory reveals that the saturation point occurs when $N_{\text{loc}} \sim \sqrt[3]{\omega/132\chi}$.

C. Collective quantum advantage in CQS

Let us now summarize the main results of this work. We perform a comparison between an ensemble of M independent critical sensors, with a chain of parametrically-coupled resonators (see Fig. 1). The QFI is additive when using M independent sensors or, equivalently, M uncorrelated repetitions of the parameter-estimation protocol. Accordingly, when M independent critical sensors are used, we straightforwardly obtain $\mathcal{I}_{\text{ind}} \sim 2M\eta^2 T^2 N_{\text{loc}}^2$, with the subscript “ind” standing for “independent”. This result can also be rewritten in terms of the total number of photons $N = MN_{\text{loc}}$, as $\mathcal{I}_{\text{ind}} \sim 2\eta^2 T^2 N^2/M$.

Our collective critical sensor is defined as a coupled-resonator chain with Hamiltonian

$$H = \sum_{j=1}^M \left[\omega a_j^\dagger a_j + \frac{\epsilon}{2} (a_j a_{j+1} + a_j^\dagger a_{j+1}^\dagger) + \chi a_j^\dagger a_j^\dagger a_j a_j \right]. \quad (2)$$

In Sec IV, we derive analytical solutions for the QFI for the estimation of ω , as well as for the photon number and the energy gap, under Gaussian approximation. The result for the QFI is shown in Eqs. (21) and (22) for even and odd M , respectively. We then obtain approximate closed-form analytical solutions in two relevant limits: in proximity of the phase transition and in the continuous limit (large M). The analysis unveils interesting features of this many-body model composed of locally critical con-

stituents. In this summary, we focus on the most effective regime for quantum sensing, which is the critical scaling of a chain with an odd number of resonators. As shown in Sec. IV A, the QFI achievable in this case scales as $I_O = 2\eta^2 M^2 T^2 N_{\text{loc}}^2 = 2\eta^2 T^2 N^2$. Thus, the critical chain presents a quadratic enhancement in the number of resonators M with respect to the non-interacting case. The critical chain is advantageous even when considering a finite value of the Kerr nonlinearity. Indeed, since it can be effectively described by a single-mode resonator with a diluted nonlinearity χ/M (see Appendix D), the saturation point is reached when $N \sim \sqrt[3]{M\omega/132\chi}$.

III. LOCAL MODELS

In this section, we analyze the performance of single- (Sec. IIIA) and two-mode (Sec. IIIB) critical quantum sensors. In both cases, we compute the QFI and express it in terms of the resources involved, i.e., the average number of photons in each mode and the duration time of the sensing protocol. The results will be used as a benchmark, as well as an intermediate step to solve the interacting chain.

A. Single-mode critical quantum system

A single-mode critical quantum sensor can be modeled, under Gaussian approximation, by the Hamiltonian

$$H_{\text{sm}} = \omega a_1^\dagger a_1 + \frac{\epsilon}{2} \left(a_1^2 + a_1^{\dagger 2} \right), \quad (3)$$

where ϵ is the squeezing parameter, ω is the mode frequency and a_1 is the bosonic mode. The full model presented in Eq. (1) undergoes a second-order phase transition in proximity of the critical point $\epsilon_c = \omega$ in the limit $\chi \rightarrow 0$. For $\epsilon < \epsilon_c$ the model is well approximated by the Hamiltonian (3), which is diagonalizable via a Bogoliubov Transformation (BT), obtaining

$$H_{\text{sm}} = \lambda d_1^\dagger d_1 + \frac{1}{2} (\lambda - \omega). \quad (4)$$

We defined $\lambda = \sqrt{\omega^2 - \epsilon^2}$, and d_1 is the normal mode related to a_1 via the transformation

$$a_1 = t d_1 - s d_1^\dagger, \quad (5)$$

where the parameters t and s are

$$t = \frac{(\lambda + \omega)}{\sqrt{2(\omega^2 - \epsilon^2) + 2\omega\lambda}}, \quad (6)$$

$$s = \frac{\epsilon}{\sqrt{2(\omega^2 - \epsilon^2) + 2\omega\lambda}}. \quad (7)$$

Let us focus now on the estimation of the parameter ω . Given a manifold of pure states $|\psi\rangle$, the QFI can be evaluated as [2]

$$\mathcal{I} = 4 \left(\langle \partial_\omega \psi | \partial_\omega \psi \rangle - |\langle \partial_\omega \psi | \psi \rangle|^2 \right), \quad (8)$$

where $|\partial_\omega \psi\rangle$ is the partial derivative with respect to the parameter to be estimated. To assess the performance of the adiabatic CQS protocol, we will compute the QFI associated with the system ground state $|g\rangle$. The ground state of the Hamiltonian (3) is the single-mode squeezed vacuum $|g\rangle = S_1(\xi)|0\rangle$. We define the single-mode squeezing operator $S_i(\xi) = e^{\frac{1}{2}(\xi^* a_i^2 - \xi a_i^{\dagger 2})}$ and the squeezing parameter $\xi = |\xi| = \ln(s + \sqrt{s^2 + 1})$. This state has an average number of photons $N_{\text{loc}} = \sinh^2(|\xi|) = s^2$, where “loc” stands for “local”. By inserting in (8) the expression of the system ground state, we obtain the single-mode QFI for the estimation of ω :

$$\mathcal{I}_{\text{sm}} = \frac{\epsilon^2}{2(\omega^2 - \epsilon^2)^2}. \quad (9)$$

The dominant contribution to the duration of the sensing protocol is given by the time T required to perform the adiabatic sweep, which must be larger [25] than the inverse of the gap between the ground and first-excited state. From (4), it immediately follows that the energy gap is $\lambda = \sqrt{\omega^2 - \epsilon^2}$. Consequently, the protocol duration time can be expressed as $T \approx (\eta\lambda)^{-1}$, with $\eta < 1$. To analyze the critical scaling approached when $\epsilon \rightarrow \epsilon_c$, we define $\epsilon = (1-x)\omega$ with $0 < x < 1$, and compute the QFI in the asymptotic regime $x \rightarrow 0$. Close to the critical point, we have $\mathcal{I}_{\text{sm}} \approx (2\sqrt{2}\omega x)^{-2}$, $T \approx (\omega\sqrt{2x})^{-1}$ and $N_{\text{loc}} \approx (8x)^{-\frac{1}{2}}$. Using the latter, we express x in terms of the number of photons as $x \approx (2\sqrt{2}N_{\text{loc}})^{-2}$, which implies $T \approx 2N_{\text{loc}}/\eta\omega$. Finally, in terms of the resources involved the single-mode QFI scales as

$$\mathcal{I}_{\text{sm}} \sim 2\eta^2 T^2 N_{\text{loc}}^2. \quad (10)$$

B. Two-mode critical quantum system

Similarly to the single-mode case, a two-mode critical quantum system can be modeled with a two-mode squeezing Hamiltonian. The single-resonator critical sensor can be generalized to a two-mode critical system as follows:

$$H_{\text{tm}} = \omega \left(a_1^\dagger a_1 + a_2^\dagger a_2 \right) + \epsilon \left(a_1 a_2 + a_1^\dagger a_2^\dagger \right), \quad (11)$$

where ϵ is the parametric-coupling strength, ω is the frequency of each mode and a_1 and a_2 are two independent bosonic modes. Here, the subscript “tm” means “two-mode”. Also in this case, the model undergoes a second-order phase transition, when a vanishingly small Kerr nonlinearity is included [27]. The normalization of the coupling parameter is chosen to have the same critical point $\epsilon_c = \omega$ of the Hamiltonian (1). When $\epsilon < \epsilon_c$, the Hamiltonian (11) can be diagonalized with the following BT:

$$\begin{aligned} a_1 &= t d_1 - s d_2^\dagger, \\ a_2 &= t d_2 - s d_1^\dagger, \end{aligned} \quad (12)$$

where d_1, d_2 are the normal modes. We can then write

$$H_{\text{tm}} = \lambda(d_1^\dagger d_1 + d_2^\dagger d_2) + \lambda - \omega. \quad (13)$$

The ground state of the system is a two-mode squeezed vacuum state $|g\rangle = S_{1,2}(\xi)|0\rangle$, with $S_{i,j}(\xi) = e^{(\xi^* a_i a_j - \xi a_i^\dagger a_j^\dagger)}$ a two-mode squeezing operator. Notice that the squeezing parameter $\xi = |\xi|$ is the same as in the single-mode case. The average number of photons in the ground state is $N = 2 \sinh^2(|\xi|) = 2s^2$ since now there are two normal modes with the same eigenvalue λ and the same number of photons $N_{\text{loc}} = \langle a_1^\dagger a_1 \rangle = \langle a_2^\dagger a_2 \rangle = s^2$. Using (8), we can compute the two-mode QFI for the estimation of ω :

$$\mathcal{I}_{\text{tm}} = \frac{\epsilon^2}{(\omega^2 - \epsilon^2)^2}. \quad (14)$$

In the asymptotic limit $x \rightarrow 0$, we have $\mathcal{I}_{\text{tm}} \approx (2\omega x)^{-2}$ and $N_{\text{loc}} \approx (8x)^{-\frac{1}{2}}$. Since the energy gap is still λ , the protocol duration time remains unchanged with respect to the single mode case. Indeed, expressing x as $x \approx (2\sqrt{2}N_{\text{loc}})^{-2}$ leads to $T \approx 2N_{\text{loc}}/\eta\omega$. Finally, in terms of the involved resources the two-mode QFI scales as

$$\mathcal{I}_{\text{tm}} \sim 4\eta^2 T^2 N_{\text{loc}}^2. \quad (15)$$

Comparing this result with Eq. (10), we see that \mathcal{I}_{tm} is simply twice what is obtained with a single critical sensor. A two-resonator chain is then equivalent to two independent resonators, suggesting that there is no advantage in considering coupled critical sensors. However, in the following, we demonstrate that this conclusion does not hold for larger chains.

IV. CRITICAL RESONATOR CHAIN

Let us finally consider the critical resonator chain. Under Gaussian approximation, it is described by the Hamiltonian

$$H = \sum_{j=1}^M \left[\omega a_j^\dagger a_j + \frac{\epsilon}{2} (a_j a_{j+1} + a_j^\dagger a_{j+1}^\dagger) \right]. \quad (16)$$

Assuming periodic boundary conditions $a_{M+j} = a_j$, the Hamiltonian (16) is diagonalizable in the reciprocal space. By applying the discrete Fourier transform $a_j = (M)^{-\frac{1}{2}} \sum_k a_k e^{-ikj}$, we obtain

$$H = H_0 + \delta_{M,2m} H_\pi + \sum_{k>0} \omega (a_k^\dagger a_k + a_{-k}^\dagger a_{-k}) + \sum_{k>0} \epsilon \cos(k) (a_k a_{-k} + a_k^\dagger a_{-k}^\dagger). \quad (17)$$

Here, $\delta_{M,2m}$ is a Kronecker delta which is nonzero only for even M . The terms H_0, H_π are single-resonator critical models defined as in Eq. (3), but replacing a_1

with a_0 and a_π , respectively. The sign of the coupling for the a_π mode must be inverted, a phase difference that has no consequence for the current analysis. The periodic boundary conditions imply that we can restrict ourselves to the first Brillouin zone (FBZ), and $k = 2n\pi/M$, with $n \in \mathbb{Z}$. The FBZ is defined as follows: if M is even, $n \in [-M/2 + 1, M/2]$; if M is odd, $n \in [-(M-1)/2, (M-1)/2]$. We point out that the mode a_π is present only in the case in which the number of modes M is even.

In reciprocal space, the multi-mode Hamiltonian (17) consists of M independent critical systems. Specifically, two single-mode systems defined by a_0 and a_π , along with a series of two-mode systems defined by pairs $\{a_k, a_{-k}\}$. Each element can be independently diagonalized using the single- and two-mode BTs derived in Sec. III, by replacing $\epsilon \rightarrow \epsilon \cos(k)$. Notice that the BT can be performed only if $\epsilon < \epsilon_c$. In the multi-mode case, we have many critical points defined by the dispersion relation $\epsilon_c = \omega / \cos(k)$. This means that the multi-mode Hamiltonian is diagonalizable via a BT only if the parameter ϵ is smaller than the lowest among the critical points. Therefore, we restrict ourselves to the case $\epsilon < \omega$. If this condition is satisfied, applying the BT to (17) results in

$$H = \sum_{k \in FBZ} \lambda_k \left(d_k^\dagger d_k + \frac{1}{2} \right) - \frac{M}{2} \omega, \quad (18)$$

where $\lambda_k = \sqrt{\omega^2 - \epsilon^2 \cos^2(k)}$, and d_k are independent normal modes. The modes d_0, d_π are defined as in (5), while the modes $d_{\pm k}$ as in (12), with the substitution $\epsilon \rightarrow \epsilon \cos(k)$. The ground state of the system is then a multi-mode squeezed vacuum state

$$|g\rangle = S_0(\xi_0) S_\pi(\xi_\pi) \bigotimes_{k>0} S_{k,-k}(\xi_k) |0\rangle. \quad (19)$$

We recall that $S_i(\xi_i) = e^{\frac{1}{2}(\xi_i^* a_i^2 - \xi_i a_i^{\dagger 2})}$ with $i = \{0, \pi\}$ is a single-mode squeezing operator associated with the modes a_0, a_π . Conversely, $S_{k,-k}(\xi_k) = e^{(\xi_k^* a_k a_{-k} - \xi_k a_k^\dagger a_{-k}^\dagger)}$ is a two-mode squeezing operator related to pairs of coupled modes $\{a_k, a_{-k}\}$. The squeezing parameters are $\xi_k = |\xi_k| = \ln(s_k + \sqrt{s_k^2 + 1})$ for $k \in [0, \pi/2]$ and $\xi_k = |\xi_k| e^{i\pi}$ for $k \in [\pi/2, \pi]$. The parameter s_k is obtained from (7) with the substitution $\epsilon \rightarrow \epsilon \cos(k)$. The total average number of photons in the multi-mode ground state is given simply by the sum of the photons in each mode a_k . Given that the ground state is a multi-mode squeezed vacuum state, each mode contains $N_k = \sinh^2(|\xi_k|) = s_k^2$ photons. Since $s_k^2 = s_{-k}^2$, the total number of photons in the ground state can be expressed as

$$N = \sum_k N_k = s_0^2 + \delta_{M,2m} s_\pi^2 + \sum_{k>0} 2s_k^2. \quad (20)$$

Moreover, the local number of photons at each site j is determined by $N_{\text{loc}} = \langle a_j^\dagger a_j \rangle = \sum_k \langle a_k^\dagger a_k \rangle / M = N/M$.

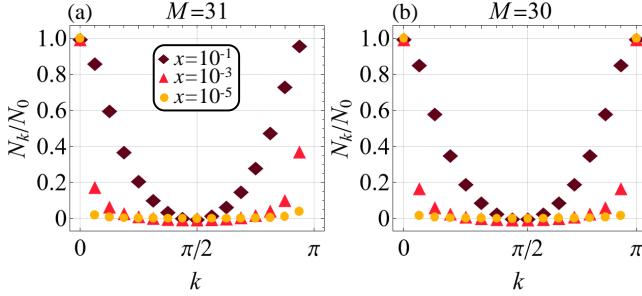


FIG. 2. Local number of photons. We plot the ratio N_k/N_0 for a) $M = 31$ and b) $M = 30$ for different values of $x = 1 - \epsilon/\omega$. As the critical point is approached ($x \rightarrow 0$), the majority of photons is strongly confined in the mode a_0 when M is odd, and in the modes a_0, a_π when M is even (see yellow circles). The fraction of the photons populating the remaining modes is vanishingly small.

A detailed derivation is provided in Appendix A, where we show that N_{loc} is independent of the specific site, i.e., each resonator has the same number of photons. It is important to stress that in the reciprocal space, when the critical point is approached, the photons are strongly confined only in the mode a_0 when M is odd and in the modes a_0, a_π when M is even. The fraction of photons populating all the other modes vanishes as $\epsilon \rightarrow \epsilon_c$ (see Fig. 2). In this scenario, we can effectively approximate the critical chain as a single-mode system with $N = MN_{\text{loc}}$ photons if M is odd, or as a two-mode system with $N = MN_{\text{loc}}/2$ photons per mode when M is even.

Having obtained its ground state, we can now use (8) to calculate the QFI of the critical chain. From Eq. (19) we see that the ground state is separable in reciprocal space, hence the QFI of the critical chain is given by the sum of the QFIs of each individual mode a_0 and a_π , and of each pair $\{a_k, a_{-k}\}$. We obtain

$$\mathcal{I}_E = \sum_{n=0}^{\frac{M}{2}-1} \frac{\epsilon^2 \cos^2(\frac{2\pi n}{M})}{[\omega^2 - \epsilon^2 \cos^2(\frac{2\pi n}{M})]^2}, \quad (21)$$

$$\mathcal{I}_O = \frac{\epsilon^2}{2(\omega^2 - \epsilon^2)^2} + \sum_{n=1}^{\frac{M}{2}-1} \frac{\epsilon^2 \cos^2(\frac{2\pi n}{M})}{[\omega^2 - \epsilon^2 \cos^2(\frac{2\pi n}{M})]^2}, \quad (22)$$

where \mathcal{I}_E is the QFI for an even number of modes, while \mathcal{I}_O is the QFI for an odd number of modes. Let us now provide closed-form analytical expressions of the multi-mode QFI in relevant asymptotic regimes. First, in Sec. IV A we consider a finite number of modes and compute explicitly the QFI close to the critical point. Then, in Sec. IV B we compute the QFI in the continuous limit $M \rightarrow \infty$. In both cases, we express the QFI in terms of the involved resources, i.e., the local number of photons N_{loc} and the total protocol duration time T .

A. Asymptotic regimes: Critical scaling

In this section, we provide the scaling behavior of the critical chain QFI in the asymptotic regime where $\epsilon \rightarrow \epsilon_c$, assuming a finite M . To analyze the critical scaling we define $\epsilon = (1 - x)\omega$ with $0 < x < 1$, where the critical point is approached for $x \rightarrow 0$. We then expand Eqs. (21) and (22) in power series, up to the first order in x . Following this expansion, the QFI separates into two terms: the first depends solely on x and diverges as the critical point is approached, while the second term is a discrete sum dependent only on the number of modes M . This sum can be analytically computed using established results from trigonometric series, allowing us to derive a closed expression for the asymptotic QFI. See Appendix B for further mathematical details. Thus, in the limit $x \rightarrow 0$, up to the terms of order x , the QFI for even and odd critical chains can be expressed as

$$\mathcal{I}_E \approx \frac{1}{4\omega^2} \left(\frac{1}{x^2} - \frac{1}{x} - \frac{1}{4} + \frac{M^4 - 20M^2 + 64}{180} \right), \quad (23)$$

$$\mathcal{I}_O \approx \frac{1}{8\omega^2} \left(\frac{1}{x^2} - \frac{1}{x} - \frac{1}{4} + \frac{4M^4 - 20M^2 + 16}{45} \right). \quad (24)$$

A similar analysis applies to the average number of photons in the multi-mode ground state. Notice that the expression for the number of photons is given by (20), and the only distinction between odd and even M lies in the range of k values in the discrete sum, as specified in Sec. III A. Then, expanding (20) up to terms of order \sqrt{x} yields

$$N \approx \left\{ \begin{array}{l} \frac{1}{\sqrt{2x}} + \frac{M}{\pi} \left[\log\left(\frac{M}{\pi}\right) + \gamma - \frac{\pi}{2} \right], \quad M \text{ even} \\ \frac{1}{\sqrt{8x}} + \frac{M}{\pi} \left[\log\left(\frac{2M}{\pi}\right) + \gamma - \frac{\pi}{2} \right], \quad M \text{ odd} \end{array} \right. \quad (25)$$

$$N \approx \left\{ \begin{array}{l} \frac{1}{\sqrt{2x}} + \frac{M}{\pi} \left[\log\left(\frac{M}{\pi}\right) + \gamma - \frac{\pi}{2} \right], \quad M \text{ even} \\ \frac{1}{\sqrt{8x}} + \frac{M}{\pi} \left[\log\left(\frac{2M}{\pi}\right) + \gamma - \frac{\pi}{2} \right], \quad M \text{ odd} \end{array} \right. \quad (26)$$

where γ is the Euler-Mascheroni constant. See Appendix B for further details on the derivation of Eqs. (25)-(26). Our aim now is to express the multi-mode QFI in terms of the resources involved, namely the local number of photons N_{loc} and the protocol duration time T . To achieve this, we first express the parameter x as a function of the local number of photons using (25)-(26). Specifically, we find $x \approx (\sqrt{2}MN_{\text{loc}})^{-2}$ when M is even and $x \approx (2\sqrt{2}MN_{\text{loc}})^{-2}$ when M is odd. It is important to note that these relations hold true under the condition that the local number of photons is sufficiently large, satisfying $\pi N_{\text{loc}} \gg |\ln(M/\pi) + \gamma - \pi/2|$ for M even, and $\pi N_{\text{loc}} \gg |\ln(2M/\pi) + \gamma - \pi/2|$ for M odd. This assumption is reasonable in the considered asymptotic regime, where M is finite and near the criticality. On the other hand, from (18) we have that the lowest energy gap is λ_0 , which means that the protocol duration time is estimated as $T \approx (\eta\lambda_0)^{-1} \approx (\eta\omega\sqrt{2x})^{-1}$. In terms of the local number of photons we write $T \approx MN_{\text{loc}}/\eta\omega$ for M even and $T \approx 2MN_{\text{loc}}/\eta\omega$ for M odd. Finally, inserting the expression of x in (23)-(24) and taking into account

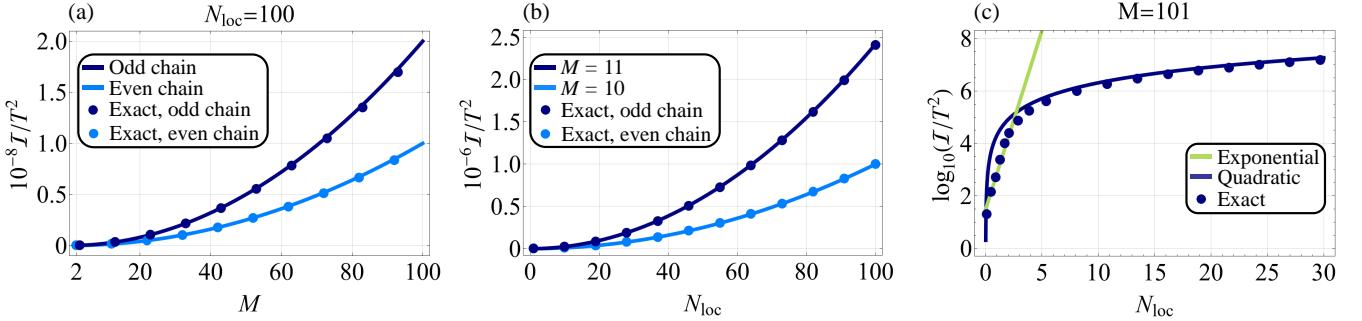


FIG. 3. Quantum Fisher information. We show the ratio \mathcal{I}/T^2 as a function of **a)** the number of modes M , for M odd (dark blue) and M even (light blue); and as a function of **b)** the local number of photons N_{loc} , for $M = 11$ (dark blue) and $M = 10$ (light blue). The quadratic scaling (27)-(28) (solid lines) fits well with the exact ratio (small circles) computed numerically from (21)-(22). **c)** Plot of $\log_{10}(\mathcal{I}/T^2)$ as a function of N_{loc} , for $M = 101$. The exact ratio computed numerically (small circles) is well approximated by the exponential scaling (32) (solid green line) only when $N_{\text{loc}} = \mathcal{O}(1)$. However, as soon as N_{loc} becomes comparable to M , the quadratic scaling (28) fits well the exact ratio (solid blue line), while the exponential scaling is no longer valid. The same behavior holds also in the case of an even M . To obtain the plots, for each value of M we have optimized the QFI under the constraint $N_{\text{loc}} = N_{\text{max}}$, where in **a)** $N_{\text{max}} = 100$, while in **b)** $N_{\text{max}} \in (0, 100]$ and in **c)** $N_{\text{max}} \in (0, 30]$. We have set $\omega = 1$.

the protocol duration time, the QFI scales as

$$\mathcal{I}_E \sim \eta^2 M^2 T^2 N_{\text{loc}}^2, \quad (27)$$

$$\mathcal{I}_O \sim 2\eta^2 M^2 T^2 N_{\text{loc}}^2. \quad (28)$$

These solutions then predict a quadratic scaling of the QFI with the number of resonators M , which implies a collective quantum enhancement of the estimation precision. The difference between the even and odd cases arises from the fact that, for an odd M , only the mode with $k = 0$ becomes critical whereas, for an even M , both the modes with $k = 0$ and $k = \pi$ become critical (see Fig. 2). In Fig. 3(a) and (b), we plot the optimized ratio \mathcal{I}/T^2 as a function of M and as a function of N_{loc} , respectively. We show that the asymptotic expressions (27)-(28) align perfectly with the full analytical solutions of Eqs. (21)-(22), in the asymptotic regime where $x \rightarrow 0$ and M is finite.

B. Asymptotic regimes: Continuous limit

As M increases, the separation between adjacent modes decreases. This is due to the periodic boundary conditions, which impose $\Delta k = 2\pi/M$, so that $\Delta k \rightarrow 0$ when $M \rightarrow \infty$. Consequently, in this regime we have a continuum of modes a_k . Therefore, we can compute the QFI in (21)-(22) by approximating the discrete sums with definite integrals over this continuum, obtaining

$$\mathcal{I} = \frac{M}{4\omega^2} \frac{(1-x)^2}{(x(2-x))^{\frac{3}{2}}}. \quad (29)$$

Notice that, of course, by making this approximation, a numerical error is introduced. This error becomes negligible when the number of modes is sufficiently large, such

that $\Delta k \rightarrow dk$. We estimate the absolute error incurred when approximating a discrete sum with an integral via

$$\left| \sum_{k>0} f(k)\Delta k - \int_0^\pi f(k)dk \right| < \frac{M}{2} \Delta k^2 \max\{\partial_k f(k)\}. \quad (30)$$

By imposing $M\Delta k^2 \max\{\partial_k f(k)\}/2 \ll 1$, we can evaluate how large M should be so that (29) is a good approximation of the QFI. In the limit of $x \rightarrow 0$, we find that $\max\{\partial_k f(k)\} \approx x^{-\frac{5}{2}}/5$, which implies that, to have a negligible absolute error, we require $M \gg 2\pi^2 x^{-\frac{5}{2}}/5$. This indicates that the closer we are to the critical point, the larger the M needed to ensure the validity of (29).

The same reasoning can be applied to the average number of photons in the ground state (20). However, the resulting integral cannot be computed analytically as we did with the QFI. An approximated analytical solution in the limit $x \rightarrow 0$ is

$$N \approx M \left[\frac{1}{2\pi} \log \left(\frac{8}{x} \right) - \frac{1}{2} \right]. \quad (31)$$

Notice that in this regime, the local number of photons $N_{\text{loc}} = N/M$ is independent of the number of modes but depends only on the proximity to the critical point. In this case, the parameter x can be expressed in terms of N_{loc} as $x \approx 8e^{-(1+2N_{\text{loc}})\pi}$. Furthermore, the protocol duration time is approximately $T \approx (\eta\lambda_0)^{-1} \approx (\eta\omega\sqrt{2x})^{-1}$, which in terms of N_{loc} becomes $T \approx e^{(\frac{1}{2}+N_{\text{loc}})\pi}/4\eta\omega$. Using these results, in terms of the involved resources we express the QFI as

$$\mathcal{I} \sim \frac{e^{\frac{\pi}{2}}}{16} \eta^2 M T^2 e^{\pi N_{\text{loc}}}. \quad (32)$$

By looking at (32), one might conclude that the optimal operating regime of the critical chain is the continuous limit, given that the QFI scales exponentially

with the photon number resource. However, the validity of this expression is related to (30). As mentioned above, approximating the discrete sums defining the multi-mode QFI with definite integrals is a good approximation only if $M \gg 2\pi^2 x^{-\frac{5}{2}}/5$. Considering the expression of x in terms of N_{loc} , we find that M must satisfy $M \gg C e^{5\pi N_{\text{loc}}}$, with $C = \pi^2 e^{\frac{5\pi}{2}}/320\sqrt{2}$. This reveals that even for a low number of photons, the required M to ensure a negligible error is extremely large. In Fig. 3(c), we plot the ratio $\log_{10}(\mathcal{I}/T^2)$ as a function of N_{loc} , and compare the exact optimized ratio obtained numerically from (22) with the quadratic scaling of (28) and the exponential scaling of (32). As expected, the QFI goes exponentially with N_{loc} only when $N_{\text{loc}} = \mathcal{O}(1)$. Indeed, as soon as N_{loc} becomes comparable with M , the QFI is well approximated by the quadratic regime of (28). The exponential in Eq. (32) is then only an apparent super-Heisenberg scaling.

V. PERFORMANCE ANALYSES

We can now compare the performance of the critical resonator chain with that of an ensemble of M independent single-mode critical resonators. In Sec. V A, we discuss the collective precision enhancement of the array of critical resonators in the relevant asymptotic regime of Sec. IV A. Additionally, in Sec. V B, we assess the validity of the Gaussian approximation of the critical models and provide an estimate of the saturation of the QFI due to finite Kerr nonlinearity. Finally, in Sec. V C we discuss the experimental relevance and future perspectives.

A. Asymptotic scaling

Let us first examine the case of an ensemble of M uncoupled critical resonators. This scenario corresponds to applying the CQS protocol with a single-mode system M times, independently. For independent measurements, the QFI is additive, meaning that overall $\mathcal{I}_{\text{ind}} = M\mathcal{I}_{\text{sm}}$. For an array of independent critical sensors, from (10) we obtain

$$\mathcal{I}_{\text{ind}} \sim 2\eta^2 T^2 M N_{\text{loc}}^2 = 2\eta^2 T^2 \frac{N^2}{M}. \quad (33)$$

Conversely, for a critical chain of M interacting resonators, the optimal QFI scales as

$$\mathcal{I}_E \sim \eta^2 T^2 M^2 N_{\text{loc}}^2 = \eta^2 T^2 N^2, \quad (34)$$

$$\mathcal{I}_O \sim 2\eta^2 T^2 M^2 N_{\text{loc}}^2 = 2\eta^2 T^2 N^2, \quad (35)$$

for an even and odd M , respectively. The critical resonator chain presents a collective enhancement in sensing precision, resulting in a quadratic scaling of the QFI with respect to M . In Fig. 4, we show the growth of the QFI with respect to M for fixed values of N_{loc} , comparing even and odd critical chains with an equivalent ensemble

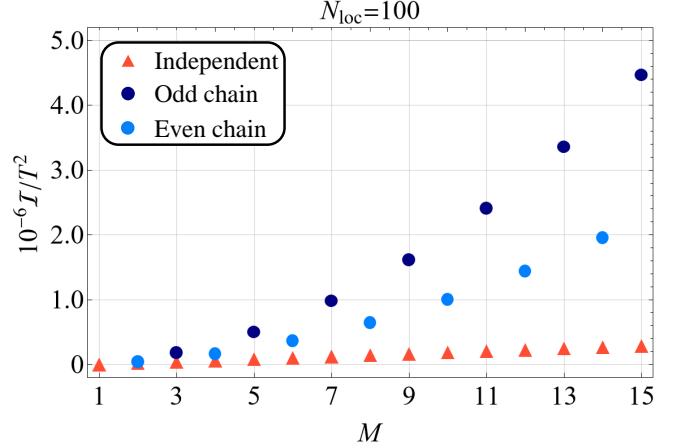


FIG. 4. **QFI: comparison between independent ensemble and critical chain.** We plot the ratio \mathcal{I}/T^2 as a function of M in the case of independent modes (9) (triangles) and coupled modes (21)-(22) (circles). While the QFI scales linearly with M for decoupled modes, interactions between the bosonic modes enhance the QFI, leading to a quadratic scaling with respect to M . To obtain the plot, for each value of M we have optimized the QFI under the constraint $N_{\text{loc}} = 100$, and we have set $\omega = 1$.

of uncoupled resonators. The plot is numerically computed from Eq. (9) and Eqs. (21)-(22).

This collective enhancement can be understood by analyzing the photon distribution in the modes k in the reciprocal space. Let us focus first on a chain with an odd M . As discussed in Sec. IV and shown in Fig. 2, in proximity of the critical point the photons are strongly confined in the single mode a_0 . As can be seen in Eq. (18), this mode is effectively the closest to its critical point. Given that the QFI is nonlinear with respect to the photon number, it is more favorable to concentrate all the photons into the most critical mode, rather than distributing them uniformly. This also explains the origin of the $1/2$ prefactor in the scaling of the QFI for an even chain: when M is even the photons are confined into the two modes $\{a_0, a_\pi\}$, so the collective enhancement is achieved for two uniformly populated modes. Let us now show that confining all photons into a single critical mode is advantageous also when considering saturation.

B. Saturation for finite nonlinearity

Let us now assess the validity of the Gaussian approximation, which is strictly valid for a vanishing nonlinearity ($\chi = 0$) [see Eqs. (1) and (2)]. In any physical implementation, nonlinear effects will inevitably play a role at high energies. For finite values of χ , as the critical point is approached, the Kerr term induces saturation in the photon number and, consequently, in the QFI. For a single critical sensor, the maximum achievable QFI for finite χ can be evaluated numerically [52]. However, nu-

merical simulations of the full quantum model for large chains are beyond the capabilities of classical computers. We then take a perturbative approach to assess the saturation point of the QFI as a function of the ratio between the frequency ω and the nonlinearity χ . For the single-mode system, we applied perturbation theory directly on the QFI, the details can be found in Appendix C. In this case, the validity region of the Gaussian approximation is given by

$$N_{\text{loc}} \lesssim \sqrt[3]{\frac{\omega}{132\chi}}, \quad (36)$$

where N_{loc} is the number of photons in the unperturbed single-mode ground state. Using the bound in (36), we can establish an upper limit also for the QFI in relation to the physical parameters ω and χ . Indeed, recalling that, for a single-mode critical resonator $T \approx 2N_{\text{loc}}/\eta\omega$, the scaling of the QFI can be written solely in terms of the number of photons as $\mathcal{I}_{\text{sm}} \sim 8N_{\text{loc}}^4/\omega^2$. Thus, for an ensemble of M independent critical resonators, the saturation of the QFI is determined by the bound

$$\mathcal{I}_{\text{ind}} \sim \frac{8MN_{\text{loc}}^4}{\omega^2} \lesssim \frac{M}{100} \sqrt[3]{\frac{1}{\omega^2\chi^4}} \lesssim \frac{M}{100\omega^2} \left(\frac{\omega}{\chi}\right)^{4/3}. \quad (37)$$

For the critical chain, we focus exclusively on the case where M is odd, as it yields the most favorable QFI scaling. Applying perturbation theory in this case is much more complex, as the Kerr terms in (2) are spatially local, but highly non-local in the reciprocal space. However, close to the criticality (when saturation effects are relevant), only a single critical mode is effectively populated. In Appendix D, we show that when $x \rightarrow 0$ the Hamiltonian of the critical chain can be mapped onto a single critical mode, with a rescaled nonlinearity χ/M . In other words, for what concerns the critical scaling, the nonlinearity of the interacting chain is effectively diluted. Under single-mode approximation, we can calculate perturbations to the QFI in spite of the nonlocality of the Kerr terms in the reciprocal space. In addition, in Appendix D we confirm also the validity of this single-mode approximation by evaluating perturbative corrections to the energy gap between the ground and first-excited states in the multi-mode case. This perturbative analysis reveals that, for the critical chain, the Gaussian approximation holds true if

$$N \lesssim \sqrt[3]{\frac{\omega M}{132\chi}}, \quad (38)$$

where now $N = MN_{\text{loc}}$ is the number of photons in the multi-mode ground state. It follows that, given $T \approx 2MN_{\text{loc}}/\eta\omega$, the multi-mode QFI (28) is asymptotically upper-bounded by

$$\mathcal{I} \sim \frac{8M^4N_{\text{loc}}^4}{\omega^2} \lesssim \frac{1}{100} \sqrt[3]{\frac{M^4}{\omega^2\chi^4}} \lesssim \frac{1}{100\omega^2} \left(\frac{M\omega}{\chi}\right)^{4/3}. \quad (39)$$

Note that the asymptotic bound (39) is $\sqrt[3]{M}$ times larger than the single-mode bound (37). This shows that the collective enhancement of the critical chain is not canceled by saturation effects. On the contrary, the critical chain presents a scaling advantage also with respect to the ratio ω/χ .

C. Discussion and perspectives

We have demonstrated that a critical quantum sensor based on a chain of M parametrically coupled Kerr resonators exhibits a quadratic enhancement of the QFI with respect to both photon and resonator numbers. Beyond asymptotic scalings, the coupled chain offers practical advantages in terms of physical parameters. This collective quantum advantage in critical quantum sensing is significant even for a small number of resonators. We have analyzed a variety of parameter regimes, and the following table summarizes the QFI scaling for the most interesting cases relevant to quantum sensing applications:

Scaling of the QFI		
	Independent Sensors	Coupled chain
N_{loc} scaling	$2\eta^2 T^2 M N_{\text{loc}}^2$	$2\eta^2 T^2 M^2 N_{\text{loc}}^2$
N scaling	$2\eta^2 T^2 \frac{N^2}{M}$	$2\eta^2 T^2 N^2$
Saturation	$\frac{M}{100\omega^2} \left(\frac{\omega}{\chi}\right)^{4/3}$	$\frac{1}{100\omega^2} \left(\frac{M\omega}{\chi}\right)^{4/3}$

This work paves the way to the exploration of collective quantum advantage in critical quantum sensing, both at experimental and theoretical levels. The considered class of finite-component phase transitions has already been experimentally observed using atomic [43], polaritonic [44, 45] and circuit-QED [46–50] devices. More generally, chains of parametrically-coupled Kerr resonators can be implemented using a broad variety of solid-state platforms, including polaritonics, magnonics, opto- and electro-mechanical systems and superconducting quantum circuits. Further theoretical work is needed to properly support the development of collective CQS protocols. The driven-dissipative case should be addressed considering platform-specific couplings with the environment, which will also depend on the chosen readout mechanism. The need for adiabatic processes, responsible for the $\eta < 1$ prefactor in the scaling of the QFI, can be lifted by designing dynamical protocols [35, 36]. Continuous-measurement schemes [40, 41] can be designed to optimize information retrieval from the output signals. Finally, topological properties have been shown [67] to enhance the sensing precision in sensors based on non-Hermitian lattice dynamics. The inclusion of topological properties in critical chains could potentially enhance the precision and resilience to parameter fluctuations of collective CQS protocols. Be-

sides quantum sensing and metrology, this work presents a phenomenological study of the ground-state properties of an unconventional quantum many-body model, which is critical both at the local and global levels. We analyzed the regime of weak nonlinearity before symmetry breaking occurs. Numerical simulations of the full quantum model are challenging due to large photon numbers and long-range correlations. Phase-space methods appear to be the most suitable for analyzing lattices of critical resonators with weak but finite nonlinearity, where the ground state of the system is approximately Gaussian.

ACKNOWLEDGMENTS

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Appendix A: Local number of photons

In this section, we provide details on the calculation of the local number of photons in the critical resonator chain of Sec. III. The local number of photons in the ground state is given by

$$N_{\text{loc}}^j = \langle g | a_j^\dagger a_j | g \rangle, \quad (\text{A1})$$

where $|g\rangle$ is defined in (19). We express the operators a_j in the reciprocal space as

$$a_j = \frac{1}{\sqrt{M}} \sum_k a_k e^{-ikj}, \quad (\text{A2})$$

$$a_j^\dagger = \frac{1}{\sqrt{M}} \sum_k a_k^\dagger e^{ikj}, \quad (\text{A3})$$

and rewrite the expectation value in (A1):

$$N_{\text{loc}}^j = \langle g | \frac{1}{M} \sum_{k,k'} a_k^\dagger a_{k'} e^{-i(k'-k)j} | g \rangle. \quad (\text{A4})$$

Thus, to obtain N_{loc} , we have now to compute the expectation values $\langle g | a_k^\dagger a_{k'} | g \rangle$, knowing that the ground state is the multi-mode squeezed vacuum state of Eq. (19). Let us consider first the case $(k, k') \neq 0$ and $(k, k') \neq \pi$. Then, the expectation value becomes

$$\begin{aligned} \langle g | a_k^\dagger a_{k'} | g \rangle &= \langle 0 | \bigotimes_{k''>0} S_{k'',-k''}^\dagger a_k^\dagger a_{k'} \bigotimes_{k''>0} S_{k'',-k''} | 0 \rangle = \\ &= \langle 0 | S_{k,-k}^\dagger a_k^\dagger a_k S_{k,-k} | 0 \rangle = \\ &= \langle g | a_k^\dagger a_k | g \rangle, \end{aligned} \quad (\text{A5})$$

which is valid since, in the case in which $k \neq k' \neq k''$, all the operators commute, and the expectation values are null. Following the same reasoning, if k or k' are equal to 0 or π , the expectation value is always zero unless $k = k'$. It follows that

$$N_{\text{loc}}^j = \langle g | \frac{1}{M} \sum_k a_k^\dagger a_k | g \rangle = N_{\text{loc}}, \quad (\text{A6})$$

which does not depend on the particular site location j . The photons in the system are uniformly distributed across each cavity. We can further simplify (A6) by expanding the discrete sum over $k > 0$:

$$\begin{aligned} N_{\text{loc}} = \frac{1}{M} \langle g | a_0^\dagger a_0 + \delta_{N,2m} a_\pi^\dagger a_\pi | g \rangle + \\ + \frac{1}{M} \langle g | \sum_{k>0} (a_k^\dagger a_k + a_{-k}^\dagger a_{-k}) | g \rangle. \end{aligned} \quad (\text{A7})$$

From Eq. (19), the photon number in each collective mode a_k is $N_k = \sinh^2(|\xi_k|) = s_k^2$, where s_k is defined as in (7) with the substitution $\epsilon \rightarrow \epsilon \cos(k)$. Finally, since $s_k^2 = s_{-k}^2$, we write the local number of photons as

$$N_{\text{loc}} = \frac{1}{M} \sum_k N_k = \frac{1}{M} \left(s_0^2 + \delta_{M,2m} s_\pi^2 + \sum_{k>0} 2s_k^2 \right). \quad (\text{A8})$$

Appendix B: Series expansions

In this section, we present the mathematical details concerning the trigonometric series that appear when computing the QFI and N_{loc} in the asymptotic limit $x \rightarrow 0$ (see Sec. IV A). Let us first consider the QFI when M is odd. Expanding the QFI (22) up to the first order in x yields

$$\begin{aligned} \mathcal{I}_0 = \frac{1}{8\omega^2} \left(\frac{1}{x^2} - \frac{1}{x} - \frac{1}{4} \right) + \\ + \frac{1}{\omega^2} \sum_{n=1}^{\frac{M-1}{2}} \left[\csc^4 \left(\frac{2\pi n}{M} \right) - \csc^2 \left(\frac{2\pi n}{M} \right) \right] + o(x). \end{aligned} \quad (\text{B1})$$

A closed analytical expression for the QFI in this limit can be obtained using the following identities on trigonometric series [68]:

$$\sum_{n=1}^{\frac{M-1}{2}} \csc^2 \left(\frac{2n\pi}{M} \right) = \frac{M^2 - 1}{6}, \quad (\text{B2})$$

$$\sum_{n=1}^{\frac{M-1}{2}} \csc^4 \left(\frac{2n\pi}{M} \right) = \frac{M^4 + 10M^2 - 11}{90}. \quad (\text{B3})$$

Thus, in our case we have

$$\sum_{n=1}^{\frac{M}{2}-1} \left[\csc^4\left(\frac{2n\pi}{M}\right) - \csc^2\left(\frac{2n\pi}{M}\right) \right] = \frac{M^4 - 5M^2 + 4}{90}, \quad (\text{B4})$$

which then leads to (24). On the other hand, when M is even, expanding the QFI (21) leads to

$$\begin{aligned} \mathcal{I}_E &= \frac{1}{4\omega^2} \left(\frac{1}{x^2} - \frac{1}{x} - \frac{1}{4} \right) + \\ &+ \frac{1}{\omega^2} \sum_{n=1}^{\frac{M}{2}-1} \left[\csc^4\left(\frac{2\pi n}{M}\right) - \csc^2\left(\frac{2\pi n}{M}\right) \right] + o(x). \end{aligned} \quad (\text{B5})$$

From the symmetry properties of the cosecant functions and their periodicity, it follows

$$\sum_{n=1}^{\frac{M}{2}-1} \left[\csc^4\left(\frac{2n\pi}{M}\right) - \csc^2\left(\frac{2n\pi}{M}\right) \right] = \frac{M^4 - 20M^2 + 64}{720}, \quad (\text{B6})$$

which then leads to (23).

Let us now focus on the average number of photons. For an even M , expanding (20) up to the first order in x yields

$$N = \frac{1}{\sqrt{2x}} - \frac{M}{2} + \sum_{n=1}^{\frac{M}{2}-1} \csc\left(\frac{2n\pi}{M}\right) + o(\sqrt{x}). \quad (\text{B7})$$

According to Ref. [69] it is

$$\sum_{n=1}^{\frac{M}{2}-1} \csc\left(\frac{2n\pi}{M}\right) = \frac{M}{\pi} \left[\gamma + \log\left(\frac{M}{\pi}\right) \right] + o\left(\frac{1}{M}\right), \quad (\text{B8})$$

where γ is the Euler-Mascheroni constant. Thus, we obtain Eq. (25).

Conversely, for an odd M , expanding in series (20) results in

$$N = \frac{1}{\sqrt{8x}} - \frac{M}{2} + \sum_{n=1}^{\frac{M-1}{2}} \csc\left(\frac{2n\pi}{M}\right) + o(\sqrt{x}). \quad (\text{B9})$$

Due to the symmetry properties of the cosecant function and its periodicity, we have

$$\sum_{n=1}^{\frac{M-1}{2}} \csc\left(\frac{2n\pi}{M}\right) = \frac{M}{\pi} \left[\gamma + \log\left(\frac{2M}{\pi}\right) \right] + o\left(\frac{1}{M}\right), \quad (\text{B10})$$

from which we retrieve Eq. (26).

Appendix C: Finite non-linearity: single-mode

In this section, we go through the details of the perturbation theory applied to assess the validity of the Gaussian approximation.

The squeezing Hamiltonian (3) is an approximation of the general model:

$$H = \omega a^\dagger a + \frac{\epsilon}{2} (a^2 + a^{\dagger 2}) + \chi a^\dagger a^\dagger aa, \quad (\text{C1})$$

with χ representing the Kerr non-linearity. The Gaussian approximation is obtained by setting $\chi = 0$. However, when the number of photons in the system state increases beyond a certain limit, dependent on χ and ω , the nonlinear term becomes significant and cannot be neglected. We take a perturbative approach to identify an upper bound on the allowed number of photons, and, consequently, estimate the saturation of the QFI.

Let us start by analysing the energy gap ΔE between the ground and the first excited state. Up to the first order in χ , it is defined as

$$\Delta E = \Delta E^{(0)} + \chi \Delta E^{(1)} + o(\chi^2), \quad (\text{C2})$$

where $\Delta E^{(0)} = \lambda = \sqrt{\omega^2 - \epsilon^2}$ and $\Delta E^{(1)} = \langle e | H_{\text{NL}} | e \rangle - \langle g | H_{\text{NL}} | g \rangle$. Here, $|g\rangle = S(\xi)|0\rangle$ is the ground state, $|e\rangle = S(\xi)|1\rangle$ is the first excited state and $H_{\text{NL}} = a^\dagger a^\dagger aa$. We recall that the squeezing parameter is defined as $\xi = \ln(s + \sqrt{s^2 + 1})$, with s given by (7). Since $|g\rangle$ and $|e\rangle$ are single-mode squeezed Fock states, $\Delta E^{(1)}$ can be easily computed as

$$\begin{aligned} \Delta E^{(1)} &= \langle e | H_{\text{NL}} | e \rangle - \langle g | H_{\text{NL}} | g \rangle = \\ &= (9N + 15N^2) - (N + 3N^2) = \\ &= 8N + 12N^2, \end{aligned} \quad (\text{C3})$$

with $N = \sinh^2(|\xi|)$ the number of photons in the ground state. In the Gaussian approximation, we have $\Delta E = \Delta E^{(0)}$. Thus, when $\chi > 0$, it must hold that

$$\frac{\chi \Delta E^{(1)}}{\Delta E^{(0)}} = \frac{\chi (8N + 12N^2)}{\lambda} \ll 1. \quad (\text{C4})$$

By expressing λ in terms of N as $\lambda = \omega/(1 + 2N)$, we obtain an estimate of the desired upper bound on the number of photons in the ground state, i.e.,

$$N \lesssim \sqrt[3]{\frac{\omega}{24\chi}}. \quad (\text{C5})$$

This means that, as far as N is well below the upper bound (C5), the Gaussian approximation is valid and the non-linearity can be safely neglected.

The same procedure may be applied directly to the QFI. In the following, we show that this leads to an upper bound with the same scaling in ω/χ as in (C5) but slightly tighter. To apply the perturbation theory to the QFI, we use (8) with the ground state defined as

$$|g\rangle = |g^{(0)}\rangle + \chi |g^{(1)}\rangle + o(\chi^2). \quad (\text{C6})$$

Here, $|g^{(0)}\rangle = S(\xi)|0\rangle$ is the unperturbed ground state, while $|g^{(1)}\rangle$ is the first-order correction to the unperturbed ground state. Up to the first order in χ , the QFI can be expressed as

$$\mathcal{I} = \mathcal{I}^{(0)} + \chi \mathcal{I}^{(1)} + o(\chi^2), \quad (\text{C7})$$

where $\mathcal{I}^{(0)}$ is the unperturbed QFI defined in (9), while $\mathcal{I}^{(1)} = 8 \text{Re} [\langle \partial_\omega g^{(0)} | \partial_\omega g^{(1)} \rangle]$ is the first order correction. Perturbation theory allows us to easily compute $|g^{(1)}\rangle$. Indeed, in our case

$$|g^{(1)}\rangle = \sum_{j=1}^4 \frac{\langle j^{(0)} | H_{\text{NL}} | g^{(0)} \rangle}{-j\lambda} |j^{(0)}\rangle, \quad (\text{C8})$$

with $|j^{(0)}\rangle = S(\xi) |j\rangle$. The sum is performed for $j \in [1, 4]$, since when $j > 4$ all the expectation values are null due to the orthogonality of the Fock states. The transition elements in (C8) can be then analytically evaluated as $|j^{(0)}\rangle$ are single-mode squeezed Fock states. The first order correction to the ground state is then

$$|g^{(1)}\rangle = v_{\omega,\epsilon} S(\xi) |2\rangle + h_{\omega,\epsilon} S(\xi) |4\rangle. \quad (\text{C9})$$

The amplitude probability functions $v_{\omega,\epsilon}$ and $h_{\omega,\epsilon}$ are defined as

$$\begin{aligned} v_{\omega,\epsilon} &= \frac{(1+6N)\sqrt{2(N+N^2)}}{2\lambda}, \\ h_{\omega,\epsilon} &= -\frac{\sqrt{24}(N+N^2)}{4\lambda}, \end{aligned} \quad (\text{C10})$$

and depend on ω and ϵ according to $N = N(\omega, \epsilon)$ and $\lambda = \lambda(\omega, \epsilon)$. If we now rewrite the squeezing operator as $S(\xi) = e^{f_{\omega,\epsilon} O}$, where $f_{\omega,\epsilon} = |\xi|/2$ and $O = a^2 - a^{\dagger 2}$, the derivatives with respect to the parameter ω read

$$\begin{aligned} |\partial_\omega g^{(0)}\rangle &= f'_{\omega,\epsilon} O |g^{(0)}\rangle, \\ |\partial_\omega g^{(1)}\rangle &= (v'_{\omega,\epsilon} + v_{\omega,\epsilon} f'_{\omega,\epsilon} O) |2^{(0)}\rangle + \\ &\quad + (h'_{\omega,\epsilon} + h_{\omega,\epsilon} f'_{\omega,\epsilon} O) |4^{(0)}\rangle, \end{aligned} \quad (\text{C11})$$

where the apex stands for ∂_ω . Finally, using (C11), we are able to compute the first-order correction to the QFI:

$$\begin{aligned} \mathcal{I}^{(1)} &= 8 \langle \partial_\omega g^{(0)} | \partial_\omega g^{(1)} \rangle = \\ &= -8\sqrt{2} f'_{\omega,\epsilon} \left(v'_{\omega,\epsilon} + \sqrt{12} f'_{\omega,\epsilon} h_{\omega,\epsilon} \right). \end{aligned} \quad (\text{C12})$$

From the definition $N = s^2$, we obtain an expression for ϵ in terms of the unperturbed ground state number of photons, i.e., $\epsilon = 2\omega\sqrt{N+N^2}/(1+2N)$. With this in mind, the leading terms of the zeroth-order QFI and its first-order correction are

$$\mathcal{I}^{(0)} \sim \frac{8N^4}{\omega^2}, \quad \mathcal{I}^{(1)} \sim -\frac{1056N^7}{\omega^3}. \quad (\text{C13})$$

In the Gaussian approximation, it holds that $\mathcal{I} = \mathcal{I}^{(0)}$. Thus, when $\chi > 0$, we have to impose $|\chi \mathcal{I}^{(1)}| \ll |\mathcal{I}^{(0)}|$, from which we obtain a new estimate for the upper bound on the number of photons in the ground state:

$$N \lesssim \sqrt[3]{\frac{\omega}{132\chi}}. \quad (\text{C14})$$

This bound has the same scaling in ω/χ as that in (C5), but is tighter by roughly a factor of two.

Appendix D: Finite non-linearity: resonator chain

For the multi-mode case, the Hamiltonian in (16) can be thought of as the Gaussian approximation of the general model

$$H = \sum_{j=1}^M \left[\omega a_j^\dagger a_j + \frac{\epsilon}{2} (a_j a_{j+1} + a_j^\dagger a_{j+1}^\dagger) + \chi a_j^\dagger a_j^\dagger a_j a_j \right], \quad (\text{D1})$$

where χ represents the local Kerr non-linearity. To assess the validity of the Gaussian approximation, we can apply perturbation theory, as we did for the single-mode case. In the following, we consider M odd, since this choice leads to the most favorable scaling of the QFI. Close to the criticality, the majority of the photons are then confined in the mode a_0 , leaving the other modes $a_{\pm k}$ unpopulated (see Sec. IV). Thus, we approximate the unperturbed ground state defined in (19) as

$$|g^{(0)}\rangle \approx S_0(\xi_0) |0\rangle. \quad (\text{D2})$$

This allows us to map the multi-mode system to a single-mode system. The low-energy excited states can be written as $|j^{(0)}\rangle \approx (j!)^{-\frac{1}{2}} S_0(\xi_0) (a_0^\dagger)^j |0\rangle$. The single-mode system has now $N = MN_{\text{loc}}$ photons in the mode a_0 , where N_{loc} is the local number of photons in each site j . Let us now focus on the effect of the nonlinear terms, which can be written in the reciprocal space as

$$\begin{aligned} \chi H_{\text{NL}} &= \chi \sum_j^M a_j^\dagger a_j^\dagger a_j a_j = \\ &= \frac{\chi}{M^2} \sum_j^M \sum_{k_1, k_2, k_3, k_4} a_{k_1}^\dagger a_{k_2}^\dagger a_{k_3} a_{k_4} e^{-i(k_3+k_4-k_1-k_2)j} = \\ &= \frac{\chi}{M} \sum_{k_1, k_2, k_3} a_{k_1}^\dagger a_{k_2}^\dagger a_{k_3} a_{k_1+k_2-k_3}. \end{aligned} \quad (\text{D3})$$

Given the structure of the low-excited states, the expected value and transition elements of the nonlinear term are nonvanishing only for $k_1 = k_2 = k_3 = k_4 = 0$. In this single-mode approximation, the Hamiltonian (D1) can be mapped onto a single critical resonator with a rescaled nonlinear term, such that

$$\chi H_{\text{NL}} \longrightarrow \frac{\chi}{M} a_0^\dagger a_0^\dagger a_0 a_0. \quad (\text{D4})$$

We can then perform the same calculations we did for the perturbation of the QFI in the single-mode case, but with the ground state (D2) and a rescaled nonlinear Hamiltonian. It follows that the bounds found for the single-mode case are still valid for the multi-mode one, provided that we replace $N \rightarrow MN_{\text{loc}}$ and $\chi \rightarrow \chi/M$. Finally, to ensure the validity of the Gaussian approximation in the multi-mode scenario, we require

$$N \lesssim \sqrt[3]{\frac{\omega M}{132\chi}}. \quad (\text{D5})$$

So far, we have assumed that the unperturbed multi-mode ground state can be written as (D2). This allowed us to evaluate the first-order correction to the QFI. In this section, we analytically compute the first-order correction to the multi-mode energy gap, without resorting to Eq. (D2). In the limit where $N \approx N_0$, with N the number of photons in the ground state and N_0 the number of photons in the mode a_0 , the single-mode bound (C5) is retrieved, thus endorsing the validity of the single-mode approximation. The multi-mode energy gap between the ground and the first excited state is defined as in (C2), with $\Delta E^{(0)} = \lambda_0 = \sqrt{\omega^2 - \epsilon^2}$ and $\Delta E^{(1)} = E_e^{(1)} - E_g^{(1)}$, where $E_e^{(1)}$ and $E_g^{(1)}$ are the first order corrections to the eigenvalues of the first-excited and ground state, respectively. Let us start by computing the first-order correction to the ground state $E_g^{(1)}$. By using (D3) and the relabeling $k_1 \rightarrow k' + q$, $k_2 \rightarrow k - q$, $k_3 \rightarrow k'$, we write $E_g^{(1)}$ as

$$\begin{aligned} E_g^{(1)} &= \frac{1}{M} \sum_{k,k',q} \langle g | a_{k'+q}^\dagger a_{k-q}^\dagger a_{k'} a_k | g \rangle = \\ &= \frac{1}{M} \sum_{k,k'} \langle g | a_{k'+q}^\dagger a_{k'} a_{k-q}^\dagger a_k | g \rangle - N, \end{aligned} \quad (\text{D6})$$

where we have used $[a_{k'}, a_{k-q}^\dagger] = \delta_{k', k-q}$ and $MN = \sum_{k,q} \langle g | a_k^\dagger a_k | g \rangle$. Since the ground state $|g\rangle$ is defined as in (19), any expectation value $\langle g | a_k | g \rangle$ is null. This means that the non-zero terms in (D6) are those in which the ladder operators refer to the same mode in pairs, leading to the constraints: (i) $k' + q = k'$ and $k - q = k$, meaning $q = 0$; (ii) $k' + q = k - q$ and $k' = k$, meaning $q = 0$ and $k = k'$; (iii) $k' + q = k$, meaning $q = k - k'$. Implementing these constraints in (D6) allows us to obtain a much simpler expression:

$$\begin{aligned} E_g^{(1)} &= \frac{1}{M} \left[2 \sum_{k,k'} \langle g | a_k^\dagger a_{k'} a_k^\dagger a_k | g \rangle + \right. \\ &\quad \left. - \sum_k \langle g | a_k^\dagger a_k a_k^\dagger a_k | g \rangle - N \right]. \end{aligned} \quad (\text{D7})$$

On the other hand, the first excited state in the multi-mode system is

$$|e\rangle = S(\xi_0) |1\rangle \bigotimes_{k>0} S_{k,-k}(\xi_k) |0_k\rangle. \quad (\text{D8})$$

Thus, the expression (D7) holds also for $E_e^{(1)}$, provided

that $|g\rangle \rightarrow |e\rangle$ and $N \rightarrow N+1+2N_0$ (it is straightforward to verify that the number of photons in $|e\rangle$ is indeed $N+1+2N_0$). The first-order correction to the multi-mode energy gap is then

$$\begin{aligned} \Delta E^{(1)} &= E_e^{(1)} - E_g^{(1)} = \\ &= \frac{1}{M} \left[2 \sum_{k,k'} \left(\langle e | a_{k'}^\dagger a_{k'} a_k^\dagger a_k | e \rangle - \langle g | a_{k'}^\dagger a_{k'} a_k^\dagger a_k | g \rangle \right) + \right. \\ &\quad \left. - \sum_k \left(\langle e | a_k^\dagger a_k a_k^\dagger a_k | e \rangle - \langle g | a_k^\dagger a_k a_k^\dagger a_k | g \rangle \right) + \right. \\ &\quad \left. - 1 - 2N_0 \right]. \end{aligned} \quad (\text{D9})$$

Since the first excited state $|e\rangle$ differs from ground state $|g\rangle$ only in the terms related to the mode a_0 , the differences between expectation values in (D9) will always give zero unless $k = 0$ or $k' = 0$. Then, we can further simplify $\Delta E^{(1)}$:

$$\begin{aligned} \Delta E^{(1)} &= \frac{1}{M} \left[4(1+2N_0)(N-N_0) + \right. \\ &\quad \left. + \langle e | a_0^\dagger a_0 a_0^\dagger a_0 | e \rangle - \langle g | a_0^\dagger a_0 a_0^\dagger a_0 | g \rangle + \right. \\ &\quad \left. - 1 - 2N_0 \right]. \end{aligned} \quad (\text{D10})$$

The remaining expectation values in (D10) can be computed easily, being $|e\rangle$ and $|g\rangle$ single-mode squeezed Fock states for what concerns a_0 . We have that $\langle e | a_0^\dagger a_0 a_0^\dagger a_0 | e \rangle = 1+12N_0+15N_0^2$ and $\langle g | a_0^\dagger a_0 a_0^\dagger a_0 | g \rangle = 2N_0+3N_0^2$. Finally, we write

$$\Delta E^{(1)} = \frac{1}{M} [4N + 8N_0N + 4N_0 + 4N_0^2], \quad (\text{D11})$$

and, up to the first order in χ , the multi-mode energy gap is given by

$$\Delta E = \lambda_0 + \frac{\chi}{M} [4N + 8N_0N + 4N_0 + 4N_0^2] + o(\chi^2). \quad (\text{D12})$$

Close to the critical point, we have $N = MN_{\text{loc}} \approx N_0$, so that

$$\Delta E \approx \lambda_0 + \frac{\chi}{M} [8N + 12N^2]. \quad (\text{D13})$$

We retrieve the same result obtained in the single-mode case in (C5), with a larger number of photons $N = MN_{\text{loc}}$ and a rescaled non-linearity χ/M .

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