

Theory of Subradiant States of a One-Dimensional Two-Level Atom Chain

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Recently, the subradiant states of one-dimensional two-level atom chains coupled to light modes were found to have decay rates obeying a universal scaling, and an unexpected fermionic character of the multiply excited subradiant states was discovered. In this Letter, we theoretically obtain the singly excited subradiant states, and by eliminating the superradiant modes, we demonstrate a relation between the multiply excited subradiant states and the Tonks-Girardeau limit of the Lieb-Liniger model which explains the fermionic behavior. In addition, we identify a new family of subradiant states with correlations different from the fermionic ansatz.

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To achieve controllable and deterministic photon-atom interfaces for applications in quantum information processing and quantum sensing, large atom ensembles may be used to enhance the coupling to photons [1]. The photons induce both coherent and dissipative atom-atom interactions that yield collective phenomena of super- or subradiance [2], wherein a collective excitation of the atom ensemble decays faster or slower than individual atomic excitations. While superradiance has been extensively studied since the seminal work of Dicke [3], subradiance of a large ensemble was observed only very recently in cold atom clouds [4,5] and metamaterial arrays [6]. Comprehensive theoretical tools for the subradiance are still elusive [7–10] due to the complicated long-range interactions and many-body features of the atomic ensembles [11–13]. A one-dimensional (1D) chain of equally spaced two-level atoms offers the simplest geometry to gain insight in the collective decay mechanisms, and implementation of such chains coupled to nanofibers [14], 1D waveguides [15–18], and the full vacuum electromagnetic field in 3D free space [8,13,19–21] has attracted considerable attention. Super- and subradiance phenomena are in these systems supplemented by further interesting properties and applications such as atomic mirrors [22], photon Fock state synthesis [23], enhancement of cooperativity [24], and applications in quantum computation [25].

Recently, the subradiant states of such 1D chains of N qubits in 3D free space and coupled to 1D waveguide were numerically found to have a series of seemingly universal properties [8–10]: in the one-excitation sector where only one of the N atoms is excited, if we sort all eigenstates (to be elaborated) by increasing decay rates with integer labels from $\xi = 1$ to $\xi = N$, the most subradiant states ($\xi \ll N$) have decay rates $\gamma_\xi \propto \xi^2/N^3$. In the multiexcitation sectors, the most subradiant states have a fermionic character, e.g., a subradiant state with two excitations is given by

$|F_{1,2}\rangle \propto \sum_{i<j} (c_{1,i}c_{2,j} - c_{1,j}c_{2,i})|e_i, e_j\rangle$, built from subradiant states $|\psi_{1(2)}\rangle = \sum_i c_{1(2),i}|e_i\rangle$ in the one-excitation sector, where $|e_i\rangle$ ($|e_i, e_j\rangle$) represents the state with the i th (i th and the j th) atom excited to $|e\rangle$ while all other atoms are in the ground state $|g\rangle$. The decay rate of $|F_{1,2}\rangle$ is the sum of the decay rates of $|\psi_1\rangle$ and $|\psi_2\rangle$. The infidelity of the fermionic ansatz $|F_{1,2}\rangle$ to exact numerical results scale as N^{-2} and N^{-1} for two different classes of states [9].

It is intriguing why these properties appear for both the infinite range atom-atom interactions mediated by the 1D guided fields [9,10], and the long-range ($\sim 1/r$) or short-range ($\sim 1/r^3$) interactions mediated by the 3D free-space field [8,10]. A thorough theoretical understanding is needed to guide further experimental studies and applications of subradiance. In this Letter, we provide such understanding based on the theoretical treatment of the physics summarized in Fig. 1.

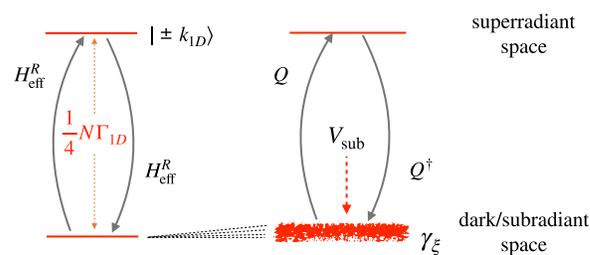


FIG. 1. Outline of the theory: the imaginary part of the effective atomic Hamiltonian, H_{eff}^I , identifies the superradiant ($|\pm k_{1D}\rangle$) and dark subspaces of states (left panel), coupled perturbatively by H_{eff}^R to produce the subradiant modes with decay rates $\gamma_\xi \propto \xi^2/N^3$ ($\xi \ll N$). The Holstein-Primakoff transformation bosonizes the super- and subradiant modes and introduces a coupling (Q and Q^\dagger) between them (right panel). The coupling effectively yields a strong interaction V_{sub} among the multiply excited subradiant modes leading to the formation of states with fermionic character.

Spin model.—For light-matter interactions where the Markovian approximation is applicable, we can eliminate the light modes to obtain a master equation describing only the atoms [26]. The master equation is equivalent to a Monte Carlo wave function formalism [27], where the atomic state evolves stochastically under quantum jumps and deterministically under $H = H_0 + H_{\text{eff}}$, where H_0 is the bare Hamiltonian of the atoms and the *non-Hermitian* H_{eff} describes both coherent and dissipative atom-atom interactions mediated by the vacuum field. The right eigenstates of H , or equivalently of H_{eff} [8], in each manifold of states with any given number of atomic excitations have decay rates that are twice the imaginary parts of the corresponding right eigenvalues. We focus our analysis on the qubit chain coupled to a 1D waveguide, but our treatment provides sufficient insight to also account for the case of coupling to the 3D vacuum field. For an atom chain coupled to a 1D waveguide, we have [9,22]

$$H_{\text{eff}} = -\frac{i}{2}\Gamma_{\text{1D}} \sum_{m,n=1}^N e^{ik_{\text{1D}}|z_m - z_n|} \sigma_m^\dagger \sigma_n, \quad (1)$$

where Γ_{1D} is the decay rate of a single atom coupled to the waveguide [22], k_{1D} is the wave number of the waveguide mode resonant with the atomic transition, and $\sigma_m = |g\rangle_m \langle e|$ acts on the m th atom. We assume the atoms are equidistantly spaced by d . For convenience, we shall denote $H_{\text{eff}} = H_{\text{eff}}^R - iH_{\text{eff}}^I$.

One-excitation sector.—The eigenstates of H_{eff}^I divide the one-excitation sector into a two-dimensional superradiant subspace spanned by the Bloch states $|\pm k_{\text{1D}}\rangle = N^{-1/2} \sum_{m=1}^N e^{\pm ik_{\text{1D}}z_m} |e_m\rangle$ with eigenvalue $N\Gamma_{\text{1D}}/4$ and an $(N-2)$ -dimensional dark space with an eigenvalue of 0; see Fig. 1. The dark states acquire weak (subradiant) decay rates because of their admixture of superradiant states induced by the perturbation from H_{eff}^R .

While the perturbation view is informative, a more direct approach to the subradiant states applies the following exact result for the Bloch states $|k\rangle$ ($k \neq \pm k_D$)

$$H_{\text{eff}}|k\rangle = \omega_k|k\rangle - i\frac{\Gamma_{\text{1D}}}{2}(g_k|k_{\text{1D}}\rangle - h_k| - k_{\text{1D}}\rangle), \quad (2)$$

where $\omega_k = (\Gamma_{\text{1D}}/4) \sum_{\epsilon=\pm} \cot[(k_{\text{1D}} + \epsilon k)d/2]$, and the ‘‘tails’’ $g_k = [e^{i(k-k_{\text{1D}})z_1}]/[1 - e^{i(k-k_{\text{1D}})d}]$ and $h_k = [e^{i(k+k_{\text{1D}})z_N}]/[e^{-i(k+k_{\text{1D}})d} - 1]$. It follows that a superposition of two degenerate states, $|k\rangle$ and $| - k\rangle$, is an eigenstate of H_{eff} with eigenvalue ω_k and has no tails if k is a solution to the equation $g_k h_{-k} = g_{-k} h_k$. This equation has only solutions for complex values of k . In the regimes $k \approx 0$ or $\pm\pi/d$ (center or edges of the first Brillouin zone), supposing $k_\xi = 0 + \delta_\xi$ and $k_\xi = -\pi/d + \delta_\xi$, respectively, we find to the order of N^{-2}

$$\delta_\xi = \frac{\xi\pi}{Nd} \times \begin{cases} 1 - i\frac{1}{N} \cot(\frac{k_{\text{1D}}}{2}d), & k \approx 0 \\ 1 + i\frac{1}{N} \tan(\frac{k_{\text{1D}}}{2}d), & k \approx -\pi/d, \end{cases} \quad (3)$$

with $\xi = 1, 2, 3, \dots, \xi \ll N$. Note that Eq. (3) amounts to an $1/N^2$ -order imaginary correction to the real Bloch wave number.

Next, we substitute Eq. (3) into the expression for ω_k , which is parabolic near $k \approx 0$ and $\pm\pi/d$, i.e., $\omega_k \propto \delta_\xi^2$. Then the imaginary corrections directly yield the ξ^2/N^3 scaling of the decay rates [9]

$$\gamma_\xi = \Gamma_{\text{1D}} \frac{\pi^2}{2} \frac{\xi^2}{N^3} \times \begin{cases} \frac{\cos^2(k_{\text{1D}}d/2)}{\sin^4(k_{\text{1D}}d/2)}, & k \approx 0 \\ \frac{\sin^2(k_{\text{1D}}d/2)}{\cos^4(k_{\text{1D}}d/2)}, & k \approx -\pi/d. \end{cases} \quad (4)$$

The eigenstates are written as

$$\begin{aligned} |\phi_{k_\xi}\rangle &\propto g_{-k_\xi}|k_\xi\rangle - g_{k_\xi}| - k_\xi\rangle \\ &= \frac{1}{\sqrt{2}}(|k_\xi^{(0)}\rangle - | - k_\xi^{(0)}\rangle) + O\left(\frac{\xi}{N}\right), \end{aligned} \quad (5)$$

where $k_\xi^{(0)} = \xi\pi/(Nd)$ or $-\pi/d + \xi\pi/(Nd)$.

Universality.—The ξ^2/N^3 scaling has also been numerically found for 1D atom chains coupled to 3D free-space modes [8–10], where the effective Hamiltonian $H_{\text{3D,eff}}$ is determined by the vacuum Green’s dyadic tensor [28]. Fourier transformation of the Green’s tensor reveals a hidden similarity between the coupling to the 1D and 3D quantized radiation fields: $H_{\text{3D,eff}}$ can be written as weighted integrals of terms resembling H_{eff} with real-valued $k_{\text{1D}} \in [0, k_0]$ and imaginary-valued $k_{\text{1D}} \in [i0, +i\infty]$

$$\begin{aligned} H_{\text{3D,eff}} &= -i\frac{3\gamma_0}{4k_0} \int_0^{k_0} \frac{d\tilde{k}}{2\pi} \rho_+(\tilde{k}) \sum_{m,n=1}^N e^{i\tilde{k}|z_m - z_n|} \sigma_m^\dagger \sigma_n \\ &\quad - \frac{3\gamma_0}{4k_0} \int_0^{+\infty} \frac{d\tilde{k}}{2\pi} \rho_-(\tilde{k}) \sum_{m,n=1}^N e^{-\tilde{k}|z_m - z_n|} \sigma_m^\dagger \sigma_n, \end{aligned} \quad (6)$$

where γ_0 is the spontaneous emission rate and k_0 is the resonant wave number. If the atoms are polarized parallel to the chain, $\rho_\pm(\tilde{k}) = 2\pi(1 \mp \tilde{k}^2/k_0^2)$ and the atom-atom interaction is short-range ($\sim 1/r^3$). If the atoms are polarized transverse to the chain, $\rho_\pm(\tilde{k}) = \pi(1 \pm \tilde{k}^2/k_0^2)$ and the atom-atom interaction is long-range ($\sim 1/r$).

In combination with the two exact features of our analytical results for H_{eff} , (1) the leading order solutions of δ_ξ and $|\phi_{k_\xi}\rangle$ are independent of the values of k_{1D} ; see Eq. (5). (2) the proportionality $\delta_\xi \propto \xi$ and the parabolic dispersion relation $\omega_k \propto \delta_\xi^2$ hold to order N^{-2} , for all values of k_{1D} . This explains the universality of the ξ^2/N^3 scaling: feature 1 implies that the leading order eigenstates of H_{eff} , shown in Eq. (5), are shared simultaneously by all terms

integrated in $H_{3D,eff}$, and thus by the full $H_{3D,eff}$ due to linearity. Feature 2, hence implies that the corresponding decay rates, scaling as ξ^2/N^3 , also apply to the subradiant states of $H_{3D,eff}$. The prerequisite is that $H_{3D,eff}^I$ must have dark states with $k \approx \pm\pi/d$. Hence, we require $k_0 < \pi/d$ which implies that the ensemble is only subradiant in the 3D field if the atom-atom distance is less than half the resonant wavelength [8,13].

Subradiant multiply excited states.—When the number of atomic excitations $n_e \ll N$, the leading order Holstein-Primakoff (HP) approximation [29] usually applies and one may replace $\sigma_m^\dagger \sigma_n$ of Eq. (1) with the bosonic operators $b_m^\dagger b_n$ and obtain a quadratic bosonic H_{eff} . It works well for the superradiant modes with wave number $\pm k_{1D}$. But for the subradiant multiply excited states, the bosonic creation operators prepare exchange symmetric combinations of subradiant one-excitation states with decay rates scaling as N^{-1} [8,9] which is much larger than the numerically observed N^{-3} scaling [8–10]. Instead, the numerical results were found to favor fermionic exchange antisymmetric combinations of the subradiant one-excitation states [8–10].

This somewhat surprising result inspires a closer scrutiny of the HP transformation. Including second order corrections due to saturation, the HP transformation reads $\sigma_m = (1 - b_m^\dagger b_m/2)b_m$ so that we can write $H_{eff}^I = H_{SR} + Q + Q^\dagger$ with the quadratic term $H_{SR} = N\Gamma_{1D}/4 \sum_{\epsilon=\pm} b_{\epsilon k_{1D}}^\dagger b_{\epsilon k_{1D}}$ and the quartic term

$$Q = -\frac{\Gamma_{1D}}{8} \sum_{\epsilon=\pm} \sum_{p,q} b_{\epsilon k_{1D}}^\dagger b_{p+q-\epsilon k_{1D}}^\dagger b_p b_q. \quad (7)$$

Here, $b_k^\dagger = N^{-1/2} \sum_m e^{ikz_m} b_m^\dagger$ and the summation over wave number is taken over an orthonormal basis $\{|k\rangle\}_k$ containing $|\pm k_{1D}\rangle$. The quadratic term H_{SR} has a prefactor N times larger than those of Q and Q^\dagger , but to assess their influence, we should take account of not only the prefactors but also the magnitudes of the operator terms. For H_{SR} , the magnitude of $b_{\epsilon k_{1D}}^\dagger b_{\epsilon k_{1D}}$ can be estimated by its typical expectation value over the relevant Hilbert space, i.e., the subradiant states.

Reasonably, one may expect that a subradiant state contains no excitation of superradiant modes, i.e., typically $\langle b_{\epsilon k_{1D}}^\dagger b_{\epsilon k_{1D}} \rangle \approx 0$. Thus, the magnitude of H_{SR} is suppressed. Meanwhile, Eq. (7) shows that Q annihilates a two-boson state $b_p^\dagger b_q^\dagger |\emptyset\rangle$, which is dark with respect to H_{SR} , and generates a superradiant two-boson state $b_{\epsilon k_{1D}}^\dagger b_{p+q-\epsilon k_{1D}}^\dagger |\emptyset\rangle$ where $|\emptyset\rangle$ denotes the vacuum state. This demonstrates that the saturation correction to the first order HP transformation plays a significant role even in the low excitation regime ($n_e \ll N$), in contrast to its role in many other applications.

An effective theory for how Q couples the dark states to superradiant states and hereby determines their subradiant behavior is illustrated in the right panel of Fig. 1. The effect

of Q and Q^\dagger is distilled by eliminating the superradiant states, in a manner similar to the adiabatic elimination of excited state manifolds to restrict the effective dynamics of quantum systems to their ground state manifold [30].

Note that the subset of superradiant states with only a single excitation of the superradiant modes and thus the eigenvalue (decay rate) $N\Gamma_{1D}/4$ has the strongest coupling to the dark or subradiant states. We hence disregard the coupling to other superradiant states and the effective coupling between the subradiant states reduces to $V_{sub} = (4/N\Gamma_{1D})P_{DS}Q^\dagger P_{SRS}QP_{DS}$, with projection operators $P_{DS(SRS)}$ on the dark and superradiant spaces, respectively. To evaluate this expression, we use the operator relation that $(b_{p'+q'-\epsilon'k_{1D}} b_{\epsilon'k_{1D}})^\dagger (b_{\epsilon k_{1D}}^\dagger b_{p+q-\epsilon k_{1D}}^\dagger) = \delta_{\epsilon,\epsilon'} \delta_{p'+q',p+q}$, i.e., no population of the superradiant two-boson modes within the dark or subradiant space. Finally, we obtain

$$\begin{aligned} V_{sub} &= \frac{1}{8N} \Gamma_{1D} \sum_{p,q,k} b_{-p+q+k}^\dagger b_p^\dagger b_q b_k \\ &= \frac{1}{8} \Gamma_{1D} \sum_{m=1}^N (b_m^\dagger)^2 (b_m)^2. \end{aligned} \quad (8)$$

That is, V_{sub} induces decay with rate $O(\Gamma_{1D})$ of nominally subradiant states having more than a single HP boson excitation at the same site.

In the absence of V_{sub} , approximate multiply excited states are created by the operators $b_\xi^\dagger = \sum_m \langle e_m | \phi_{k_\xi} \rangle b_m^\dagger$, with $|\phi_{k_\xi}\rangle$ the one-excitation eigenstates Eq. (5) of H_{eff} . Since V_{sub} cannot be treated as a perturbation, we study the effective Hamiltonian $\mathcal{H} = \frac{1}{2} \sum_\xi \gamma_\xi b_\xi^\dagger b_\xi + V_{sub}$, where only the most subradiant states ($\xi \ll N$) are included in the sum. In the continuous limit [28], \mathcal{H} can be written as the second-quantized form of the Hamiltonian

$$\mathcal{H} = \sum_{i=1}^{n_e} \left(\frac{-\partial_{x_i}^2}{2m_*} + V(x_i) \right) + 2c_{LL} \sum_{i<j=1}^{n_e} \delta(x_i - x_j), \quad (9)$$

where $c_{LL} = d\Gamma_{1D}/8$ and $V(x_i)$ is a 1D box potential for bosons in the interval $[0, Nd]$ with one-excitation eigenstates $|\phi_{k_\xi}\rangle$ given by Eq. (5). The observation behind Eq. (9) is that the ξ^2/N^3 scaling of γ_ξ takes the form of a kinetic energy $\gamma_\xi = k_\xi^2/(2m_*)$ when $k \approx 0$; or the kinetic energy in a gauge field $\gamma_\xi = (k_\xi + \pi/d)^2/(2m_*)$ when $k \approx -\pi/d$. With the parametrization of the model, the effective mass in the kinetic energy term reads $m_* = \xi^2 \pi^2 / (N^2 d^2 \gamma_\xi) \propto N$.

We recognize Eq. (9) as the Lieb-Liniger model [31] originally proposed for 1D gases of hard-core bosons. As the effective mass m_* diverges in the large N limit, the kinetic energy-like part of Eq. (9) vanishes. This implies that Eq. (9) reaches the Tonks-Girardeau regime [32,33] of the Lieb-Liniger model, where the eigenstates of \mathcal{H} can be

obtained via a fermion-boson mapping [33,34]: for a free fermion model described by the Hamiltonian $\sum_{\xi} \frac{1}{2} \gamma_{\xi} f_{\xi}^{\dagger} f_{\xi}$, where $f_{\xi}^{\dagger} = \sum_m \langle e_m | \phi_{\xi} \rangle f_m^{\dagger}$, we write down its eigenstates (e.g., two-fermion states) $f_{\xi_1}^{\dagger} f_{\xi_2}^{\dagger} |\emptyset\rangle$, and replace $f_m^{\dagger} f_n^{\dagger}$ with $\text{sign}(n-m) b_m^{\dagger} b_n^{\dagger}$, where $\text{sign}(n-m)$ is necessary to ensure the consistency with the fermionic commutation relation. This yields a fermion-like bosonic state $|F_{\xi_1, \xi_2}\rangle = \sum_{m < n} [\phi_{k_{\xi_1}}(z_m) \phi_{k_{\xi_2}}(z_n) - e^{i\varphi} \phi_{k_{\xi_2}}(z_m) \phi_{k_{\xi_1}}(z_n)] b_m^{\dagger} b_n^{\dagger} |\emptyset\rangle$ ($e^{i\varphi} = 1$ is introduced for later convenience). Because $b_m^{\dagger} b_n^{\dagger} |\emptyset\rangle = |e_m, e_n\rangle$, we recover the fermionic ansatz of the two-excitation sector [8–10]. As a direct consequence of their representation as noninteracting fermions, the decay rates of the most subradiant multiply excited states are merely the sum of the decay rates of their one-excitation constituents, e.g., $\gamma_{\xi_1} + \gamma_{\xi_2}$. This explains the numerical observations of Refs. [8–10]. The above mapping also applies to states with more excitations [33,34].

For a finite atom chain, m_* is finite and Eq. (9) deviates from the Tonks-Girardeau limit; hence the fermionic ansatz, e.g., the two-excitation state $|F_{\xi_1, \xi_2}\rangle$, deviates slightly from the numerical eigenstates denoted by $|\psi_{\xi_1, \xi_2}^{\text{num}}\rangle$. The deviation quantified by the infidelity $1 - |\langle F_{\xi_1, \xi_2} | \psi_{\xi_1, \xi_2}^{\text{num}} \rangle|^2$, is numerically found to scale as N^{-2} when both components, $|\phi_{k_{\xi_1}}\rangle$ and $|\phi_{k_{\xi_2}}\rangle$, come from the same branch of the one-excitation subradiant states, i.e., $k_{\xi_1}, k_{\xi_2} \approx \pm\pi/d$ (or both ≈ 0); otherwise, the infidelity scales as N^{-1} (when $k_{\xi_1} \approx \pm\pi/d$ and $k_{\xi_2} \approx 0$) [9]. These behaviors can also be explained from the Lieb-Liniger model of Eq. (9). The fermion-boson mapping is not exact and the phase factor $e^{i\varphi}$ introduced

above deviates from unity by a factor in the form of $(k_{\xi_1} - k_{\xi_2}) / (m_* c_{LL})$ [31]. Because $k_{\xi_1} - k_{\xi_2}$ is $O(N^{-1})$ or $\approx \pi/d$ in the two cases considered, while $m_* c_{LL}$ scales as N , their ratio scales exactly in the same manner as the numerically observed infidelities [9]. Larger discrepancies with the fermionic ansatz are detectable when the decay rates increase.

Universality.—The mapping to the Lieb-Liniger model and the Tonks-Girardeau gas can also be extended to the 1D atomic chain coupled to 3D free-space modes described by Eq. (6). Here $H_{\text{3D,eff}}^I$ possesses short-lived eigenstates $|\tilde{k}\rangle$ with $\tilde{k} \in [-k_0, k_0]$ and different decay rates $\gamma_{\tilde{k}}$. Each of them will contribute to V_{sub} a term with prefactor $\gamma_{\tilde{k}}/N^2$. Hence, we have $V_{\text{sub}} \propto \sum_{\tilde{k}} \gamma_{\tilde{k}}/N^2 = \gamma_0/N$, similar to the coefficient in the first line of Eq. (8). Because the ξ^2/N^3 scaling decay rates apply in the one-excitation sector of the 3D free space case, the fermionic ansatz also applies here. Because only the branch of subradiant states with $k \approx \pm\pi/d$ appear here, the pertaining N^{-2} scaling applies to the infidelities of all states given by the fermionic ansatz. This matches the numerical results [8].

Conclusion and discussion.—In this Letter, we have developed a theory to explain the ξ^2/N^3 scaling of subradiant decay rates and the fermionic behavior of multiply excited subradiant states identified in numerical calculations on 1D atom chains coupled to both 1D and 3D radiation reservoirs [8–10]. We find that the universal ξ^2/N^3 scaling results from a parabolic dispersion relation of the atomic excitation and imaginary corrections of the Bloch quasi-momentum eigenstates of the non-Hermitian Hamiltonian. For multiply excited systems, quartic corrections to the HP expansion of the effective spin Hamiltonian for the atom chain dominate the coupling of the subradiant states and lead to a formulation equivalent to the Lieb-Liniger model of a 1D bosonic quantum gas in the Tonks-Girardeau regime [32,33]. The fermionic ansatz solution of that problem explains the decay rates and the properties of the solutions found in Refs. [8,9]. There is a high current interest and many potential applications of subradiance [22–25] and the analytical findings presented here may inspire further study of subradiance in light-matter interactions of more complex geometries, e.g., systems with higher dimensional atom arrays [35], chiral waveguides that break the parity symmetry [36] and setups with topological effects [37,38].

Let us conclude by discussing a remaining theoretical issue. We recall our effective separation of the Hamiltonian into an interaction term, V_{sub} , based on H_{eff}^I and an expansion on subradiant eigenmodes b_{ξ} for which H_{eff}^R contributes the decay rates γ_{ξ} . Like the numerical calculations, a more rigorous analytical approach should incorporate H_{eff}^R and H_{eff}^I on an equal footing. The fact that our separate treatment applies may be understood from the perturbation view. The leading order approximation of the

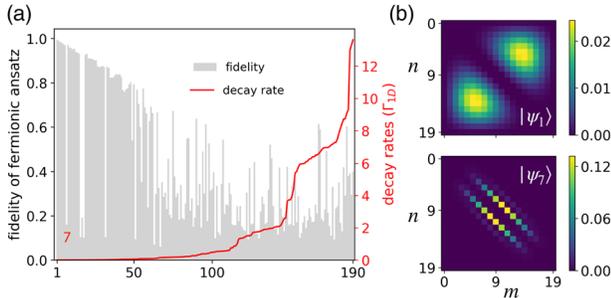


FIG. 2. (a) The two-excitation eigenstates of the system with $k_{1D} = 0.2\pi/d$ and $N = 20$ are sorted by increasing decay rates. The bars show the maximal fidelity that a fermionic ansatz can achieve for each eigenstate. The fermionic ansatz fits a broad range of the most subradiant states while a few exceptional states (the dips in the fidelity, e.g., state No. 7) show distinct non-fermionic behavior. (b) Position distributions of the atomic excitations, $|\langle \psi_{1(7)} | e_m, e_n \rangle|^2$ with $m(n) = 0, 1, \dots, 19$ of a typical fermionic subradiant state, $|\psi_1\rangle$ (upper panel), and the non-fermionic state $|\psi_7\rangle$ (lower panel). The lower panel feature at $|z_m - z_n| \approx 2d$ indicates that $|\psi_7\rangle$ supports dimerlike bound excitations.

subradiant states are the dark states of H_{eff}^I . They are also approximate eigenstates of H_{eff}^R when restricted to the most subradiant states. It means that the fermionic ansatz, as the leading order approximation, is shared by both H_{eff}^I and H_{eff}^R . Therefore, analyzing V_{sub} from the simpler H_{eff}^I is sufficient to capture the salient fermionic behavior. This is also verified by a direct construction of the fermionic ansatz without using the HP transformation, for both H_{eff}^I and H_{eff}^R [28]. Interestingly, we find that the fermionic ansatz does not exhaust all the most subradiant eigenstates. For a medium-size ensemble of $N = 20$ atoms, we obtained numerical eigenstates of H_{eff}^R with very low fermionic state fidelity. The subradiant states of this different character have well-defined “center of mass” wave number, and well-defined spatial separation, as illustrated in Fig. 2(b). Further discussion of these states is beyond the scope of this Letter, but may be of interest for future work possibly together with the interesting prospects for studying quantum fluctuations [39–41] in the Tonk-Girardeau gas theory by detection of the excited state correlations among atoms in a subradiant chain.

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