Multimode N00N states in driven atomtronic circuits

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We propose a method to generate multimode N00N states with arrays of ultracold atoms. Our protocol requires a strong relative offset among the wells and a drive of the interparticle interaction at a frequency resonant with the offsets. The proposal is demonstrated by a numerical and a Floquet analysis of the quantum dynamics of a ring-shaped atomtronics circuit made of M weakly coupled optical traps. We generate a hierarchy of energy scales down to very few low-energy states where N00N dynamics takes place, making multimode N00N states appear at nearly regular time intervals. The production of multimode N00N states can be probed by time-of-flight imaging. Such states may be used to build a multiple beam splitter.

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I. INTRODUCTION

Entanglement is at the root of quantum technology [1]. Depending on the number of subsystems involved, such genuine quantum correlations can be of very different nature [2,3]. While the picture for the pairwise case has been cleared up, multipartite entanglement remains challenging for both its mathematical analysis and quantum state engineering (see Refs. [4-6]). Nonetheless, multipartite entangled states were recognized as an important resource in different areas of quantum technology, including quantum information theory, quantum cryptography, and noteworthy quantum sensing [7–10]. As an example, nonclassical states of several identical particles possess intrinsic quantum correlations that can be exploited in technological applications. Among the most well-known nonclassical states, the bipartite "N00N" states correspond to the macroscopic superposition $\psi = (|N, 0\rangle +$ $|0, N\rangle)/\sqrt{2}$, where $|n_1, n_2\rangle$ have n_1, n_2 bosons in the modes 1 and 2, respectively. Such states are maximally entangled and highly sought for applications, e.g., to high-precision interferometry [11–13].

We focus on N00N states with multipartite entanglement [14]. The suggested platform is provided by cold atoms. They can be controlled with high precision and flexibility of the operating conditions [15]. In these systems, the physical con-

ditions, as for instance the atom-atom interaction, can be tuned in a wide range of values or as function of time [16-20]. Control of the atom number in small traps is achieved by fluorescence [21]. It is now possible to handle the atomic confinement on local spatial scales (micrometers) [22,23], and on timescales much shorter than atom's dynamics timescales [22,24–30]. Atomtronics exploits these progresses with the goal to widen the scope of cold atom simulators and to achieve new concepts of devices of practical value [31,32]. Roads to formation of bipartite N00N states with ultracold atoms were proposed either by phase imprinting followed by a free evolution [33-35], via rotation of the condensate [36], or using time-dependent fields [18,37–39]. Very few experiments with NOON states exist with ultracold atoms [40], and other nonclassical states have been reported such as squeezed [8] and oversqueezed states [41].

In this work, we propose a protocol for the generation of multimode N00N states, i.e. the states of the type $|N000...\rangle + |0N00...\rangle + |00N...\rangle + ...$ Such states are relevant to different contexts of quantum technology such as multiport interferometry [14]. Here we show how the interplay between periodic driving of the interaction and a strong offset can generate multimode N00N states with high fidelity at nearly periodic time intervals. As examples, we consider mesoscopic atom numbers (from N = 2 to 10) and number of modes M = 2-5.

Our atomtronic circuit is made of *N* neutral bosons trapped in *M* optical traps arranged in a ring geometry at zero temperature (Fig. 1). All the well's minima are offset by an energy $\pm \mu$ with respect to a reference well (the "source"), with $\mu \gg J_{ij}$ where the J_{ij} are the interwell tunneling amplitudes. The interaction is driven periodically [16–20,42,43] around 0 with frequency ω , chosen to be resonant with the offset, $\hbar \omega = 2\mu$. This makes both single-atom transitions

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FIG. 1. Scheme of the three-well geometry: bosons are confined in a triple-well ring configuration, where *J* is the tunneling rate, $\mu_1 = \mu$, $\mu_2 = 0$, and $\mu_3 = -\mu$ are respectively the local energy bias in sites 1,2, and 3. The blue arrows represent motion of bosons in pairs, triggered by the periodic drive. (Insets) M = 4-site and M = 5-site configurations with corresponding bias.

between offset wells and pair transitions from the source well resonant, with very different energy scales. A Floquet analysis shows that this generates very few low-energy states connecting the N00N state components. Optimal parameters for NOON states are found close to coherent destruction of tunnelling (CDT [16,18,44,45]) from the initial state, which we explain by a "Russian doll" fitting of five different energy (or time) scales, down to the N00N recurrence time. This turns out to be a powerful scenario to create multimode NOON states ($M \ge 3$) on timescales much shorter than those associated to the free evolution of the system (e.g., after a quench) [46-48]. Our analysis involves an exact solution of the quantum dynamics from the time-dependent Hamiltonian, as well as a high-frequency expansion of the Floquet Hamiltonian, that gives excellent agreement with the direct solution. Finally, we discuss how to detect experimentally such a correlated transport: second-order momentum correlations, as obtained after a time-of-flight expansion, can distinguish between coherent and localized states achieved in the transfer.

Section II presents the model and the analytical or numerical methods. Section III presents in detail the correlated boson transfer obtained from the driven dynamics. Section IV discusses the setup constraints and provides an explanation of multipartite N00N state formation. Section V explains the readout protocol. Section VI concludes, and appendices provide more details.

II. THE MODEL AND ITS ANALYSIS

The system is described by the M-mode Bose-Hubbard Hamiltonian subjected to a periodic modulation of the inter-

action strength:

$$H(t) = -\sum_{i,j=1 \ (i \neq j)}^{M} \left(J_{ij} \ a_{j}^{\dagger} \ a_{i} + \text{H.c.} \right) + \left(U_{0} + \frac{U_{1}}{2} \sin(\omega t) \right) \sum_{i=1}^{M} \hat{n}_{i} (\hat{n}_{i} - 1) - \sum_{i=1}^{M} \mu_{i} \ \hat{n}_{i}, \quad (1)$$

where a_i^{\dagger} creates a boson in site *i* and $\hat{n}_i = a_i^{\dagger} a_i$ is the number operator. The parameters J_{ij} quantify the hopping strengths, U_0 is the average interparticle interaction, U_1 the amplitude of its periodic modulation, and μ_i the local chemical potential, i.e., the well offset of the site *i*.

Let us first briefly mention the behavior of a three-atom ring with strong offset but static interaction U_0 only and $\mu_1 = \mu$, $\mu_3 = -\mu$. Details can be found in Appendix A. A second-order perturbative analysis in J/μ shows how to transfer a single pair with high fidelity from site 2 to sites (1,3) altogether, i.e., obtain a nearly perfect Rabi resonance between states $|n_1, n_2, n_3\rangle$ and $|n_1 + 1, n_2 - 2, n_3 + 1\rangle$. Such a highly correlated transition requires a fine tuning of the offset μ_2 . Importantly, this does not allow to transfer many pairs simultaneously, and the presence of a strong residual interaction that competes with the pair hopping prevents from creating a coherent state made of boson pairs.

Our aim is now to investigate how a correlated pair transport of ultracold bosonic atoms can be triggered with $U_0 = 0$ by a suitable choice of μ_i , U_1 , and ω . To illustrate the analysis in presence of a driven interaction, let us consider a three-well ring (see Fig. 1), with equal and real couplings $J_{ij} = J$, and offsets $\mu_{1,2,3} = \mu, 0, -\mu$. We drive the interaction strength as in Eq. (1) with $\omega \gg J/\hbar$ (see Refs. [49,50]). A key parameter for the dynamics is $K_0 = U_1/(\hbar\omega)$. Given the values of the relative offsets between the bosonic islands (1,2,3), one encounters two different Josephson frequencies for noninteracting atoms in a ring geometry: $\omega_I = \mu/\hbar$ for single boson transitions between 1 (or 3) and 2, and $\omega_I = 2\mu/\hbar$ for boson pairs hopping between 1 (or 3) and 2, as well as for single boson transitions between 1 and 3. Therefore coherent pair correlations are expected to be formed for resonant driving frequencies $\omega = 2\mu/\hbar$.

We study the dynamics through exact diagonalization of the many-body Hamiltonian (1) and the evolution of an initial state chosen to be $|0N0\rangle$. We complement this study by the analysis of an effective time-averaged Hamiltonian, following the method proposed by Dalibard and Goldman [51,52] (see also Ref. [53]). Our derivation (Appendix B) involves elimination of the second and third terms in Eq. (1) by a canonical transformation. The dressed $1 \leftrightarrow 3$ transitions can be directly averaged on time while the $1 \leftrightarrow 2$ and $3 \leftrightarrow 2$ transitions require a Floquet expansion of order $1/\omega$ yielding:

$$H_{\text{eff}} = -J \left[\mathcal{J}_1[K_0(\hat{n}_1 - \hat{n}_3 - 1)] a_1^{\dagger} a_3 + \text{H.c.} \right] + \frac{J^2}{\mu} \left[\mathcal{L}(K_0, \hat{n}_1, \hat{n}_2) (a_1^{\dagger} a_2)^2 + \mathcal{L}(K_0, \hat{n}_2, \hat{n}_3) (a_2^{\dagger} a_3)^2 + \mathcal{M}(K_0, \hat{n}_1, \hat{n}_2, \hat{n}_3) a_1^{\dagger} a_3^{\dagger} (a_2)^2 + \mathcal{N}(K_0, \hat{n}_1, \hat{n}_2, \hat{n}_3) a_1^{\dagger} a_3 + \text{H.c.} \right]$$

$$+ \frac{J^2}{\mu} \left[\hat{n}_2(\hat{n}_1 + 1) \mathcal{P}(K_0, \hat{n}_1, \hat{n}_2) - \hat{n}_1(\hat{n}_2 + 1) \mathcal{P}(K_0, \hat{n}_2, \hat{n}_1) + \hat{n}_2(\hat{n}_3 + 1) \mathcal{P}(K_0, \hat{n}_2, \hat{n}_3) - \hat{n}_3(\hat{n}_2 + 1) \mathcal{P}(K_0, \hat{n}_3, \hat{n}_2) \right]$$

$$(2)$$



FIG. 2. Generation of two-mode N00N states. (Top) Map of the N00N fidelity indicator $4F_{N0}F_{0N}$ with the interaction K_0 and reduced time. (Bottom) Time variation of the fidelities for a N00N state ($F_S = F_{N0} + F_{0N}$). N = 10 particles, $\mu = 15J$, and $K_0 = 2.81$.

 $\mathcal{L}, \mathcal{M}, \mathcal{N}, \mathcal{P}$ being infinite series of Bessel functions of the arguments $K_0(\hat{n}_i - \hat{n}_i)$ (see Appendix B).

III. CORRELATED ATOM PAIR TRANSPORT

The initial state involves *N* atoms in the unbiased trap 2. This choice aims at a redistribution in time of the N atoms into traps 1 and 3. Traps 2, and 1,3 can be respectively considered as the "source" and "output" modes of an atomic "beam splitter." The exact evolution of the initial state $\Psi(t = 0) = |0N0\rangle$ under the time-dependent Hamiltonian, or under the Floquet expansion Hamiltonian H_{eff} , is obtained by discretization of the Trotter formula. Convergence is checked with respect to the time step, much smaller than the smallest timescale in the Hamiltonian, i.e., the drive period.

A. The two-mode model

Let us benchmark our protocol with a driven two-mode model. In contrast with the models studied so far [18,34,35], driving of the interaction is combined with a large static offset. Starting in the state $|N, 0\rangle$ at time t = 0, Fig. 2 show a fidelity map for the N00N state as a function of K_0 and reduced time t (in units of \hbar/J). We choose as an indicator the product $4F_{N0}F_{0N}$, where $F_{N0} = |\langle N0|\Psi(t)\rangle|^2$, $F_{0N} = |\langle 0N|\Psi(t)\rangle|^2$. It shows that N00N states are created in sizable intervals of K_0 (Fig. 2), nearly with time periodicity. The recurrence time of N00N states is controlled by the interplay of small energy scales and is of the order of a few hundred times \hbar/J . The protocol operates for even values of N. Its efficiency comes from the atom pair dynamics triggered by the resonant drive, which considerably reduces the effective Hilbert space and helps atoms to bunch equally in both wells.





FIG. 3. Optimum transfer and three-mode N00N states. (Top) Comparison of the fidelity (maximized over $t \sim 2000 \hbar/J$) for transfer from $|0, 4, 0\rangle$ to $|4, 0, 0\rangle$, calculated from the exact dynamics (blue) and from the Floquet expansion (red). (Middle) Interaction-time map of the fidelity F_W . (Bottom) Time evolution of the fidelities F_W (in black) and for states $|N, 0, 0\rangle$, $|0, N, 0\rangle$, $|0, 0, N\rangle$, with N = 8 and $K_0 = 2.59$, $\mu = 17J$.

B. Three modes and tripartite entanglement

Remarkably, the two-mode result generalizes to larger well arrays, despite the growth of the Hilbert space. Let us consider the three-well symmetric case. Equation (2) shows that atom pair transfers $2 \leftrightarrow 1$ and $2 \leftrightarrow 3$ as well as $2 \leftrightarrow (1, 3)$ are dynamically generated.

Directed transfer of N atoms. Let us first consider full transfer from the "source" trap 2 to trap 1. Figure 3 shows the time dynamics of the transfer fidelity $F_{N00} = |\langle N00 | \Psi(t) \rangle|^2$ from state $|0, N, 0\rangle$ to state $|N, 0, 0\rangle$ as a function of the interaction, as well as its maximum on a large time interval (middle and top panels). The agreement between the exact dynamics and the one resulting from the effective Hamiltonian (3) is remarkable. For specific values of K_0 , the fidelity displays marked peaks. Such a targeted optimal transfer requires a fine tuning of the interaction parameter, especially for large (even) N. The system achieves a coherent transfer of N bosons from trap 2 to trap 1 (or to trap 3, at the same K_0 value but at different times). Moreover, for $K_0 \gg 1$, the asymptotic expansion of the Bessel functions gives $\mathcal{L} \sim \frac{\sin(2K_0)}{2K_0}$ (Appendix B). Therefore the transfer of pairs from the islands 2 to 1 or 3 is suppressed for $K_{0,n} \sim n\pi/2$, $n \in \mathbb{N}$. One verifies that this makes the system trapped in the initial state $|0, N, 0\rangle$, by a CDT mechanism [44], as visible on the fidelity map of Fig. 3. Three-mode generalized NOON states. We now demonstrate that, exploiting the ring geometry, coherent superpositions of the states $|N, 0, 0\rangle$, $|0, N, 0\rangle$, $|0, 0, N\rangle$ can be dynamically created for a wider range of system's parameters. To illustrate this, we represent the state of the system as $|\Psi_S(t)\rangle = |\Psi_W(t)\rangle + |\Psi_W(t)\rangle$ with

$$|\Psi_W(t)\rangle = a_1(t)|N, 0, 0\rangle + a_2(t)|0, N, 0\rangle + a_3(t)|0, 0, N\rangle$$
(3)

with $|\Psi_{W\perp}(t)\rangle$ an orthogonal vector to $|\Psi_W(t)\rangle$, and compute the fidelity on $|\Psi_S(t)\rangle$

$$F_W(t) = |a_1(t)|^2 + |a_2(t)|^2 + |a_3(t)|^2.$$
(4)

For all the explored values of N ($2 \le N \le 10$), we find that $F_W(t)$ maxima are nearly periodic and close to 1 (see Fig. 3, bottom panel and Appendix C). In particular, in specific intervals of K_0 , a three-component W-like superposition Ψ_W [4] with comparable amplitudes $|a_1| \simeq |a_2| \simeq |a_3|$ is generated with high fidelity at nearly regular times. Notice that forming a W-state $a_1|0N0\rangle + a_2|N00\rangle + a_3|00N\rangle$ is equivalent to a three-mode beam splitter operation.

One can check the effect of asymmetries (see Appendix C). First, taking the couplings J_{21} and J_{23} from the "source" well 2 as different, a similar efficiency is achieved for moderate asymmetry $\frac{|J_{21}-J_{23}|}{J_{21}+J_{23}}$. Second, varying the coupling J_{13} has a stronger effect. If $J_{13} = 0$, i.e., a linear configuration of the three wells, no three-mode N00N state is achieved, only two-mode N00N states are attained, either between wells 2,1 or between wells 2,3. Strikingly, the ring configuration turns out to be essential to create multimode N00N states.

C. More than three modes

Remarkably, our protocol can be generalized to engineer *M*-mode *W* states with M > 3. A four-mode and a five-mode *N*00*N* generator circuit are represented on Fig. 1. The latter is scalable to any odd *M*. Figure 4 shows fidelities up to 90% for M = 4 and 80% for M = 5. For *N* atoms in *M* modes, the size of the Hilbert space is $N(N, M) = \frac{(N+M-1)!}{N!(M-1)!}$, e.g., for instance, N(4, 3) = 15, N(10, 3) = 66, N(4, 5) = 70. Achieving N00N states with the latter numbers requires a precision of about 1% in the choice of parameter K_0 . This precision grows with N(N, M) but the average time (recurrence time) Δt separating the occurrence of N00N states does not increase. For the values of μ taken in this study, the recurrence time is of the order of a few hundreds times $\frac{\hbar}{J}$. This time scales linearly with the ratio $\frac{\mu}{J^2}$, as shown by the amplitude of the pair atom transfer in the Floquet expansion (3).

IV. DISCUSSION

A. Energy scale separation

The formation of multimode N00N states stems from the Floquet effective model, Eq. (2). It relies on energy scale separation. The resonant drive eliminates the largest scale μ from the averaged motion, leaving a dressed hopping of order J between wells 1,3, and pair hoppings as well as potential terms of order $J^2/\mu << J$. As a result, the spectrum of H_{eff} shows a few low-energy states with splittings $\sim J^2/\mu$, among



FIG. 4. Generation of *M*-mode N00N states. (Top) Four sites, $K_0 = 3.08$, $\mu = 20$, and N = 4. (Bottom) Five sites, $K_0 = 1.57$, $\mu = 15$, and N = 4. The total fidelity F_S generalizes Eq. (4) to *M* components.

high-energy states at scale $\sim J$. During time evolution from state $|0N0\rangle$ (N even) the wave function mostly stays within the low-energy sector made of paired states $|n_1, n_2, n_3\rangle$, i.e., with n_i even, while most Fock states other than $|N00\rangle$, $|00N\rangle$ participate in the high-energy one. This eventually makes the fully unbalanced states $|N00...\rangle$, $|0N0...\rangle$ linked together by a connecting chain of about $\frac{N}{2}$ states *only*. This drastically reduces the transition time between those states, compared to an undriven system. The process can be further optimized by noting that close to values yielding CDT, $K_0 \simeq n\frac{\pi}{2} + \delta K_0$, some of the transition amplitudes within the connecting chain are very small, of order $\delta J = (J^2/\mu) \delta K_0/K_0$. A perturbative analysis eventually makes a much lower energy scale emerge, that controls the transition between states $|N00...\rangle$, $|0N0...\rangle$,...To summarize, a kind of renormalization scheme produces a hierarchy of energy scales $\mu \gg J \gg J^2/\mu \gg$ $\delta J \gg (\delta J/J)^2 \mu \sim \hbar/\tau_{\rm rec}$, where $\tau_{\rm rec}$ is the typical NOON recurrence time (see Appendix D for a detailed analysis). The latter time is still experimentally accessible, as it is way shorter than the recurrence time predicted from a standard undriven dynamics of an interacting system.

B. The setup constraints

Let us now comment on the setup constraints. First, the multimode N00N states generated at times t_i are perfectly coherent. Yet, the relative phases of the components of each state change from a time t_i to another recurrence (see Appendix C). Notice that this is not detrimental to the multipartite entanglement of each state, and allows its use as a multiphase probe after release. Second, our protocol requires even filling of



FIG. 5. Read-out maps (a) $n(\mathbf{k})$ for a W state; (b) $\sigma(\mathbf{k}, -\mathbf{k})$ for $|0, 0, N\rangle$, (c) $\sigma(\mathbf{k}, -\mathbf{k})$ for $|0, N, 0\rangle$, (d) $\sigma(\mathbf{k}, -\mathbf{k})$ for the (nonideal) W-type state; $\sigma(\mathbf{k}, \mathbf{k})$ for a W-type state, either perfect (e) or with a component $|\Psi_{W\perp}(t)\rangle$ [Eq. (3)]. (f) N = 8.

the trap array, and the atom number to be constant during a time interval Δt , where $\Delta t \simeq |t_{i+1} - t_i|$. This requires an evaporation time $\tau_{ev} > \Delta t$.

The protocol assumes a drive resonant with the offsets. If this resonance is not perfect, i.e., $\omega = 2\mu/\hbar + \delta\omega$, the many-body interference making multimode N00N states is destroyed at long times. This can be estimated by comparing the "beating" frequency $\delta\omega$ to the recurrence time Δt . We indeed find numerically that N00N states still occur at times $t << \delta\omega^{-1}$, which requires the condition $\delta\omega << \Delta t^{-1}$ for the protocol to operate. This implies in practice that the offset μ should not be taken too large.

V. READOUT OF MULTIMODE NOON STATES

The diagnostic of the system relies on the time-of-flight technique: in our dynamical scheme, the state is prepared, driven and then released at the time $t = \bar{t}$. The long-time density in real space can be accessed by the Fourier transform at the releasing time \bar{t} [54,55]

$$\hat{n}(\mathbf{k}) = \sum_{i,j} e^{i\mathbf{k}\cdot(\mathbf{R}_i - \mathbf{R}_j)} a_i^{\dagger} a_j, \quad n(\mathbf{k},\bar{t}) = \langle \Psi(\bar{t}) | \hat{n}(\mathbf{k}) | \Psi(\bar{t}) \rangle.$$
(5)

While $n(\mathbf{k})$ is featureless [Fig. 5(a)], pair and many-boson pair transfers can be probed through the correlations in $\hat{n}(\mathbf{k})$:

$$\sigma(\mathbf{k}, \mathbf{k}', \bar{t}) = \langle \Psi_{\mathcal{S}}(\bar{t}) | \, \hat{n}(\mathbf{k}) \, \hat{n}(\mathbf{k}') \, | \Psi_{\mathcal{S}}(\bar{t}) \rangle - n(\mathbf{k}, \bar{t}) \, n(\mathbf{k}', \bar{t}).$$
(6)

A perfect $|0N0\rangle$ state [Fig. 5(c)] is characterized by a pattern in $\sigma(\mathbf{k}, \mathbf{k}')$ with $\mathbf{k} = -\mathbf{k}'$, due to coherent virtual

single-atom transitions towards one of the unoccupied sites (Appendix E). Rotated but similar patterns are obtained for states $|0, 0, N\rangle$ [Fig. 5(b)] and $|N, 0, 0\rangle$. The combination thereof gives rise to the pattern corresponding to the optimal *W* state: $\frac{1}{\sqrt{3}}(|0, N, 0\rangle + |N, 0, 0\rangle + |0, 0, N\rangle)$. The pattern for a dynamically achieved *W*-type state [see Fig. 5(d)] is indeed very similar to that of an optimal one (see Appendix E). Remarkably, it is also possible to probe the overall fidelity of the *W* state: indeed, the patterns taken instead at $\mathbf{k} = \mathbf{k}'$ are structureless for ideal *W* states. Therefore they allow probing other states contributing to $F_W < 1$ [Figs. 5(e) and 5(f)] (see Appendix E for details).

VI. CONCLUSIONS

We have considered a class of atomtronic circuits to study the correlated dynamics in bosonic networks: M coupled mesoscopic dots of interacting particles in-this is essentiala ring geometry. The correlated transfer results from the combination of suitable offset potentials and strong resonant driving of the interaction. The dynamics of such a system is dictated by short chains of states connecting the NOON components, making multimode N00N states form on realistic timescales. The diagnosis of the system states is carried out through the analysis of the momentum distribution which, in a cold atoms setting, corresponds to time of flight images. Our protocol continously produces multimode N00N states and is only limited by the lifetime of the optically trapped atoms. These states could be used to bring parallelism in multiple phase-imaging protocols [14]. Our study provides a new route to engineer complex correlations in quantum networks. The physical system we studied is within the current experimental capabilities in the atomtronics field.

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APPENDIX A: NONLOCAL PAIR TRANSPORT COMBINING BIAS AND STATIC REPULSION

Let us consider the effect of the strong offset μ with a *static* interaction only, with the Hamiltonian:

$$H_{=}\frac{U_{0}}{2}\sum_{j=1}^{3}n_{j}(n_{j}-1)-\sum_{j=1}^{3}\mu_{j}n_{j}-J\sum_{i,j=1(i\neq j)}^{3}(a_{j}^{\dagger}a_{i}+\text{H.c.}).$$
(A1)

This simple system indeed triggers correlated pair hopping between site 2 and sites (1,3) altogether. We choose for this purpose $\mu_1 = -\mu_3 = \mu$. Specifically, starting from the initial state $|\psi_i\rangle = |n_{1i}, n_{2i}, n_{3i}\rangle$, the correlated transport of *m* pairs brings the system coherently through the states $|\psi_f\rangle =$ $|n_{1i} + m, n_{2i} - 2m, n_{3i} + m\rangle$. N_T being the total number of pairs, an entangled target state is a superposition of such states:

$$|\psi_T\rangle = \sum_{m=1}^{N_T} c_m |n_{1i} + m, n_{2i} - 2m, n_{3i} + m\rangle.$$
 (A2)

The Hilbert space described by the states $|n_1, n_2, n_3\rangle$, can be split into subspaces with fixed difference $n_1 - n_3$, that contain the target state (A2). Correlated processes are favoured by enforcing energy separation of these subspaces, i.e., we set $\mu \gg J, U_0, \mu_2$. We use both exact diagonalization and second-order perturbation theory in J/μ to analyze how a pair of bosons is transferred with high fidelity within one subspace. Two-atom and higher-order virtual processes allow to identify an effective model for single pair transfer:

$$H_{\rm eff,hop} = J_{\rm eff}[(a_2^{\dagger})^2 a_1 a_3 + {\rm H.c.}],$$
 (A3)

Consider the transfer of a single pair from the state $|\psi_i\rangle = |n_1, n_2, n_3\rangle$ to the state $|\psi_f\rangle = |n_1 + 1, n_2 - 2, n_3 + 1\rangle$. Following a second-order effective hopping process, the intermediate virtual states are respectively $|\psi_{l_1}\rangle = |n_1 + 1, n_2 - 1, n_3\rangle$ and $|\psi_{l_2}\rangle = |n_1, n_2 - 1, n_3 + 1\rangle$. By mean of a quasidegenerate perturbation theory approach, the amplitude probability of the process is given by the sum over the contributions of all intermediate paths. We find that the effective hopping is given by

$$J_{\text{eff}}^{II} = -\frac{J^2 U_0 \left(\mu^2 + \mu_2^2 + 2U_0^2 + 3\mu_2 U_0\right)}{(-\mu + \mu_2 + U_0)(-\mu + \mu_2 + 2U_0)(\mu + \mu_2 + U_0)(\mu + \mu_2 + 2U_0)}.$$
 (A4)

For $\mu_2 = 0$ and small interaction, $J_{\text{eff}}^{II} \simeq -J^2 U_0/\mu^2$. More generally, the second-order hopping probability is zero for noninteracting particles $U_0 = 0$, due to a destructive interference between the virtual paths such as

$$\begin{aligned} |0, N, 0\rangle &\to |1, N - 1, 0\rangle &\to |1, N - 2, 1\rangle, \\ |0, N, 0\rangle &\to |0, N - 1, 1\rangle &\to |1, N - 2, 1\rangle. \end{aligned} (A5)$$

Optimization of the resonance between two states differing by one nonlocal atom pair happens when two levels of the spectrum encounter an anticrossing. In the vicinity of the anticrossing, these states are little different from states $|0, N, 0\rangle$ and $|1, N - 2, 1\rangle$. This generates a Rabi oscillation with nearly perfect fidelity on very long timescales (Fig. 6).



FIG. 6. (Top) Rabi-type oscillations of $F_{|1,8,1\rangle} = |\langle 1, 8, 1|e^{-iHt}|0, 10, 0\rangle|^2$ and $F_{|0,10,0\rangle} = |\langle 0, 10, 0|e^{-iHt}|0, 10, 0\rangle|^2$ (see text in Appendix A). Anticrossing makes a very long timescale emerge on top of fast single particle processes. Parameters are $\mu = 48$, $U_0 = 0.9$, $\mu_2 = 7.635$ (in *J* units). (Bottom) Dotted lines indicate the offset $\mu(N)$ required to achieve through tuning μ_2 a given fidelity transfer from $|0, N, 0\rangle$ to $|1, N - 2, 1\rangle$.

Notice that in this correlated transport regime, the interaction U_0 must be large compared to the transfer rates J_{eff}^{II} . As a result, coherent combinations involving a large number of states in Eq. (A2) cannot be achieved with the present static protocol. Single pair correlated transfer, though, is obtained with high fidelity by fine tuning the potential μ_2 to compensate the total interaction energy. It connects two states well-separated from the rest of the spectrum. High fidelity transfer can hold for any number of atoms $N = 2N_T$ for large enough μ , under the condition that the manifolds of states with the same $n_1 - n_3$ are well-separated from each other (in order that second-order processes in J/μ is well-defined). Large μ leads to a nearly perfect resonance (Fig. 6). Figure 6 also shows the corresponding scaling $\mu(N_T, F)$ of the minimum μ necessary to obtain a given fidelity F. Readout of the resulting superposition can be achieved by time-of-flight (see Appendix E).

APPENDIX B: DRIVEN INTERACTION, FLOQUET EXPANSION, AND THE EFFECTIVE MODEL

We consider a resonantly driven three-mode model ($\omega = 2\mu/\hbar$) with a strong offset between sites 1 and 3.

$$H = -J \sum_{i,j=1 \ (i \neq j)}^{3} \left(a_{j}^{\dagger} a_{i} + \text{H.c.} \right)$$

+ $\frac{U_{1}}{2} \sin(\omega t) \sum_{i=1}^{3} \hat{n}_{i} (\hat{n}_{i} - 1) - \sum_{i=1}^{3} \mu_{i} \hat{n}_{i} , \qquad (B1)$

We apply the analysis carried out by Goldman *et al.* [52] (see also Ref. [53]) to obtain an effective model in the strongly driven regime. Since $U_1 = K_0 \omega$ where $K_0 \sim 1$ and $\omega = 2\mu/\hbar$, two terms diverge in the limit of $\omega \rightarrow \infty$. They can be eliminated away by a unitary transformation. We first rewrite the Hamiltonian in a reference frame rotating with the driving term as

$$\tilde{H} = \mathcal{R}^{\dagger}(t)H(t)\mathcal{R}(t) - i\mathcal{R}^{\dagger}(t)\partial_t \mathcal{R}(t)$$
(B2)

with

$$\mathcal{R}(t) = -\exp\left(i\int_0^t O(\tau)d\tau\right),\tag{B3}$$

$$O(t) = \mu(\hat{n}_1 - \hat{n}_3) + \frac{U_1}{2}\sin(\omega t)\sum_j \hat{n}_j(\hat{n}_j - 1).$$
 (B4)

The Hamiltonian takes the form

$$\begin{split} \tilde{H}(t) &= -J \left(a_2^{\dagger} \exp \left[i K_0 \left(\hat{n}_2 - \hat{n}_1 \right) \sin \omega t - i \frac{\omega}{2} t \right] a_1 \\ &+ a_2^{\dagger} \exp \left[i K_0 \left(\hat{n}_2 - \hat{n}_3 \right) \sin \omega t + i \frac{\omega}{2} t \right] a_3 \\ &+ a_1^{\dagger} \exp \left[i K_0 \left(\hat{n}_1 - \hat{n}_3 \right) \sin \omega t + i \omega t \right] a_3 \right) + \text{H.c.} \\ &= \tilde{H}^{(21)}(t) + \tilde{H}^{(23)}(t) + \tilde{H}^{(13)}(t). \end{split}$$
(B5)

Following the high-frequency expansion, the effective Hamiltonian at order zero is given by

$$H_{\rm eff}^{(0)} = \frac{1}{T} \int_0^T \tilde{H}(t) dt.$$
 (B6)

The term $\tilde{H}^{(13)}(t)$ obviously has a nonzero time average on a period $T = \frac{2\pi}{\omega}$, yielding the following term corresponding to transitions between sites 1 and 3:

$$H_{\rm eff}^{(13)} = -J a_1^{\dagger} \mathcal{J}_1[K_0(\hat{n}_1 - \hat{n}_3)] a_3 + \text{H.c.}$$
(B7)

On the other hand, due to the half-frequency factor in the transitions from site 2 to sites 1 and 3, the corresponding terms have a zero average on the period 2T, which is the actual period of the transformed Hamiltonian \tilde{H} . Thus one needs to perform a Floquet expansion to first order in $1/\omega$, yielding terms of order J^2/μ :

$$H_{\rm eff}^{(21)} + H_{\rm eff}^{(23)} = \sum_{p>0} \frac{1}{p\mu} \left[\tilde{\mathcal{H}}_{12}^p + \tilde{\mathcal{H}}_{23}^p , \ \tilde{\mathcal{H}}_{12}^{-p} + \tilde{\mathcal{H}}_{23}^{-p} \right], \ (B8)$$

where \tilde{H}_{ii}^{p} denotes the *p*th harmonic of \tilde{H}_{ij} .

The calculation of the harmonics and of the commutators is straightforward and yields Eq. (2) with the definitions:

$$\mathcal{L}(K_{0},\hat{n},\hat{n}') = \sum_{m\geq 0} \frac{1}{(2m+1)} \\
\times (\mathcal{J}_{m+1}[K_{0}(\hat{n}-\hat{n}'-1)]\mathcal{J}_{-m}[K_{0}(\hat{n}-\hat{n}'-3)] - \mathcal{J}_{-m}[K_{0}(\hat{n}-\hat{n}'-1)]\mathcal{J}_{m+1}[K_{0}(\hat{n}-\hat{n}'-3)]), \\
\mathcal{M}(K_{0},\hat{n}_{i=1,2,3}) = \sum_{m\geq 0} \frac{1}{(2m+1)} \\
\times (\mathcal{J}_{m}[K_{0}(\hat{n}_{2}-\hat{n}_{1}-1)]\mathcal{J}_{-m}[K_{0}(\hat{n}_{2}-\hat{n}_{3}-2)] - \mathcal{J}_{m}[K_{0}(\hat{n}_{2}-\hat{n}_{1}-2)]\mathcal{J}_{-m}[K_{0}(\hat{n}_{2}-\hat{n}_{3}-1)] \\
+ \mathcal{J}_{m+1}[K_{0}(\hat{n}_{1}-\hat{n}_{2}-1)]\mathcal{J}_{-(m+1)}[K_{0}(\hat{n}_{3}-\hat{n}_{2}-2)] - \mathcal{J}_{m+1}[K_{0}(\hat{n}_{1}-\hat{n}_{2}-2)]\mathcal{J}_{-(m+1)}[K_{0}(\hat{n}_{3}-\hat{n}_{2}-2)]), \\
\mathcal{N}(K_{0},\hat{n}_{i=1,2,3}) = \sum_{m\geq 0} \frac{1}{(2m+1)} \\
\times [(\mathcal{J}_{m+1}[K_{0}(\hat{n}_{1}-\hat{n}_{2}-1)]\mathcal{J}_{-m}[K_{0}(\hat{n}_{2}-\hat{n}_{3})] - \mathcal{J}_{-m}[K_{0}(\hat{n}_{1}-\hat{n}_{2})]\mathcal{J}_{m+1}[K_{0}(\hat{n}_{2}-\hat{n}_{3}+1)])(\hat{n}_{2}+1) \\
+ \mathcal{J}_{-m}[K_{0}(\hat{n}_{1}-\hat{n}_{2}+1)]\mathcal{J}_{m+1}[K_{0}(\hat{n}_{2}-\hat{n}_{3})] - \mathcal{J}_{m+1}[K_{0}(\hat{n}_{1}-\hat{n}_{2})]\mathcal{J}_{-m}[K_{0}(\hat{n}_{2}-\hat{n}_{3}-1)]\hat{n}_{2})], \\
\mathcal{P}(K_{0},\hat{n},\hat{n}') = \sum_{m\geq 0} \frac{1}{(2m+1)} (\mathcal{J}_{m}[K_{0}(\hat{n}-\hat{n}'-1)]^{2} - \mathcal{J}_{m+1}[K_{0}(\hat{n}-\hat{n}'-1)]^{2}). \tag{B9}$$

This effective Hamiltonian can be further simplified in the limit of large interaction $K_0 \gg 1$, where the asymptotic expansion holds:

$$J_n(x) \sim \sqrt{\frac{2}{\pi x}} \cos\left(x - \frac{n\pi}{2} - \frac{\pi}{4}\right) \tag{B10}$$

yielding after a few steps the following form of the second-order terms in J/μ :

$$H_{\rm eff}^{\rm Asympt,(2)} = \left[\mathcal{H}^{(12)}(a_1^{\dagger} a_2)^2 + \mathcal{H}^{(23)}(a_2^{\dagger} a_3)^2 + \mathcal{H}^{(123)}(a_2^{\dagger})^2 a_1 a_3 \right] + \text{H.c.} + \mathcal{H}_{\rm pot}^{(123)}$$
(B11)

with

$$\begin{aligned} \mathcal{H}^{12} &= \frac{J^2}{2K_0\mu} [\delta_{\hat{h}_{12},2} \cos(2K_0) + (1 - \delta_{\hat{h}_{12},1})(1 - \delta_{\hat{h}_{12},2})(1 - \delta_{\hat{h}_{12},3}) \sin(2K_0)] \frac{1}{\sqrt{|\hat{h}_{12} - 1||\hat{h}_{12} - 3|}} \\ &+ \frac{J^2}{\mu} [\delta_{\hat{h}_{12},1} + \delta_{\hat{h}_{12},3}] \mathcal{J}_1(2K_0), \\ \mathcal{H}^{23} &= \frac{J^2}{2K_0\mu} [\delta_{\hat{h}_{23},2} \cos(2K_0) + (1 - \delta_{\hat{h}_{23},1})(1 - \delta_{\hat{h}_{23},2})(1 - \delta_{\hat{h}_{23},3}) \sin(2K_0)] \frac{1}{\sqrt{|\hat{h}_{23} - 1||\hat{h}_{23} - 3|}} \\ &+ \frac{J^2}{\mu} [\delta_{\hat{h}_{23},1} + \delta_{\hat{h}_{23},3}] \mathcal{J}_1(2K_0), \end{aligned}$$



FIG. 7. Trajectories and readout maps for (from top to bottom) N = 4, $K_0 = 2.4$, $\mu = 15$; N = 6, $K_0 = 2.7$, $\mu = 19$; and N = 10, $K_0 = 1.16$, $\mu = 11.7$, starting at t = 0 from state $|0N0\rangle$: (left) fidelities for states $|0N0\rangle$, $|N00\rangle$, $|00N\rangle$ and their sum F_W ; (right) two-atom correlation $\sigma(\mathbf{k}, -\mathbf{k})$ (see Sec. V and Appendix E) for a W-type state realized at times indicated by a vertical dotted line on the left panels.

$$\mathcal{H}^{123} = \frac{J^2}{2K_0\mu} \left[\frac{(1-\delta_{\hat{n}_{32},1})(1-\delta_{\hat{n}_{12},2})}{\sqrt{|\hat{n}_{12}-1||\hat{n}_{32}-2|}} \cos\left(K_0(\hat{n}_{13}-1)\right) - \frac{(1-\delta_{\hat{n}_{32},2})(1-\delta_{\hat{n}_{12},1})}{\sqrt{|\hat{n}_{12}-2||\hat{n}_{32}-1|}} \cos\left(K_0(\hat{n}_{13}+1)\right) \right] \\ + \left[\delta_{\hat{n}_{12},2}\mathcal{J}_0(K_0(\hat{n}_{32}-1)) + \delta_{\hat{n}_{32},1}\mathcal{J}_0(K_0(\hat{n}_{12}-2)) - \delta_{\hat{n}_{12},1}\mathcal{J}_0(K_0(\hat{n}_{32}-2)) - \delta_{\hat{n}_{32},2}\mathcal{J}_0(K_0(\hat{n}_{12}-1)) \right], \\ \mathcal{H}^{123}_{\text{pot}} = \frac{J^2}{2\mu} \left[\operatorname{sinc}(2K_0(\hat{n}_{21}-1)) \hat{n}_1 (\hat{n}_2+1) - \operatorname{sinc}(2K_0(\hat{n}_{21}+1)) \hat{n}_2 (\hat{n}_1+1) + \operatorname{sinc}(2K_0(\hat{n}_{32}-1)) \hat{n}_2 (\hat{n}_3+1) - \operatorname{sinc}(2K_0(\hat{n}_{32}+1)) \hat{n}_3 (\hat{n}_2+1) \right],$$
(B12)

where $\operatorname{sin}(x) = \frac{\sin x}{x}$. Inspection of the successive terms helps to understand several features of the exact numerical solution. First, the term \mathcal{H}^{123} vanishes when the wells 1,3 have the same occupation. Second, terms \mathcal{H}^{12} , \mathcal{H}^{23} vanish when $K_{0,n} = \frac{n\pi}{2}$, therefore coherent destruction of tunneling (CDT) is obtained starting from state $|0, N, 0\rangle$ with K_0 close to those $K_{0,n}$ values. Despite the approximation contained in this asymptotic expansion, it explains very well the map of Fig. 3, middle panel.

APPENDIX C: ADDITIONAL DATA WITH A DRIVEN INTERACTION

We present here some wave-function trajectories for other even atom numbers. For N = 4, 6, 10 one finds the same trends as in Fig. 3, bottom panel: tuning the interaction K_0 allows to periodically achieve a high-fidelity superposition of states $|0N0\rangle$, $|N00\rangle$, $|00N\rangle$, in particular *W*-type states (Fig. 7). Notice the nearly periodical appearance of such states, with a very long period (several $100J^{-1}$) related to anticrossings in the Floquet pseudoenergy spectrum. The larger *N*, the finer the tuning of K_0 necessary to obtain high fidelities.

We have also tested the effect of coupling asymmetries in the ring geometry (M = 3). Writing the intersite coupling as $H_J = \sum_{i \neq j} J_{ij} a_j^{\dagger} a_i$, three-mode N00N states can still be generated with moderate asymmetry $J_{12} \neq J_{32}$ (Fig. 8). On the other hand, keeping $J_{12} = J_{32}$ but taking $J_{13} \neq J_{12}$, N00N states also occur unless J_{13} is too small.

An important point is the coherence of the multimode N00N states. First, as resulting from the unitary evolution of a pure state $|0N0\rangle$, they are by construction pure states. We do not discuss here the possible causes of decoherence due to external fluctuations. Let us instead focus on the relative phases of the states forming such states, i.e., the phases of



FIG. 8. Trajectories for an asymmetric three-site ring (top) N = 6, $\mu = 17$, $K_0 = 2.65$, $J_{12} = 0.9$, $J_{23} = 1.1$, $J_{13} = 1$ and (bottom) N = 6, $\mu = 17$, $K_0 = 2.7$, $J_{12} = J_{13} = 1$, $J_{13} = 0.9$.

the coefficients $a_i(t)$ of a state:

$$\Psi_{\mathcal{S}}(t_k) = \sum_i a_i(t_k) |00\dots N_i\dots 00\rangle.$$
(C1)

We find that the a_i 's are in general complex numbers, and their phases, fixed for a given time, are not correlated from an occurrence time t_k to another t_l . This comes from the nontrivial interferences occurring in the many-body wave function. As an example, Fig. 9 gives the phases of successive N00N states components.

APPENDIX D: FORMATION OF THREE-MODE N00N STATES BY THE EMERGENCE OF VERY LOW ENERGY SCALES

Let us show how our protocol defines a hierarchy of energy scales, down to very low ones where N00N states form by resonance of very few low-energy Floquet states. The analysis is performed on the effective Hamiltonian, which provides a very accurate qualitative and almost quantitative description of the motion, averaged on the period \hbar/μ . Compared to the exact dynamics, the one led by H_{eff} misses the fast motion (on times \hbar/μ) and yields some irrelevant drift at longer scales, without changing the main conclusions. The scheme

	T = 228.5 t.J	T = 684 t.J	T = 1143 t.J	T = 1600 t.J
Phase $ 006\rangle$	1,007	2,018	-0,657	2,784
Phase $ 060\rangle$	-2,744	0,850	-2,334	-1,187
Phase $ 600\rangle$	0,023	0,320	-3,092	2,207

FIG. 9. Phases (in radians) of the coefficients of states $|006\rangle$, $|060\rangle$, $|600\rangle$ for different W states obtained successively in time.

will be illustrated on N = 2, 4, 6 examples with three wells. It generalizes to any even N.

The Floquet averaging leaves two kinds of terms in $H_{\rm eff}$. Direct averaging of the transitions between wells 1 and 3 yields a dressed single-atom hopping term of order $J \ll \mu$. This is, hierarchically speaking, the second energy scale in the problem. The third energy scale is $J^2/\mu \ll J$, it governs three kinds of terms, given by series of products of Bessel functions: (i) atom pair hoppings from 2 to 1, from 2 to 3, and from 2 to (1,3) simultaneously, as a nonlocal pair; (ii) corrections to the hopping between 1 and 3; (iii) potential terms. Notice that when the occupations n_1 and n_3 are equal, pair hopping from 2 to (1,3) is forbidden by a cancellation of terms in the Bessel series. This situation reminds that encountered in the undriven case with a constant repulsive interaction (Appendix A), where such transitions are found only if the repulsive interaction is nonzero. In the driven case, the average interaction is in fact zero.

One also notices that within the Fock space $|n_1, n_2, n_3\rangle$, H_{eff} splits into two uncoupled blocks, one made with even n_2 and the other made with odd n_2 . Starting with the initial state $|0N0\rangle$ with N even, the dynamics restricts to states $|n_1, n_2, n_3\rangle$ with n_2 even but n_1, n_3 can still be even or odd. This leaves 4 states over 6 for N = 2, 9 states over 15 for N = 4, 16 states over 28 for N = 6, more generally $(\frac{N}{2} + 1)^2$ states.

1. The limit of infinite μ

Let us treat H_{eff} by solving first the infinite- μ limit, then in perturbations in J/μ . Dropping all the J^2/μ terms yields $(1 + \frac{N}{2})$ blocks of states where n_2 is constant (and even), connected by 1 \leftrightarrow 3 hopping. Block n ($n = 0, \ldots, \frac{N}{2}$) is formed by (2n + 1) Fock states:

$$|2n, N-2n, 0\rangle, |2n-1, N-2n, 1\rangle \dots |0, N-2n, 2n\rangle.$$
 (D1)

This subspace is represented by a symmetric tridiagonal matrix with all diagonal terms being zero, and successive couplings $T_{n,1}, T_{n,2}, \ldots, T_{n,2n}$, which are of order *J* and obey $T_{n,m} = -T_{n,2n-m+1}$ (change of sign by interchanging n_1 and n_3), as can be understood by inspection of the Bessel functions in Eq. (B9). As a result, the matrix is simplified by taking the combinations:

$$|\psi_{n,m,\pm}\rangle = 2^{-1/2} (|2n - m, N - 2n, m\rangle)$$

 $\pm |m, N - 2n, 2n - m\rangle).$ (D2)

The dimension of this matrix is 2n + 1 and its spectrum consists of n pairs of states with energies $\pm E_{n,\kappa}^{(0)}$ ($\kappa = 1, ...n$) of order *J*, and one state at E = 0. To show this, one considers the two cases, *n* even and *n* odd and observe that this matrix further splits into two blocks. In the first one $|\Psi_{n,0,+}\rangle$ is coupled to $|\Psi_{n,1,-}\rangle$ and successively to $|\Psi_{n,n/2,+}\rangle$ if *n* is odd, and up to the "central" state $|n, N - 2n, n\rangle$ if *n* is even. The second point is that such tridiagonal matrices possess a zero eigenvalue if their dimension is odd. One checks that this is always the case with the matrix containing the state $|\Psi_{n,0,+}\rangle =$ $2^{-1/2}(|2n, N - 2n, 0\rangle + |0, N - 2n, 2n\rangle)$. For n = N/2 this is nothing but the NOON state $2^{-1/2}(|NOO\rangle + |OON\rangle)$ built on wells 1,3. Moreover, the zero-energy state of each block is obtained by successively eliminating all odd combinations, leaving only "even" states $|n - 2p, N - 2n, 2p\rangle$. Therefore, *all zero-energy states in the infinite* μ *limit are made exclusively with even occupations of all three sites* 1,2,3, spanning a reduced Hilbert space of dimension (N + 2)(N + 4)/8 instead of (N + 1)(N + 2)/2 in the full space. We show in the following that the essential of the dynamics takes place within this restricted boson pair Fock space.

The *n*th E = 0 state can be formally written as

$$|\psi_n^{(0)}\rangle = \alpha_0(|2n, N - 2n, 0) + |0, N - 2n, 2n\rangle) + \alpha_1(|2n - 2, N - 2n, 2) + |2, N - 2n, 2n - 2\rangle) + \cdots$$
(D3)

The α 's are coefficients of order 1. One has $|\psi_0^{(0)}\rangle = |0N0\rangle$, and $|\psi_{N/2}^{(0)}\rangle$ contains the (1,3) N00N state. Those remain to be connected to form the *W* state.

Gathering all blocks thus yields for the infinite- μ limit $(1 + \frac{N}{2})$ zero-energy states and $\frac{N}{4}(1 + \frac{N}{2})$ pairs of high-energy states with opposite energies of order *J*. The latter can be written with generality:

$$|\psi_{n,\kappa}^{(H)}\rangle = \sum_{m,\pm} \beta_{n,\kappa,m,\pm} |\Psi_{n,m,\pm}\rangle \tag{D4}$$

with coefficients of order 1.

2. Perturbation in J^2/μ

Let us now sketch the perturbative effect of all pair and potential terms. First, it raises the degeneracy of the $(1 + \frac{N}{2})$ zero-energy states, yielding $(1 + \frac{N}{2})$ states whose energies are of order J^2/μ . They are separated from the remaining states that stay at order J. Without details, one can write these low-energy states as

$$|\Psi_{\zeta}^{(L)}\rangle = \sum_{n=0}^{N/2} a_{\zeta,n} |\Psi_{n}^{(0)}\rangle$$
 (D5)

with coefficients of order 1 in general (except close to a CDT, see below), and energies ε_{ζ} of order J^2/μ . On the other hand, the high-energy states mix together to form states $|\Psi_{\kappa}^{(H)}\rangle$ with energies E_{κ} of order J.

The dynamics from the initial state $|0N0\rangle$ is obtained by expanding this state at t = 0 in the basis $|\Psi_{\zeta}^{(L)}\rangle$, $|\Psi_{\kappa}^{(H)}\rangle$, which yields at time *t*:

$$|\Psi(t)\rangle = \sum_{\zeta} a_{\zeta} e^{i\varepsilon_{\zeta}t} |\Psi_{\zeta}^{(L)}\rangle + \sum_{\kappa} b_{\kappa} e^{iE_{\kappa}t} |\Psi_{\kappa}^{(H)}\rangle.$$
(D6)

The first part of the sum with coefficients a_{ζ} of order 1 contains only even occupation states, while the second part contains all states, but with small coefficients $b_{\kappa,n,\pm}$ of order J/μ . Let us consider the transition probability to state $2^{-1/2}(|N00\rangle + |00N\rangle)$, mainly contained in the $|\Psi_{\zeta}^{(L)}\rangle$'s, with a small component in the $|\Psi_{\kappa}^{((H)}\rangle$. It is essentially given by

$$P_{N00+00N}(t) = \sum_{\zeta,\zeta'} c_{\zeta,\zeta'} \cos(\varepsilon_{\zeta} - \varepsilon_{\zeta'})t$$
 (D7)

with $c_{\zeta,\zeta'}$ of order 1, plus terms of order $(J/\mu)^2$, and it oscillates slowly. This formal calculation shows that the probability to form the state $2^{-1/2}(|N00\rangle + |00N\rangle)$ is large in the course

of time. Moreover, corrections will also weakly couple to state $2^{-1/2}(|N00\rangle - |00N\rangle)$ (participating to the high-energy states), thus explaining why in the trajectories, the fidelities for states $|N00\rangle$ and $|00N\rangle$ are quite similar but not identical (small corrections produce drift at very long timescales).

The conclusion of this perturbative analysis is that the interesting dynamics merely develops in a low-energy sector emerging from slightly perturbed degenerate states, with one in each block with constant n_2 . The number of these states is 1 + (N/2). Despite one works in a three-mode system, it is similar to the number of states of a two-mode system, with in addition a perfect pair correlation making a huge reduction of the chain of states connecting $|0N0\rangle$ to states $|N00\rangle$, $|00N\rangle$.

3. Proximity to CDT

The latter reasoning holds for any value of the interaction drive K_0 , but it does not yet explain why among the (N + 2)(N + 4)/8 pair states, excellent W states made of wellbalanced superpositions of $|0N0\rangle$, $|N00\rangle$, and $|00N\rangle$ can form for some values of K_0 . For values of N larger than 4, those can indeed be found in the vicinity of CDT regimes where the system is merely blocked in the initial state. The asymptotic expansion shows that this happens when $K_0 = n\pi/2$. CDT due to vanishing of Bessel function dressing has been found in the past for coupled wells [16], and Watanabe [18] noticed the proximity of CDT to good N00N states generated in such a simpler system. We hereafter give an explanation in our more general case.

The low-energy states connecting $|0N0\rangle$ to states $|N00\rangle$, $|00N\rangle$ are successively $|\Psi_1^{(L)}\rangle$, $|\Psi_2^{(L)}\rangle$, ..., $|\Psi_{N/2}^{(L)}\rangle$. Let us call their mutual couplings (of order J^2/μ) $t_1^{(L)}, t_2^{(L)}, \ldots, t_{N/2}^{(L)}$, that result from the amplitudes of the terms in H_{eff} transferring pairs one by one from, say, well 2 to well 1. The asymptotic form of these terms does not always vanish at CDT: it does so if $n_1 - n_2 < 1$ or if $n_1 - n_2 > 3$, numbers being taken in the final state of each transition [see Eq. (B12)]. As a result, t_i 's vanish except for $n_1 - n_2 = 2$, which happens once in the connecting chain if N/2 is odd. It it is even, a few other couplings to even state will be nonzero (see Fig. 10). In practice, working close to a CDT point and expanding the cardinal sine function with $K_0 = n\pi/2 + \delta K_0$, most pair couplings t_i 's are of order of the very low scale $t^{(\nu L)} = (J^2/\mu)(\delta K_0/K_0)$, and a few ones stay of order J^2/μ .

At this stage one can repeat the scheme already used in front of two different energy scales: diagonalize the matrix at CDT, followed by a perturbative expansion. At CDT one finds that states $|0N0\rangle$, $|N00\rangle$, $|00N\rangle$ are disconnected, as well as the states of the other first blocks n < N/4, which happens only for $N \ge 8$. Other states lie at energies of order J^2/μ due to their nonvanishing couplings at CDT. Then, deviating slightly from CDT, transitions from $|0N0\rangle$ to $|N00\rangle$ or to $|00N\rangle$ involve perturbative couplings trough those states, with amplitudes $t^{(vL)}$. This eventually generates an effective coupling between N00N components, of order $(t^{(vL)})^2/(J^2/\mu) \sim (J^2/\mu)(\delta K_0/K_0)^2$. The latter fixes the typical recurrence frequency $\omega_{\rm rec}$ of W states.

At these energy and time scales, we are left with the three N00N state components, plus a few even states sitting in the first blocks (zero for N = 2-6, two for N = 8, 10, ...).



FIG. 10. Fock states for (a) N = 2, (b) 4, and (c) 6. At infinite μ , rectangles delimit decoupled blocks, each containing one zeroenergy eigenstate of H_{eff} . Thin black arrows denote single-atom transitions [couplings $T_1 = O(J)$ with $T_1 = T'_1$ if $\mu = \infty$], double black arrows denote pair transitions [couplings $t_1 = O(J^2/\mu)$] and double red arrows pair transitions vanishing at CDT ($K_0 = n\pi/2$), having couplings $t_0 = O((J^2/\mu)(\delta K_0/K_0))$ (represented only for N = 4 and 6).

Eventually, the good fidelity of the three-mode N00N state at some precise times relies on canceling the probabilities of those "parasitic" states by interferences, which is quite easy to achieve when their number is small. To summarize, our hierarchical scheme—a kind of ad hoc renormalization—has successively eliminated: (i) odd n_2 states by approximating the exact dynamics by that of H_{eff} (at order J/μ); (ii) odd n_1, n_3 states at order J/μ ; and (iii) close to CDT, most pair states except $|0N0\rangle$, $|N00\rangle$, $|00N\rangle$ and very few others.

The hierarchy of energies thus generated is

$$\mu \gg J \gg J^{2}/\mu \gg t^{(vL)} = J^{2}/\mu \ (\delta K_{0}/K_{0})$$
$$\gg \omega_{\rm rec} = J^{2}/\mu \ (\delta K_{0}/K_{0})^{2}.$$
(D8)

4. Examples: N = 2, N = 4, N = 6

For N = 2, there are only two degenerate states at infinite μ : $|020\rangle$ and $2^{-1/2}(|200\rangle + |002\rangle)$. The only remaining even n_2 state is $|101\rangle$, which will happen only at order $(J/\mu)^2$. Therefore the dynamics starting from $|020\rangle$ reaches excellent W states of two bosons, whatever K_0 and at any time. Nearly balanced W states can be found at regular times. In this simple case there is no need for CDT. The reduction of the Hilbert space to a few atom pair states suffices, which shows the power of our scheme based on strong offset and resonant drive, making atom pair correlations.

For N = 4, the three infinite- μ degenerate states are $|040\rangle$, $2^{-1/2}(|220\rangle + |022\rangle)$ and $\alpha(|400\rangle + |004\rangle) + \beta|202\rangle)$ with α , β numbers of order 1. Figure 10 shows in red the couplings that vanish at CDT, and in black the small (double arrow, order J^2/μ) and the large (simple arrow, order J) couplings that are insensitive to CDT. Close to CDT, the resulting dynamics essentially involves states $|040\rangle$, $|400\rangle$, $|004\rangle$, the others being split at higher energy. Actually, the small size of the pair Hilbert space makes possible to find N00N states even far from CDT values. This shows that our protocol is very robust already to maximally entangle two or four particles in three modes, which is a nontrivial achievement.

The case N = 6 illustrates well the above discussion. Four blocks yield four degenerate states at infinite μ . Among them, the initial state $|060\rangle$ and the combination $2^{-1/2}(|600\rangle + |006\rangle)$ are specially favoured because they are very weakly coupled to other states close to CDT. This for instance generates N00N states on timescales $\sim 10^4 \hbar/J$ for $K_0 \sim 7.73$, close to $5\pi/2$.

Figure 11 shows the spectrum of H_{eff} for N = 2, 4, 6, first in a large K_0 range: one clearly sees a quasiperiodicity and a decreasing envelope characteristic of the Bessel functions



FIG. 11. Spectrum of the effective Hamiltonian for (a) N = 2, (b) 4, and (c) 6 at different scales, showing the low-energy states emerging in the center of the spectrum: two states for N = 2, three states for N = 4, four states for N = 6 of which two are only distinguished at very low scale. Zooming is made in the region of a CDT.



FIG. 12. Characterization of a resonant pair state with a static repulsive interaction. Parameters are $\mu = 20$, $\mu_2 = 5.781$, and $U_0 = 0.9$.

forming the various components of the Hamiltonian. Second, in a more restricted K_0 range containing a CDT, the lowenergy states are plotted. One more zoom is made in the N =6 case to show the very low energy states and the anticrossing at the CDT. Notice that there is no strong anomaly in the spectrum at an optimum K_0 value for N00N states ($K_0 = 7.73$ for N = 6).

APPENDIX E: READOUT MAPS, DETAILS AND EXPLOITATION

The two states $|\Psi_W\rangle$ and $|\Psi_{\text{pair}}\rangle$ that we have shown to be generated in our system are superpositions of states that are connected by at least one pair transfer. Therefore, the usual time-of-flight observable $n(k) = \langle \hat{n}(k) \rangle$ is flat for both states

and one has to analyze higher-order correlations to probe the presence of these states. We are interested in the following quantity : $\sigma_{\Psi}(\mathbf{k}, \mathbf{k}') = \langle \hat{n}(\mathbf{k}) \hat{n}(\mathbf{k}') \rangle - \langle \hat{n}(\mathbf{k}) \rangle \langle \hat{n}(\mathbf{k}') \rangle$, with

$$\hat{n}(\mathbf{k}) = \frac{|w(\mathbf{k})|^2}{N} \sum_{i,j} e^{i\mathbf{k}.(\mathbf{r}_i - \mathbf{r}_j)} b_i^{\dagger} b_j \quad , \quad w(\mathbf{k}) = e^{-\frac{x^2 |\mathbf{k}|^2}{4}} \quad (E1)$$

x defining a phenomenological broadening. The vectors \mathbf{r}_i defines the position of the sites representing the three wells. These sites are represented as the three summits of an equilateral triangle of length a = 1. We choose the origin of the coordinates as the center of this triangle.

1. With static repulsive interaction only

(See Appendix A) Let us now consider $|\Psi_{\text{pair}}\rangle = \alpha |0, N, 0\rangle + \beta |1, N - 2, 1\rangle$. The two states are connected by operators like $b_1^{\dagger} b_2 b_3^{\dagger} b_2$ or $b_2^{\dagger} b_1 b_2^{\dagger} b_3$ so that in the case $\mathbf{k}' = +\mathbf{k}$, the phase acquired is $e^{i\mathbf{k}.(\mathbf{x}_{21}+\mathbf{x}_{23})} = e^{i\sqrt{3}k_y}$. Finally:

$$\sigma_{\Psi_{\text{pair}}}(\mathbf{k}, +\mathbf{k}) = |w(\mathbf{k})|^4 \\ \left[\frac{2}{N} + \frac{4\sqrt{N(N-1)}}{N^2} |\alpha| |\beta| \cos\left(\sqrt{3}k_y + \phi_\alpha - \phi_\beta\right)\right]$$
(E2)

where ϕ_{α} , ϕ_{β} are the phases of coefficients α , β .

We can see that the resulting pattern is made of stripes modulated along k_y with a periodicity $\frac{2\pi}{\sqrt{3}}$ (see Fig. 12). Moreover, as $|0, N, 0\rangle$ and $|1, N - 2, 1\rangle$ are connected by only one pair transfer, $\sigma_{\Psi_{pair}}(\mathbf{k}, +\mathbf{k})$ also depends on the phase difference between the coefficient α and β . Therefore one can directly probe the coherence of this state by measuring $\sigma_{\Psi_{pair}}(\mathbf{k}, +\mathbf{k})$.



FIG. 13. Readout maps for ideal cases: fully localized (top panels), ideal *W* and *W*-type state, and N00N state (bottom). Red dots denote useful points for the determination of the weights of the *W*-state components (see text).



FIG. 14. Readout maps for a dynamically generated state, as a function of the filtering parameter κ (see text). (Left) $\sigma(k, -k)$ is slightly modified by the parasitic states. (Right) $\sigma(k, k)$ is a direct fingerprint of the parasitic states.

2. With a driven interaction: ideal cases

with

$$\mathfrak{Re}[\sigma_{|0N0\rangle}(\mathbf{k}, -\mathbf{k})] = \frac{|w(\mathbf{k})|^4}{N}$$
(E6)

$$|\beta|^2 (\cos(k_x + \sqrt{3}k_y) + \cos(k_x - \sqrt{3}k_y)),$$
 (E7)

$$\mathfrak{Re}[\sigma_{|N00\rangle}(\mathbf{k}, -\mathbf{k})] = \frac{|w(\mathbf{k})|^4}{N} |\alpha|^2 (\cos(k_x + \sqrt{3}k_y) + \cos(2k_x)), \quad (E8)$$

 $\mathfrak{Re}[\sigma_{|00N\rangle}(\mathbf{k},-\mathbf{k})]$

$$=\frac{|w(\mathbf{k})|^4}{N}|\gamma|^2(\cos{(k_x-\sqrt{3}k_y)}+\cos{(2k_x)}).$$
 (E9)

Therefore the probabilities in state *W* can be recovered using the following relations:

$$|\alpha|^{2} = -\frac{N}{2|w(\frac{\pi}{2}, -\frac{\sqrt{2}\pi}{2})|^{4}} \Re \mathfrak{e}[\sigma(\mathbf{k}, -\mathbf{k}))]_{\mathbf{k} = (\frac{\pi}{2}, -\frac{\sqrt{2}\pi}{2})}, \quad (E10)$$

$$|\beta|^2 = -\frac{N}{2|w(\pi,0)|^4} \mathfrak{Re}[\sigma(\mathbf{k},-\mathbf{k})]_{\mathbf{k}=(\pi,0)}, \qquad (E11)$$

$$|\gamma|^{2} = -\frac{N}{2|w(\frac{\pi}{2}, +\frac{\sqrt{2}\pi}{2})|^{4}} \Re \mathfrak{e}[\sigma(\mathbf{k}, -\mathbf{k})]_{\mathbf{k}=(\frac{\pi}{2}, \frac{\sqrt{2}\pi}{2})}, \quad (E12)$$

$$1 = -\frac{N}{2} \mathfrak{Re}[\sigma(\mathbf{k}, -\mathbf{k})]_{\mathbf{k}=(0,0)}.$$
 (E13)

$$|\Psi_W\rangle = \alpha |N00\rangle + \beta |0N0\rangle + \gamma |00N\rangle.$$
(E3)

One easily checks that the usual time-of-flight observable $n(\mathbf{k}) = \langle \hat{n}(\mathbf{k}) \rangle$ is flat. The same holds for different components of $|\Psi_W\rangle$ individually and for $|\Psi_W\rangle$ itself. For a "perfect" state $|\Psi_W\rangle$, there is only one contribution to $\sigma_{\Psi_W}(\mathbf{k}, \mathbf{k}')$ which leads to a **k** dependence. This contribution is due to virtual processes where a boson hops from a site *i* to a site *j* and comes back so that the phase acquired is $(\mathbf{k} - \mathbf{k}') \cdot \mathbf{x}_{ij}$. Looking at the case $\mathbf{k}' = -\mathbf{k}$, we find

$$\Re e(\sigma_{\Psi_{W}}(\mathbf{k}, -\mathbf{k})) = \frac{|w(\mathbf{k})|^{4}}{N} [(|\alpha|^{2} + |\beta|^{2}) \cos(k_{x} + \sqrt{3}k_{y}) + (|\alpha|^{2} + |\gamma|^{2}) \times \cos(2k_{x}) + (|\beta|^{2} + |\gamma|^{2}) \cos(k_{x} - \sqrt{3}k_{y})].$$
(E4)

As the states $|N00\rangle$, $|0N0\rangle$, and $|00N\rangle$ are not connected by second-order hopping, we can write

$$\sigma_{\Psi_{S}}(\mathbf{k},\mathbf{k}') = \sigma_{|N00\rangle}(\mathbf{k},\mathbf{k}') + \sigma_{|0N0\rangle}(\mathbf{k},\mathbf{k}') + \sigma_{|00N\rangle}(\mathbf{k},\mathbf{k}')$$
(E5)

As we can see in Fig. 13, for a symmetric setup the three Fock states that compose the superposition state give the same pattern but with a different orientation. The perfect *W* state with $\alpha = \beta = \gamma$ yields a pattern with an hexagonal symmetry and an asymmetric amplitude with respect to zero. A NOON state (for instance, $\alpha = \beta$, $\gamma = 0$) and an asymmetric superposition give a pattern similar to that of the perfect *W* state but distorted in different directions, depending on the coefficients α , β , and γ . These characteristics allow to probe the presence of a superposition of the three states $|N, 0, 0\rangle$, $|0, N, 0\rangle$, and $|0, 0, N\rangle$.

3. Checking purity

Here we show some results for N = 8 and for superposition states that are generated dynamically. Starting from the state $|0N0\rangle$, we let it evolve and stop at $T = 198t J/\hbar$, when the system is in a good superposition state. At this time, the fidelity for such a state is 0.9. This means that the wavefunction has also sizable components over a few other states. One finds that $\sigma_{\Psi_W}(\mathbf{k}, \mathbf{k})$ is flat for a perfect superposition state, this observable can indicate the contributions of the other parasitic states. To see this effect, one can delete all the states $|p\rangle$ in the wavefunction for which the coefficient $|Cp|^2 < \kappa$ and plot the quantities $\sigma_{\Psi_W}(\mathbf{k}, -\mathbf{k})$ and $\sigma_{\Psi_W}(\mathbf{k}, \mathbf{k})$ for different values of κ (Fig. 14). For $\kappa = 0.0$, all the states are present, for $\kappa = 0.1$ there is a small fraction of the parasitic states and for $\kappa = 0.2$ there are only the three states $|N00\rangle$, $|0N0\rangle$, and $|00N\rangle$.

The parasitic states can also alter the values of $\sigma_{\Psi_W}(\mathbf{k}, -\mathbf{k})$ so that it is more difficult to recover the composition of $|\Psi_W\rangle$ with the help of the analytical formula for the perfect state.

In the case $\kappa = 0.1$, there are only two parasitic states: $|N - 2, 0, 2\rangle$ and $|2, 0, N - 2\rangle$. The only contributions to $\sigma_{\Psi_W}(\mathbf{k}, \mathbf{k})$ are the terms which transfer 2 particles between 1 and 3, in the same direction, so that the result is an oscillation in the direction k_x (the period is π , as we can see in Fig. 14). The two quantities $\sigma_{\Psi_W}(\mathbf{k}, -\mathbf{k})$ and $\sigma_{\Psi_W}(\mathbf{k}, \mathbf{k})$ are thus complementary : the first serves to probe the superposition and the second to check its purity.

As an example, let us consider that $|\Psi\rangle = |\Psi_W\rangle + \epsilon(|N - 2, 0, 2\rangle + |2, 0, N - 2\rangle)$ like in the case $\kappa = 0.1$. At order ϵ , the only contribution is obtained when the operator $b_i^{\dagger}b_jb_l^{\dagger}b_m$ connects $|N - 2, 0, 2\rangle + |2, 0, N - 2\rangle$ with $|\Psi_W\rangle$. There are only two terms so that the result is $\sigma_{\Psi}(\mathbf{k}, \mathbf{k}') = \sigma_{\Psi_W}(\mathbf{k}, \mathbf{k}') + 2\epsilon \frac{\sqrt{2N(N-1)}}{N^2} (e^{i(\mathbf{k}+\mathbf{k}').\mathbf{x}_{13}} + e^{i(\mathbf{k}+\mathbf{k}').\mathbf{x}_{31}})$. So in the case $\mathbf{k} = \mathbf{k}'$, we can characterize the parasitic states:

$$\sigma_{\Psi}(\mathbf{k}, +\mathbf{k}) = \sigma_{\Psi_W}(\mathbf{k}, +\mathbf{k}) + 4\epsilon |w(\mathbf{k})|^4 \frac{\sqrt{2N(N-1)}}{N^2}$$
$$\times \cos(2k_x) + O(\epsilon^2).$$

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