

Compact unary coding for bosonic states as efficient as conventional binary encoding for fermionic states

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(Received 28 September 2021; revised 24 January 2022; accepted 21 March 2022; published 29 March 2022)

We introduce a unary coding of bosonic occupation states based on the famous balls and walls counting for the number of configurations of N indistinguishable particles on L distinguishable sites. Each state is represented by an integer with a human readable bit string that has a compositional structure allowing for the efficient application of operators that locally modify the number of bosons. By exploiting translational and inversion symmetries, we identify a speedup factor of order L over current methods when generating the basis states of bosonic lattice models. The unary coding is applied to a one-dimensional Bose-Hubbard Hamiltonian with up to $L = N = 20$, and the time needed to generate the ground-state block is reduced to a fraction of the diagonalization time. For the ground state symmetry resolved entanglement, we demonstrate that variational approaches restricting the local bosonic Hilbert space could result in an error that scales with system size.

DOI: [10.1103/PhysRevB.105.L121116](https://doi.org/10.1103/PhysRevB.105.L121116)

Finite lattice models represent an essential simplification in quantum many-body physics, where a set of discrete low-energy degrees of freedom are sufficient to capture the relevant phases and phase transitions of a more complete high-energy description. The states of these quantum systems form a finite dimensional Hilbert space \mathcal{H} and are thus amenable to a numerical representation through the exact diagonalization (ED) of a lattice Hamiltonian [1–14]. This provides a complete description of the eigenstates; however, it is limited by the exponentially increasing cardinality $|\mathcal{H}| \sim e^L$ of \mathcal{H} for L lattice sites. This problem has motivated stochastic and variational approaches, such as quantum Monte Carlo (QMC) [15–18] and the density matrix renormalization group (DMRG) [19–22]. These methods are essential to our understanding of quantum many-body phenomena [23–29], however, they are stochastic or require an approximate truncation of the local Hilbert space, and thus ED still plays a crucial role in benchmarking novel methods, as well as in gaining access to the density matrix. While extremely efficient approaches exist for the Heisenberg model [1] and its $SU(N)$ generalizations [30] with applications to fermionic lattice models of up to ≈ 40 sites, the expanded Hilbert space of bosonic systems, allowing multiple occupancy of a single spatial mode, presents a formidable challenge [9, 10, 31, 32].

In this Letter, we introduce a unary basis (UB) coding of bosonic occupation (Fock) states using a highly compact and compositional, yet still human-readable bit string exploiting the famous *balls and walls* illustration of particles and sites used in Bose-Einstein counting. For example, for $N = L = 11$:

$$|2, 0, 1, 0, 3, 0, 1, 0, 0, 0, 4\rangle \equiv |\bullet\bullet|\bullet|\bullet\bullet\bullet|\bullet|\bullet\bullet\bullet\bullet$$

$$I_{\text{UB}} = 2541296 \equiv 1001101100011011110000, \quad (1)$$

where $|n_1, n_2, \dots, n_L\rangle$ is a characteristic basis state with n_j the occupancy of site j . In the UB coding, 1's denote sites with the following number of 0's corresponding to that site's occupation, and each basis state can be represented as an integer in base 10. In addition to reducing the memory required to store the basis by a factor L , this approach provides rapid access to n_j and significantly accelerates the action of local operators via bitwise operations on I_{UB} . An implementation in the presence of lattice symmetries yields a reduction of computational complexity by a factor of L , reducing the time needed to generate the basis for a one-dimensional (1D) Bose-Hubbard (BH) model to a fraction of that needed to obtain the ground state via iteration. This yields a practical speedup of over $20\times$ compared to current methods based on unique integer ordering of permanents [9, 10] for system sizes up to $L = N = 20$.

The utility of the UB is validated by studying entanglement at the critical point of the 1D BH model for both spatial mode and particle bipartitions as well as in the presence of a $U(1)$ symmetry fixing the total number of particles. In the latter case, we show that the common practice of restricting the number of bosons per site to a small integer n_{max} (soft-spin approach) can lead to large relative errors that scale with the system size. This finding serves to challenge the conventional wisdom that an investigation of the n_{max} dependence of the energy alone is enough to justify this approximation. To facilitate adoption of the UB in future work, a complete software implementation is included in Ref. [33].

Background. For both fermionic and bosonic lattice models, it is natural to diagonalize any local N -particle Hamiltonian in a basis of spatial modes $|n_1, n_2, \dots, n_L\rangle$. For fermions, the Pauli exclusion principle restricts $n_j = 0, 1$ and each of the $|\mathcal{H}_f| = \binom{L}{N}$ basis states can be encoded as a binary word of length L that can be stored as a $2^{\lceil \log_2 L \rceil}$ -bit integer as

implemented in commonly used ED software [11–13,34]. For bosons, the possibility of multiple occupancy, $n_j = 0, \dots, N$ on any site enlarges the Hilbert space to $|\mathcal{H}| = \binom{N+L-1}{N}$ with basis vectors for the occupation states naturally parameterized by $L|\mathcal{H}|$ integers. The memory required to store the Hamiltonian has an upper bound $\sim |\mathcal{H}|^2$ and thus dwarfs that needed to store the basis states. However, in bosonic lattice models with limited range hopping, the Hamiltonian is sparse, reducing its storage cost to $\sim L|\mathcal{H}|$ and encoding the basis as arrays of site occupations now composes a leading share of the required memory. To address this problem, permanent ordering (PO) schemes [9,10,31,32] have been introduced which assign a unique contiguous integer label $I_{\text{PO}}(n_1, \dots, n_L)$ to each occupation state via an iterative procedure of complexity $O(L)$. To further avoid the memory impact of storing $\{I_{\text{PO}}\}$, a lookup table combined with an $O(L)$ inverse function I_{PO}^{-1} [10], can be implemented to gain access to site occupations numbers n_j . While this method has proven to be effective, e.g., in studying many-body localization [35,36] it has yet to be extended to exploit lattice symmetries (e.g., translation, inversion) where the maximum number of nonzero Hamiltonian matrix elements is reduced to $\sim |\mathcal{H}|$. Here it is crucially important to suppress the now leading order memory share ($\sim L|\mathcal{H}|$) of the basis.

Unary basis. To study the largest possible systems, we introduce an integer labeling scheme motivated by the formal equivalence between occupation states of N bosons on L sites, and those of N fermions on $N + L$ lattice sites [32]. By reinterpreting bits corresponding to the presence (1) or absence (0) of a fermion with a site boundary and boson, respectively, the resulting set of noncontiguous $|\mathcal{H}|$ integers $\{I_{\text{UB}}\}$ can encode information about the occupation state directly in their bit string [see Eq. (1)] while remaining as efficient as a binary encoding for fermions. This can be quantified by defining the efficiency of the representation as $\eta = \log_2 |\mathcal{H}| / \log_2 2^{L+N}$, providing a maximum efficiency at unit filling ($\nu = N/L = 1$, with N and $L \gg 1$) of $\eta = 1 - (1/4L) \log_2 L + O(L^{-1})$, where for $L = 20$, $\eta \simeq 0.925$. Comparing with an alternative approach where N bosons are mapped to a spin model with $S = N/2$ [12,13] that uses a sequence of $L \lceil \log_2(N + 1) \rceil$ bits coding $n_j \leq N$ on each site, the corresponding efficiency at unit filling is $\eta \approx 2 / \lceil \log_2(L) \rceil - 1 / (2L) + O((L \log_2 L)^{-1})$, which is smaller by a factor of $2 / \lceil \log_2(L) \rceil$ compared to the unary coding, and for $L = 20$, $\eta \approx 0.37$.

In addition to memory compactness, the UB coding can accelerate the inverse operation I_{UB}^{-1} to obtain access to occupation numbers. Given a 64-bit integer I_{UB} representing $|n_1, \dots, n_L\rangle$, n_1 can be found by counting the trailing zeros in the bit string of I_{UB} (a compiler builtin). Shifting the bits of I_{UB} by $n_1 + 1$ allows for reading the number of particles on the next site, and by repeating these bitwise operations L times, the corresponding occupation vector (OV) $|n_1, \dots, n_L\rangle$ can be constructed. This procedure can be further sped up by viewing the 64 bits of I_{UB} as a sequence of four 16-bit integers from which the corresponding number of sites and occupation numbers can be obtained by direct lookup of the 2^{16} possibilities then recomposed to generate the corresponding OV. We find that generating OVs for $\{I_{\text{UB}}\}$ is more than $4\times$ faster than obtaining them on the fly using PO, at a cost of additional memory usage corresponding to only $1/L$ of that needed to

store the system Hamiltonian in the absence of translational symmetry.

Lattice symmetries. The utility of the unary coding can be extended to treat systems that preserve Hamiltonian symmetries such as translation and inversion (see Supplemental Material [37]). If the Hamiltonian commutes with the translation operator \hat{T} , it has a block diagonal structure where each of the L blocks has a quasimomentum index q and contains a maximum number of nonzero elements that scales as $\sim |\mathcal{H}|$ (for short-range hopping). The resulting number of translationally symmetrized basis states (the q th degenerate set of the eigenstates of \hat{T}) scales as $\sim |\mathcal{H}|/L$. For a given q and OV, an eigenvector $|\phi_{\alpha,q}\rangle$ of \hat{T} can be generated: $|\phi_{\alpha,q}\rangle = L_\alpha^{-1/2} \sum_{j=0}^{L_\alpha-1} \exp[-2i\pi jq/L_\alpha] \hat{T}^j |n_1, n_2, \dots, n_L\rangle$, where α is the index of a cycle with length $L_\alpha \leq L$ [38] generated by repeatedly acting on a given $|n_1, n_2, \dots, n_L\rangle$ with \hat{T} . Each cycle can be mapped to a set of eigenvectors of \hat{T} with the same number of elements.

Calculating matrix elements of local operators requires access to all OVs in $|\phi_{q,\alpha}\rangle$, as well as the ability to reconstruct the state $|\phi_{q,\alpha}\rangle$ where a given OV appears. The naive storage cost for direct lookup of this information is at least double that needed for $\{I_{\text{UB}}\}$. Alternatively, one can trade memory with computational complexity by storing L_α for each cycle, as well as an extremal integer I_α that represents one OV characteristic of the cycle, (i.e., I_{PO} or its unary coding I_{UB}). Thus, when inverting I_α to an OV, the rest of the cycle can be obtained by L_α translations. If $I_\alpha \in \{I_{\text{UB}}\}$, the unary coding of the OVs of the cycle can be obtained by bit shifting I_α . As a single 64-bit instruction can shift a vector of eight bytes (using parallel intraprocessor bit paths) faster than moving each byte sequentially (using multiple memory operations), the UB coding provides a reliable constant speedup.

The inverse process of finding the cycle α where a given $|n_1, \dots, n_L\rangle$ appears can be done in four steps: (i) cyclic shifting L times, (ii) converting each of the L OVs to their integer representation, (iii) finding the characteristic integer, and (iv) performing a fast search for α in the ordered list of integers I_α . While steps (i) and (iii)–(iv) require at most $O(L)$ operations, the time complexity of step (ii) is $O(L^2)$, as it consists of repeating the operation of converting an OV to an integer L times. However, by exploiting the unary coding, the asymptotically slowest step (ii) is not required, as the shifting process (i) can be directly applied to I_{UB} to obtain the targeted list of integers. This represents a significant opportunity for a reduction in complexity and speedup over PO methods.

Benchmarking. As a test case, we consider the unary coding for N bosons on a ring of $L = N$ sites ($\nu = 1$) governed by the BH Hamiltonian,

$$\hat{H} = - \sum_{i=1}^L (b_{i+1}^\dagger b_i + \text{H.c.}) + \frac{U}{2} \sum_{i=1}^L n_i(n_i - 1), \quad (2)$$

where $n_i = b_i^\dagger b_i$, $[b_i, b_j^\dagger] = \delta_{i,j}$ and $b_{L+1} = b_1$. The on-site interaction $U > 0$ is measured in units of the hopping. Equation (2) exhibits a continuous phase transition from a superfluid ($U \ll 1$) to an insulator ($U \gg 1$) at $U \approx 3.3$ [23,39,40]. Using the unary coding of the basis, as well as translational ($q = 0, \dots, L - 1$) and inversion

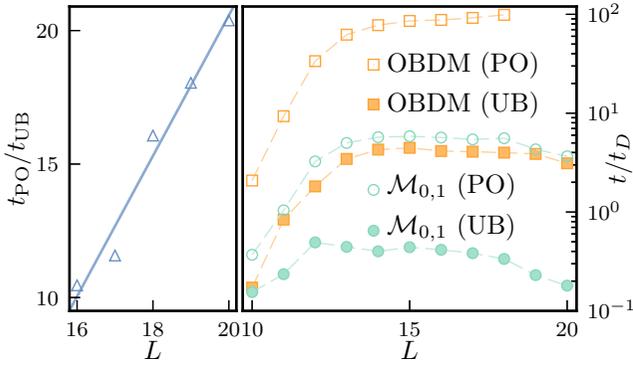


FIG. 1. Computational efficiency of the unary basis (UB) and permanent ordering (PO) encodings. Left: The ratio t_{PO}/t_{UB} of wall clock times needed to construct the ground-state block using the permanent ordering (PO) and unary basis (UB) codings as a function of system size L . The line is a linear fit demonstrating the complexity reduction discussed in the text. Right: The wall clock time ratio t/t_D versus system size L , where t_D is the time consumed in diagonalizing the ground-state block $\mathcal{M}_{q=0,r=1}$ of the Bose-Hubbard Hamiltonian and t represents the time spent in constructing $\mathcal{M}_{q=0,r=1}$ (circles) or calculating the elements of the one body density matrix (OBDM) in the ground state (squares).

($r = \pm 1$) symmetries on a ring, we have obtained the ground state $|\Psi_0\rangle$ of Eq. (2) via ED for up to $N = 20$ bosons on $L = 20$ sites.

To compare the efficiency of the PO and UB methods, we measure the times t_{PO} and t_{UB} spent in constructing the ground state block $\mathcal{M}_{q=0,r=1}$ of the sparse matrix representing \hat{H} . We find that the ratio t_{PO}/t_{UB} (Fig. 1 left), scales as $O(L)$, as expected from steps (i)–(iv) above. For $L = N = 20$, we find a speedup of $t_{PO}/t_{UB} \approx 20$. The practicality of this linear in L speedup becomes evident when measuring in units of t_D , the time required to iteratively obtain the ground state $|\Psi_0\rangle$ from $\mathcal{M}_{q=0,r=1}$. Again, for $L = N = 20$, we find $t_{BU}/t_D \approx 0.19$ compared to $t_{PO}/t_D \approx 3.96$, as illustrated in the right panel of Fig. 1. Thus, employing the unary coding can practically reduce computation time by a factor of $(t_{PO} + t_D)/(t_{UB} + t_D) \approx 4$. A larger computational speed up (≈ 20 for $L = N = 18$) is identified when calculating the elements of the one-body density matrix (OBDM) $\langle b_i^\dagger b_j \rangle$ (right panel of Fig. 1).

Application. As a further illustration of the utility of the unary coding allowing the use of an unrestricted local bosonic Hilbert space, we investigate several measures of entanglement in $|\Psi_0\rangle$. In all calculations, the required memory was less than 1.5 terabyte (TB).

By partitioning the spatial modes of the pure state $\rho = |\Psi_0\rangle\langle\Psi_0|$ into ℓ consecutive sites A_ℓ and its complement \bar{A}_ℓ , corresponding to the remaining $L - \ell$ modes, we obtain the reduced density matrix $\rho_{A_\ell} = \text{Tr}_{\bar{A}_\ell} \rho$ by tracing out all degrees of freedom in \bar{A}_ℓ . The von Neumann entanglement entropy $S(\ell) = -\text{Tr} \rho_{A_\ell} \ln \rho_{A_\ell}$ quantifying the amount of entanglement that exists between A_ℓ and \bar{A}_ℓ [41] is shown as square symbols in Fig. 2. The results demonstrate the reduction of spatial mode entanglement as the on-site repulsion U is increased and the system transitions from a superfluid to localized phase.

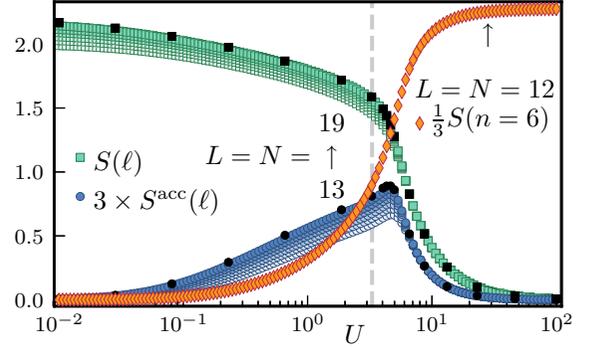


FIG. 2. Mode entanglement $S(\ell)$ and $S^{\text{acc}}(\ell)$ and particle entanglement entropy $S(n)$ in the ground state of the BH model as a function of interaction strength U at unit filling ($L = N$). For mode entanglement, the partition size is fixed at $\ell = \lceil L/2 \rceil$, where $f(x) = \lceil x \rceil$ is the least integer function.

With access to an exact representation of the ground state $|\Psi_0\rangle$, we can also study particle entanglement for a partition A_n of n particles [42–51]. This is most easily performed using the first quantized basis $|i_1, \dots, i_n\rangle$, where $i_k \in \{1, \dots, L\}$ is the lattice coordinate of the particle labeled $k \in \{1, \dots, n\}$. The familiar n -particle reduced density matrix of partition A_n is then

$$\rho_{A_n}^{\{i_k\}_a^n, \{j_k\}_a^n} = \sum_{\{i_k\}_{n+1}^N} \Psi^\dagger(\{i_k\}_1^n, \{i_k\}_{n+1}^N) \Psi(\{j_k\}_1^n, \{i_k\}_{n+1}^N),$$

where $\{i_k\}_a^b \equiv \{i_a, i_{a+1}, \dots, i_{b-1}, i_b\}$. The n -particle entanglement entropy, $S(n) = -\text{Tr} \rho_{A_n} \ln \rho_{A_n}$, is shown as diamonds in Fig. 2. In contrast to $S(\ell)$, $S(n)$ is large in the insulating phase, vanishes as $U \rightarrow 0$, and is extremely difficult to calculate via DMRG for $n > 2$ [52].

The interplay between fixed particle number and fluctuations between spatial modes, as reflected in symmetry-resolved entanglement, is a subject of growing interest [53–80] and can be potentially measured in ultracold lattice gases [81,82]. Here, the presence of conservation laws (e.g., fixed N), reduces the amount of entanglement that is accessible for quantum information processing which can be quantified using the accessible entanglement entropy $S^{\text{acc}}(\ell)$ (see Supplemental Material) [83–90]. Figure 2 demonstrates that the accessible entanglement entropy vanishes at the extremes of the two competing phases and peaks near the critical point.

Restricting the local Hilbert space by enforcing an upper limit n_{max} on the occupation number per site is a widely used approximation when performing the ED of a bosonic Hamiltonian or the DMRG representation of its ground state [23–25,39,91]. The validity of this approximation is judged based on the convergence of observables, (e.g., the ground state energy, the average occupation per site, or its fluctuations [9]) with n_{max} . While formally, the approximation is justified only in the insulating phase of the BH model, where the particle-number fluctuations are suppressed by repulsive on-site interactions, the resulting error in the ground-state energy remains small and finite, even in the superfluid phase down to the noninteracting limit [9]. This

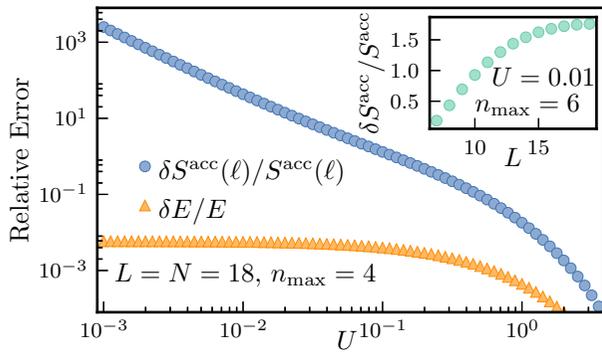


FIG. 3. The relative error in the accessible entanglement S^{acc} ($\ell = 9$) and the energy E of the BH ground state, arising from fixing the maximum occupancy at all sites to $n_{\text{max}} = 4$, as a function of interaction strength U/t at $L = N = 18$. Inset: Scaling of the relative error $\delta S^{\text{acc}}(\ell)/S^{\text{acc}}(\ell)$ with the system size L at $U = 0.01$ and $n_{\text{max}} = 6$, where $\ell = \lceil L/2 \rceil$.

empirical finding has yielded the conventional wisdom of utilizing a small n_{max} , even deep in the superfluid phase, in extended BH systems [92] and for fermions coupled to bosonic Hilbert spaces (e.g., the Hubbard-Holstein model [93]).

However, the truncated degrees of freedom *could* contribute significantly to other observables that crucially depend on particle-number fluctuations, specifically the spectrum of the n -particle reduced density matrix and entanglement in the presence of physical conservation laws. To illustrate, we calculate the error arising from restricting $n_{\text{max}} = 4$, in both the ground-state energy E and $S^{\text{acc}}(\ell)$. Figure 3 shows that while the relative error in E is less than 1%, the relative error in $S^{\text{acc}}(\ell)$ could be three orders of magnitude larger in the superfluid phase. It is expected that such large errors would also be present in the density-wave phase of the extended BH model at $U/t \ll 1$ [92]. Moreover, the relative error *increases* with L , (inset Fig. 3), even for $n_{\text{max}} = 6$, indicating that large-scale DMRG studies that employ n_{max} may contain logarithmically growing errors. To intuitively understand the origin of such scaling with L , we consider the simpler case of the bosonic

density reduced to half filling, where S^{acc} still vanishes as $U \rightarrow 0$. However, enforcing the extreme hard-core constraint of $n_{\text{max}} = 1$, which for an appropriate parity of N possesses the same $S^{\text{acc}}(\ell = L/2)$ for a partition of consecutive sites as noninteracting fermions. Here, it is known that $S^{\text{acc}} \sim \ln L$ and thus the error will have a similar scaling [78,90].

Discussion. In this Letter, we have introduced a compact unary coding for bosonic basis states that is both human and machine readable by enumerating configurations of indistinguishable particles on distinguishable sites through combinatorial counting of *balls and walls*. Having access to an efficient and compact exact representation of the ground state allows for the study of large system sizes (achieving equivalence for bosonic lattice Hamiltonians with state-of-the-art ED methods for fermions) and information quantities such as the particle partition or symmetry-resolved entanglement. These are not amenable to measurement via variational (due to the difficulty in computing n -particle density matrices or errors imposed by restricting the local Hilbert space) or stochastic (limited to Rényi entropies [94]) algorithms. Further constant speedups could be achieved using union data types that allow for the simultaneous view of an OV as an array of smaller integers. This could be useful when studying dynamics in bosonic systems, where access to asymptotic steady states requires considerable computational resources. While we have focused on a 1D BH Hamiltonian, the unary coding can be applied to any D -dimensional bosonic lattice model utilizing standard large-memory resources. By flattening a D -dimensional configuration, it can be represented by a single integer where translations can be achieved using bitwise operations on I_{UB} . Thus the reduction of time complexity in the presence of translational symmetry is expected to hold even for $D > 1$. We hope that the balls and walls coding will be rapidly incorporated into ED software, contributing to the acceleration of future studies of bosonic systems.

All code, scripts, and data used in this Letter are included in a GitHub repository [33].

Acknowledgments. This work was supported in part by the NSF under Grant No. DMR-2041995.

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