

Experimental verification of the area law of mutual information in quantum field theory

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Theoretical understanding of the scaling of entropies and the mutual information in quantum many-body systems has led to significant advances in the research of correlated states of matter, quantum field theory, and gravity. Although there have been several experimental measurements of Rényi and von Neumann entropies in systems with discrete degrees of freedom, the measurement of entropies between extensive intervals of continuous systems has remained elusive. In this work, we measure the von Neumann entropy of spatially extended subsystems in an ultra-cold atom simulator of one-dimensional quantum field theories. We experimentally verify one of the fundamental properties of equilibrium states of gapped quantum many-body systems, the area law of quantum mutual information. We also study the dependence of mutual information on temperature and the separation between the subsystems. Our work demonstrates the capability of ultra-cold atom simulators to measure entanglement in quantum field theories.

Some of the most commonly studied objects in quantum information are the (entanglement) entropy and the quantum mutual information. If a system described by the density matrix ϱ is composed of subsystems A and B , then the von Neumann (vN) entropy of subsystem A is defined as

$$S_A = -\text{Tr}(\varrho_A \ln(\varrho_A)), \quad (1)$$

where $\varrho_A = \text{Tr}_B(\varrho)$ is the reduced density matrix of subsystem A . If the state ϱ is pure ($\text{Tr}(\varrho^2) = 1$), then the vN entropy is a measure of entanglement between A and B , thus called entanglement entropy, where $S_A = S_B$. For mixed states ($\text{Tr}(\varrho^2) < 1$), the vN entropy captures both classical and quantum correlations, and it is no longer a good measure of entanglement. For mixed states, several other measures and witnesses of entanglement have been studied, with the positivity of partial transpose criterion being a prominent example [1]. In cases where, rather than entanglement, the shared amount of information between two subsystems A and B is of interest, the quantum mutual information (MI),

$$I(A : B) = S_A + S_B - S_{A \cup B}, \quad (2)$$

is a central object of study. It measures the total amount of correlation between the two subsystems, including all higher-order correlations for both pure and mixed states.

For pure states, the value of the MI is equal to twice the entanglement entropy of one of the subsystems.

Information theory measures reveal one of the fundamental properties of quantum many-body systems, the area laws. It was first noticed in gravitational physics, that surprisingly, the entropy of a black hole is proportional to the surface area of its event horizon and not its volume [2]. Interestingly, a similar property was found in quantum many-body systems: The vN entropy of ground states of systems with a gapped Hamiltonian scales with the surface area of the subsystem and not its volume [3]. Even more general, thermal states of systems with a gapped Hamiltonian exhibit an area law of mutual information [4]. This means that the information shared between parts of a quantum many-body system is only considerable over a short distance, set by the correlation length. In particular, such a bound on the required information to model a many-body system provides the foundations for the overwhelming success of tensor network-based methods [5]. In contrast, it is known that critical systems described by conformal field theory exhibit a logarithmic scaling instead of the area law [6]. In thermal states, while the mutual information has an area or a log law, the vN entropy will exhibit a volume law where it is proportional to the volume of the subsystem.

Extracting quantum information measures in quantum many-body systems has been the aim of several experi-

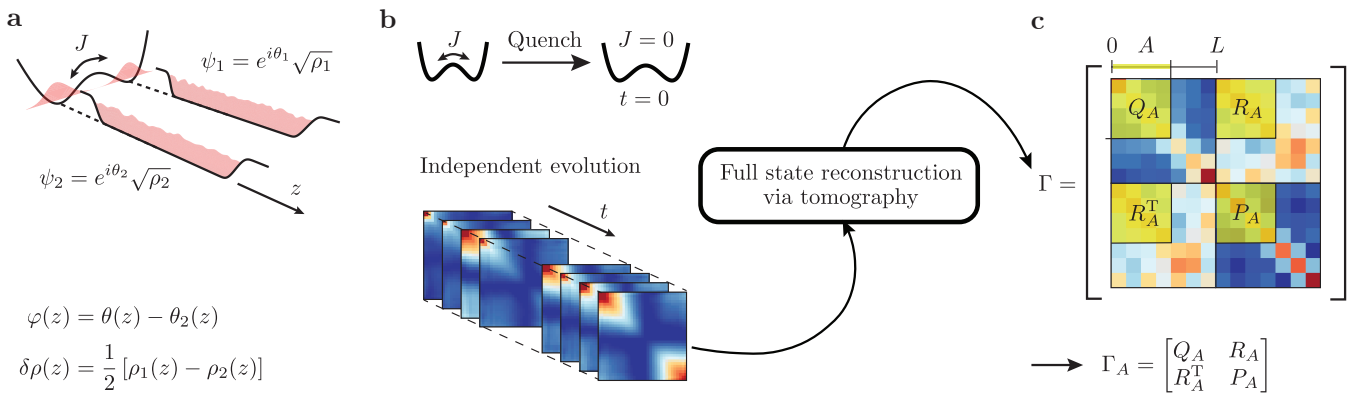


FIG. 1. **Schematic of the experimental protocol.** **a**, The experimental protocol starts by cooling down a pair of tunnelling coupled superfluids in a double-well potential with a finite single particle tunnelling rate J , typically $\approx 2\pi \times 1$ Hz. The radial trapping frequency is $\omega_{\perp} = 2\pi \times 1.4$ kHz and typical linear densities are $n_{1D} \approx 70 \mu\text{m}^{-1}$. **b**, Right after the cooling process, the tunnelling rate is changed to zero in 2 ms. At $t = 0$, the two condensates are already uncoupled and their independent evolution following the Tomonaga-Luttinger liquid Hamiltonian starts. The measured phase correlations at different times are used to fit the full covariance matrix Γ . **c**, To calculate the vN Entropy for a subsystem, $S(\Gamma_A)$, we use the correlations in that specific region (yellow shaded area).

ments. Calculating the vN entropy requires access to the density matrix of the full system, ρ , which is usually a tedious task in experimental platforms. Due to the difficulty of measuring the vN entropy, a commonly measured quantity is the second-order Rényi entropy [7, 8]. Nevertheless, several optical lattice setups are able to measure the vN entropy, which is obtained either from a single site reduced density matrix or as a classical thermodynamic entropy of the whole system [9, 10]. In special cases, when the system is diagonal enough, even many body vN entropy can be accessed [11]. However, the measurement of vN entropy between extended spatial subsystems has so far remained elusive, as well as the verification of the predicted area law scaling of the MI.

In this work, we address these challenges and study the scaling of the vN entropy and the MI with subsystem size in a continuous quantum many-body system. Our setup is a pair of tunnelling-coupled quasi-one-dimensional (1D) ultracold Bose gases (see Fig. 1a), cooled down and trapped below an atom chip [12]. Along the longitudinal axis, z , the clouds are confined in box-like potentials with hard walls, created by superposing magnetic and optical dipole potentials [13]. In one of the transverse directions, the atoms are trapped in a double-well potential, created by dressing with radiofrequency fields. The single-particle tunnelling rate J between the two condensates is adjusted by changing the amplitude of the radiofrequency fields [14].

To prepare a state in thermal equilibrium, we directly cool down an atomic cloud of ^{87}Rb in a strongly coupled double-well, using standard techniques of laser cooling and evaporative cooling [15]. The quantum fields describing each condensate can be written in phase-density representation as $\psi_n(z) = \exp(i\theta_n)\sqrt{\rho_n}$, with $n = 1, 2$. The spatially resolved relative phase between the two condensates, $\varphi(z) = \theta_1(z) - \theta_2(z)$, is extracted from in-

terference images, taken 15.6 ms after releasing the atoms from the trap and letting them fall freely. In the limit of low energy excitations, the correlations of the relative phase are well described by the sine-Gordon Hamiltonian [16]. Expanding the interaction term in the case of high tunnelling rates leads to the massive Klein-Gordon Hamiltonian,

$$H_{\text{KG}} = \int_0^L dz \left[g_{1D} \delta\rho^2(z) + \frac{\hbar^2 n_{1D}}{4m} (\partial_z \varphi(z))^2 + \hbar J n_{1D} \varphi^2(z) \right], \quad (3)$$

where L is the length of a system with uniform averaged 1D density n_{1D} , g_{1D} is the 1D interaction strength, m is the mass of an atom, and the relative density, $\delta\rho(z) = [\rho_1(z) - \rho_2(z)]/2$, is the conjugate field of the relative phase, fulfilling $[\varphi(z), \delta\rho(z')] = i\delta(z - z')$. A direct measurement of $\delta\rho(z)$ is unfeasible in our current experimental setup. Hence, to reconstruct the full covariance matrix of the initial state,

$$\Gamma = \begin{bmatrix} Q & R \\ R^T & P \end{bmatrix}, \quad (4)$$

we use a tomography procedure [17]. Here, $Q_{ij} = \langle \varphi(z_i) \varphi(z_j) \rangle$, $P_{ij} = \langle \delta\rho(z_i) \delta\rho(z_j) \rangle$, $R_{ij} = \langle \frac{1}{2} \{ \varphi(z_i), \delta\rho(z_j) \} \rangle$ with $i, j \in \{1, \dots, N\}$, where z_i denotes different points on a discrete grid with N points. The upper limit for N is given by the resolution of the imaging system, which limits our access to higher momentum modes (larger than N) and enforces an ultraviolet (UV) cutoff.

To apply this tomographic method, we ramp up the barrier between the two strongly coupled condensates in 2 ms, and let them evolve independently in the uncou-

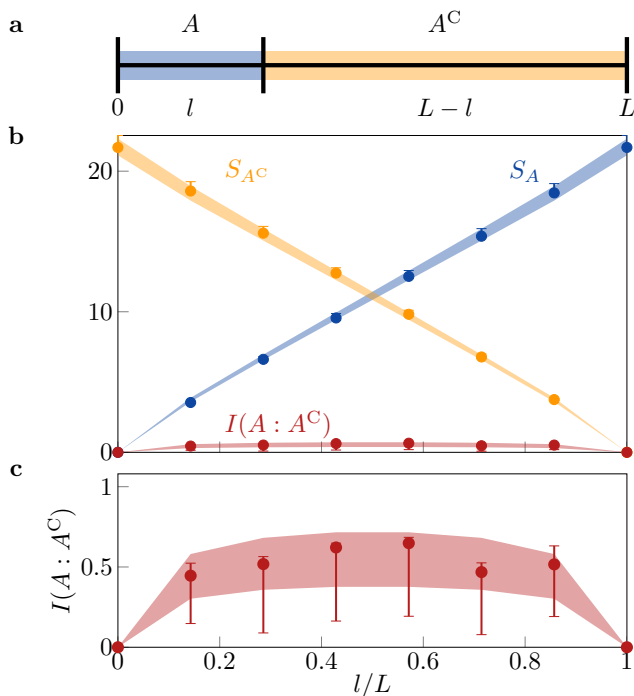


FIG. 2. **Area law of MI and volume law of vN entropy.** **a**, A system of size $L = 49 \mu\text{m}$ is divided into a subsystem A of length l and its complement subsystem of length $(L - l)$. **b**, The experimental results for $I(A : A^C)$, S_A , and S_{A^C} , calculated based on $N = 7$ modes, are plotted as bullets with error bars representing the 95% confidence intervals obtained via bootstrapping [18]. The shaded areas show the 95% confidence interval for the theory predictions, considering the uncertainty in the estimated temperature and tunnelling rate J (see SM). **c**, Close up of the measured MI.

pled double-well ($J = 0$). The post-quench dynamics follow the Tomonaga-Luttinger liquid Hamiltonian. Over time, the initial eigenmodes of the relative density rotate into the phase quadrature, enabling us to access the information about these eigenmodes by fitting the initial second-order correlation functions of phase-density and density-density to the measured evolution of the phase-phase correlations. Starting with a thermal equilibrium state of a quadratic Hamiltonian and quenching to another quadratic Hamiltonian ensures that the initial state is Gaussian and remains Gaussian under evolution [19]. To confirm this, we measure the normalized averaged connected fourth-order correlation function, $M^{(4)}$.

Having Gaussian states significantly simplifies the calculation of the quantum information measures - an otherwise highly non-trivial task in quantum field theory [20]. Gaussian states are fully described by their covariance matrix Γ . Its symplectic spectrum is obtained by diagonalizing $i\mathcal{J}\Gamma$ where $\mathcal{J} = \begin{bmatrix} 0 & I \\ -I & 0 \end{bmatrix}$ is the symplectic unit. The symplectic spectrum consists of pairs $\pm\gamma_n$, $n \in \{1, \dots, N\}$. It encodes the complete information that is contained in the covariance matrix. Consequently, it

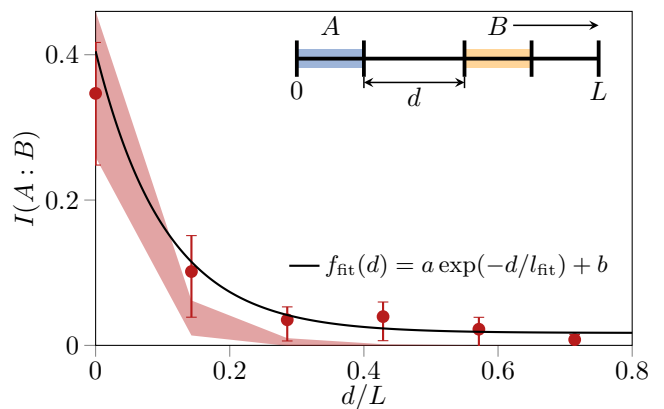


FIG. 3. **Shared information content between two spatially separated subsystems.** Here, we calculate the MI of two disjoint subsystems with the same length, $l/L = 0.15$, as a function of the distance between them, d . As demonstrated in the inset, while the subsystem A is kept fixed on the left edge of the system, B is shifted away to the other edge. Bullets and shadings represent experimental and theoretical data for $I(A : B)$ respectively (see Fig. 2 for more details on the error bars and shaded area). The solid black line is an exponential fit with $l_{\text{fit}} = 5.1 \mu\text{m}$. The finite bias b results from estimating a positive quantity, MI, using finite statistics.

can be used to reconstruct the full density matrix of the state and the measures of quantum information. In particular, the vN entropy is given by

$$S(\Gamma) = \sum_{n=1}^N \left[\left(\gamma_n + \frac{1}{2} \right) \ln \left(\gamma_n + \frac{1}{2} \right) - \left(\gamma_n - \frac{1}{2} \right) \ln \left(\gamma_n - \frac{1}{2} \right) \right] \quad (5)$$

For non-Gaussian states, neglecting higher-order correlations and estimating the vN entropy based on the covariance matrix gives a lower bound to the actual entropy [21].

Having the reconstructed initial covariance matrix at hand, we use expression (5) to calculate the vN entropy of any subsystem A , $S_A(\Gamma_A)$ (see Fig. 1c). Using equation (2), the MI between two subsystems A and B is calculated. To observe the scaling of the vN entropy and the MI, we calculate the vN entropy of subsystems with different lengths and consecutively the MI with their complement subsystems as illustrated in Fig. 2a. As expected for thermal states, the vN entropy is in the volume law regime, depending linearly on the size of the subsystem (Fig. 2b). Measuring the vN entropy allows us to study the scaling of the mutual information. We find an area law for MI, with a plateau forming in the bulk of the system (Fig. 2c). Our results represent an experimental verification of one of the elementary features of quantum many-body systems [4].

We continue by studying the dependence of MI on the distance between two subsystems. In this case, we calculate the MI of two subsystems A and B , separated by a gap of length d . The results are presented in Fig. 3.

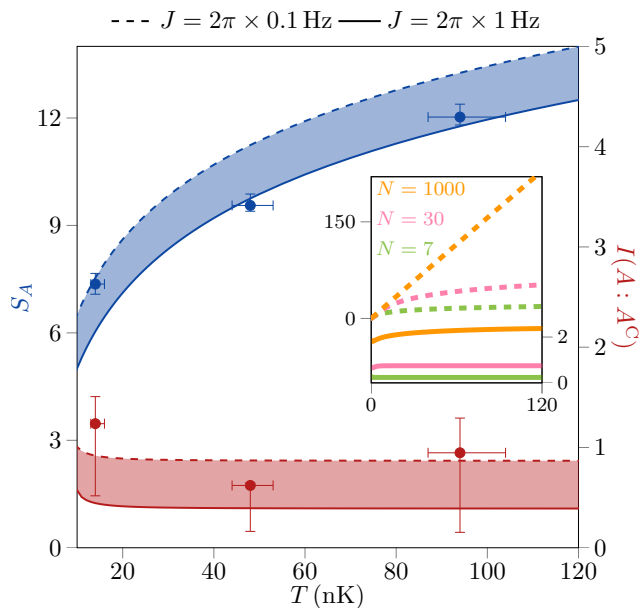


FIG. 4. **Temperature dependence of MI and vN entropy.** The bullets represent experimental data for S_A (left axis) and $I(A : A^C)$ (right axis) with $N = 7$ for measurements with three different temperatures, but otherwise similar parameters. The size of the subsystem A is $l/L = 0.4$ (see Fig. 2a). The error bars show the 95% confidence intervals achieved via bootstrapping. The shaded areas represent the calculations based on the theoretical model for different tunnelling rates, between $J = 2\pi \times 0.1$ Hz (dashed lines) and $2\pi \times 1$ Hz (solid lines). The inset shows the theoretical predictions for S_A (left axis) and $I(A : A^C)$ (right axis) for three different values of N , for $J = 2\pi \times 0.4$ Hz without considering the effect of finite optical resolution (see SM).

As expected, the MI decreases as the two subsystems get further apart. We can extract a decay length by fitting an exponential function to the experimental data. The fitted decay length, $l_{\text{fit}} = 5.1$ (3.7, 8) μm agrees with the correlation length $l_C = 6.8$ (6, 7.7) μm calculated based on the experimental parameters (see SM). The intervals represent the 95% confidence intervals obtained via bootstrapping.

For typical temperatures of our experiment, the vN entropy depends linearly on the temperature. The linear dependence changes to a logarithmic dependence by introducing a finite UV cut-off, as shown in the inset of Fig. 4 (dashed lines). As temperature increases, the symplectic eigenvalues grow, and the calculated entropy using a finite number of modes saturates. The MI, however, regardless of the UV cut-off, reaches a finite asymptotic value, given by the classical correlations [22]. Reducing the number of modes reduces the asymptotic value due to the limited available information in the modes taken into account, as presented in the inset of Fig. 4 (solid lines). In Fig. 4, measured vN entropy and MI for three different measurements are illustrated along with the theoretical predictions. The measurements agree with the theory

calculated for the extracted parameters and $N = 7$ lowest modes.

It is important to stress that the only assumption we make to calculate the vN entropy and the MI is that the post-quench dynamics follow a Tomonaga-Luttinger liquid Hamiltonian, which has been confirmed in the previous work [19]. Our measurements do not rely on any assumption related to the Gaussianity of the initial state. We rather confirm that the initial state is Gaussian by measuring the higher-order correlations. Even for non-Gaussian initial states, our results would represent a lower bound to the entropy of the full state.

The results presented here are a step towards the more ambitious goal of measuring many-body entanglement in a continuous 1D quantum system. There are two main limiting factors preventing us from detecting entanglement. The entanglement can be detected if a sufficient number of momentum modes are measured whose mode occupation numbers are close to the ground state. In our current experiments, the nonzero temperature (10-100 nK) of the Bose gas keeps the occupation numbers of the lower momentum modes too high above the value of the ground state. At the same time, the finite optical resolution introduces a soft cut-off which not only prevents us from measuring higher momentum modes, but also modifies the lower modes that can be measured. Improving any of these two aspects would make the detection of the entanglement possible.

Another promising direction for future work is to go beyond quadratic models and detect entanglement in an interacting model. It has already been demonstrated that the atom chip experiments can successfully simulate the sine-Gordon model, and higher-order correlation functions can be measured [16]. Developing a tomography procedure for this setting would give us access to entanglement properties in interacting many-body quantum systems [23, 24].

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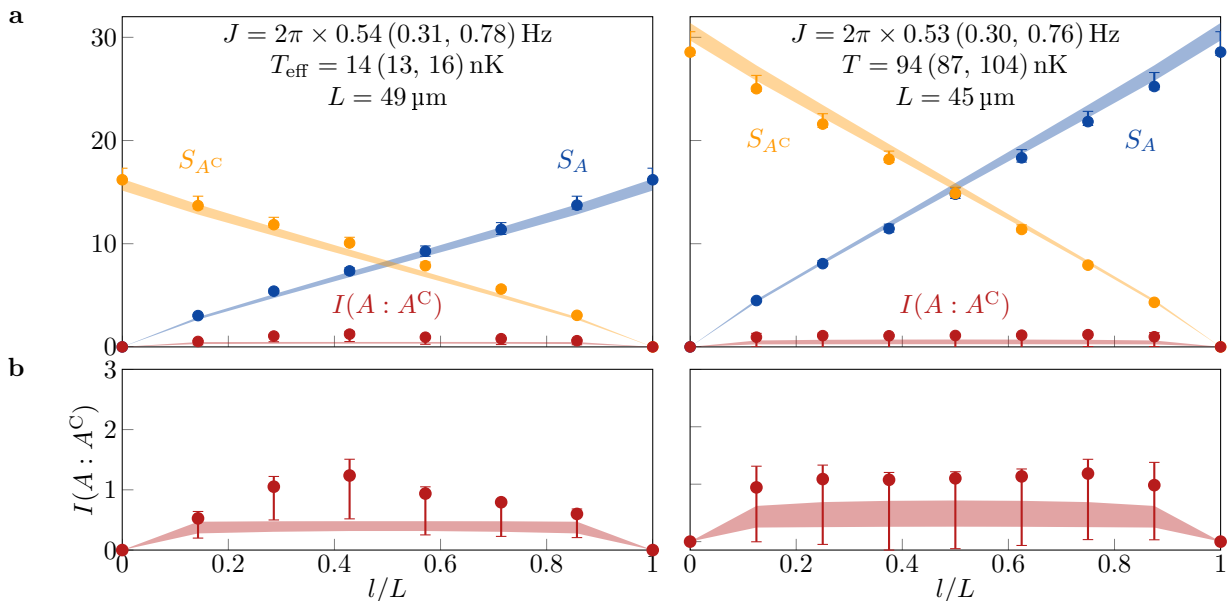


FIG. 5. **Additional results for area law of MI and volume law of vN entropy.** **a**, Experimental results for $I(A:A^C)$, S_A , and S_{A^C} , calculated based on $N = 7$ modes (left) and $N = 8$ modes (right) (see Fig. 2 for detailed explanation). The extracted parameters including their 95% confidence intervals in parenthesis are given above. **b**, Close up of the measured MI.

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SUPPLEMENTAL MATERIAL

Experimental realization and measurements

We realize a pair of strongly tunnelling coupled 1D superfluids by cooling down ^{87}Rb atoms in a double-well potential in an atom chip setup. The initial state is prepared by cooling the atoms directly into a double-well potential. The initial state in this case is a thermal equilibrium state with typical temperatures of 30-120 nK and linear atomic densities of $\approx 70 \mu\text{m}^{-1}$ in a box-like potentials with length $\approx 50 \mu\text{m}$.

To achieve lower effective temperatures, we first cool down the atoms in a dressed single-well potential, where

the first excited state is in the vacuum state. Slowly splitting the cloud into two parts, maps the ground state and the first excited state to two states with a smaller energy gap and symmetric and anti-symmetric wave functions. The resulting prethermalized state has a lower effective temperature in the anti-symmetric modes (relative degrees of freedom) [25]. This method has been deployed in the measurement with effective temperature $T_{\text{eff}} = 14$ nK presented in Fig. 4 and 5.

To probe the system, we turn off all the traps, and we let the atoms fall freely for 15.6 ms. We measure the projected 2D atomic density distributions via absorption imaging, from which we extract the relative phase between the two condensates for different points along the 1D direction z . Due to the destructive nature of the imaging process, we repeat the measurement hundreds of times to accumulate statistics. Thus, all the expectation values calculated are obtained through ensemble averaging.

Reconstruction of the initial full covariance matrix

To extract the full covariance matrix that characterises the state of the system, we use the quantum tomography method developed in [17]. Given that only one of the two canonical variables (the phase) is accessible through experimental measurements, its canonically conjugate variable (the density fluctuations) can be accessed indirectly by letting the system evolve under a harmonic Hamiltonian with known mode frequencies and measuring the phase at different times. For each of the har-

monic modes, the dynamics correspond to a rotation in phase space, so that over time, the initial density variance turns into phase variance and vice versa. Thus, we can fully reconstruct the initial covariance matrix from phase measurements at a sufficiently large number of different times. This reconstruction is done by first going to Fourier space, where the modes evolve independently and then fitting the data for the time evolution of the phase covariance of each mode to the known functions expressing this data in terms of the initial phase and density correlations.

However, one mode needs special treatment: The zero mode, i.e. the mode corresponding to zero momentum, therefore zero energy. This mode does not rotate in phase space; instead, it moves at a constant velocity. This means that the zero mode phase variance does not oscillate in time but grows as a quadratic function of time instead, an effect known as phase diffusion [26]. Moreover, because of the compactified nature of the phase-field, which means that phases differing by 2π should be considered identical, even though the phase grows in time with no bound, measurements can only observe its growth within the interval $[-\pi, +\pi]$. Therefore, the estimation of the initial phase and density correlations of the zero mode is different from that of the others. For the zero mode, we fit a quadratic function instead of an oscillatory function of time, and we restrict the fit to times before reaching the upper bound due to compactification.

More specifically, the dynamics are chosen to follow the Tomonaga-Luttinger liquid model with Hamiltonian

$$H_{\text{TLL}} = \int_0^L dz \left[g_{1\text{D}} (\delta\rho(z))^2 + \frac{\hbar^2 n_{1\text{D}}}{4m} (\partial_z \varphi(z))^2 \right] \quad (6)$$

For a hard-wall box trap the vanishing of the particle current at the edges of the system means that the effective boundary conditions are of Neumann type $\partial_z \varphi(x=0) = \partial_z \varphi(x=L) = 0$. In this case, using cosine eigenfunctions

$$f_n^\varphi(z) = \begin{cases} 2 \left(n \hbar \pi \sqrt{\frac{n_{1\text{D}}}{g_{1\text{D}} m}} \right)^{-1/2} \cos\left(n \frac{\pi}{L} z\right), & n > 0 \\ 1, & n = 0 \end{cases}$$

$$f_n^{\delta\rho}(z) = \begin{cases} -\frac{1}{L} \left(n \hbar \pi \sqrt{\frac{n_{1\text{D}}}{g_{1\text{D}} m}} \right)^{1/2} \cos\left(n \frac{\pi}{L} z\right), & n > 0 \\ -\frac{1}{L}, & n = 0 \end{cases}, \quad (7)$$

the Hamiltonian can be diagonalised in terms of cosine Fourier modes:

$$H = \frac{\hbar u}{2} \delta\rho_0^2 + \sum_{n=1}^{\infty} \frac{\hbar \omega_n}{2} [\delta\rho_n^2 + \varphi_n^2], \quad (8)$$

with

$$u = 2g_{1\text{D}}/\hbar L \quad (9)$$

$$\omega_n = ck_n. \quad (10)$$

Furthermore:

$$k_n = n\pi/L \quad (11)$$

$$c = \sqrt{g_{1\text{D}} n_{1\text{D}}/m} \quad (12)$$

$$g_{1\text{D}} = \hbar \omega_\perp a_s \frac{2 + 3a_s n_{1\text{D}}}{(1 + 2a_s n_{1\text{D}})}. \quad (13)$$

Here, c is the speed of sound and $g_{1\text{D}}$ the density-broadened 1D interaction strength [27], $a_s = 5.2 \text{ nm}$ is the three-dimensional scattering length [28] and $m = 1.44 \times 10^{-25} \text{ kg}$ the mass of a ^{87}Rb atom.

Unlike all higher modes, the zero mode, which corresponds to the constant in space eigenfunction, is not a harmonic oscillator mode of the Tomonaga-Luttinger liquid Hamiltonian (equation (6)). This is because only one of the canonical variables, $\delta\rho_0$, is present in the Hamiltonian for the zero mode. As a result, the time evolution of the zero mode is given by

$$\delta\rho_0(t) = \delta\rho_0(0) = \text{const}, \quad (14)$$

$$\varphi_0(t) = -u\delta\rho_0(0)t + \varphi_0(0), \quad (15)$$

which means that the phase variance grows with time as

$$\langle \varphi_0(t)^2 \rangle = \langle \varphi_0^2 \rangle_{t=0} - \langle \{\varphi_0, \delta\rho_0\} \rangle_{t=0} ut + \langle \delta\rho_0^2 \rangle_{t=0} u^2 t^2 \quad (16)$$

However, because of the compactified nature of the phase-field, its zero mode component φ_0 is not a well-defined, measurable operator. Only imaginary exponentials of the form $e^{in\varphi_0}$ for integer n are well-defined. Nevertheless, under the assumption the initial state is Gaussian in terms of the zero mode (as also for all other modes too) and given that it remains Gaussian under the dynamics following from H_{TLL} , we can derive the zero mode variance from the mean value of $e^{in\varphi_0}$ using the cumulant expansion formula for the special case of Gaussian random variables

$$\langle \exp(i\varphi_0) \rangle = \exp(i\langle \varphi_0 \rangle - \langle \varphi_0^2 \rangle/2) \quad (17)$$

From the above, we find

$$\langle \varphi_0^2 \rangle = -2 \log |\langle \exp(i\varphi_0) \rangle| \quad (18)$$

Therefore, in order to extract the zero mode part of the covariance matrix $\langle \varphi_0^2 \rangle_{t=0}$, $\langle \delta\rho_0^2 \rangle_{t=0}$ and $\langle \{\varphi_0, \delta\rho_0\} \rangle_{t=0}$ in the initial state we calculate the zero mode variance of the phase at each time from (17) and fit this with the theoretical formula (18).

Having the covariance matrix in the Fourier space for the first N modes, we use a discrete Fourier transformation based on the eigenfunctions (equation (8)) to calculate the covariance matrix Γ in the real space. We chose the cut-off based on the reconstructed occupation numbers. We only take into account modes with physical (positive) occupation numbers. This covariance matrix is used to calculate vN entropy and MI, as discussed in the main text and Fig. 1C.

Covariance matrix of Klein-Gordon model in thermal equilibrium

The theory predictions in our work are calculated based on the covariance matrix of the thermal equilibrium states of Klein-Gordon (KG) model (equation 3), which is given by [29]:

$$Q_{ij} = \frac{1}{2L} \frac{\hbar c}{Mc^2} \coth\left(\frac{Mc^2}{2k_B T}\right) + \frac{1}{L} \sum_{n=1}^N \frac{\hbar c}{\epsilon_n} \coth\left(\frac{\epsilon_n}{2k_B T}\right) \cos(k_n z_i) \cos(k_n z_j) \quad (19)$$

$$R_{ij} = 0 \quad (20)$$

$$P_{ij} = \frac{1}{2L} \frac{Mc^2}{\hbar c} \coth\left(\frac{Mc^2}{2k_B T}\right) + \frac{1}{L} \sum_{n=1}^N \frac{\epsilon_n}{\hbar c} \coth\left(\frac{\epsilon_n}{2k_B T}\right) \cos(k_n z_i) \cos(k_n z_j), \quad (21)$$

with the dispersion relation $\epsilon_n = \sqrt{\hbar^2 k_n^2 c^2 + M^2 c^4}$. Here, M is the KG mass, T the temperature, N the UV cut-off and the rest of the parameters are defined before. The next section will explain how M and T are estimated based on the measured data. To include the effect of the finite imaging resolution, we convolve the theoretical calculations with a Gaussian point-spread function with a standard deviation $\sigma_{\text{PSF}} \approx 3 \mu\text{m}$ [27] (see next section).

Estimation of the temperature and Klein-Gordon mass

In order to compute theoretical predictions for the mutual information in the initial state based on the assumption that these are thermal states of the KG model, we need to estimate two effective parameters, the mass and the temperature. We do this by fitting the results of the tomographic reconstruction for the mode variances to those corresponding to KG thermal states. Given that the modes are decoupled from each other both initially and throughout the tomography dynamics, the estimation of the KG mode frequency can be done independently for each mode. Having estimated the mode fre-

quencies, we can then verify if they follow the theoretical dispersion relation of the KG model and extract the corresponding mass parameter by a fit.

The relation between the post-quench quadratures and the initial (pre-quench) KG state with mode occupation number N_{0n} is given by

$$\langle \varphi_n^2 \rangle = \frac{\omega_n}{\omega_{0n}} \left(N_{0n} + \frac{1}{2} \right) \quad (22)$$

$$\langle \delta \rho_n^2 \rangle = \frac{\omega_{0n}}{\omega_n} \left(N_{0n} + \frac{1}{2} \right) \quad (23)$$

where ω_{0n} and ω_n are the pre- and post-quench mode frequencies respectively. Both $\langle \varphi_n^2 \rangle$ and $\langle \delta \rho_n^2 \rangle$ are achieved via tomography and ω_n is given by the equation (10). From the equations (22) and (23) we can calculate both ω_{0n} and N_{0n} :

$$N_{0n} = \sqrt{\langle \varphi_n^2 \rangle \langle \delta \rho_n^2 \rangle} - \frac{1}{2} \quad (24)$$

$$\omega_{0n} = \omega_n \sqrt{\frac{\langle \delta \rho_n^2 \rangle}{\langle \varphi_n^2 \rangle}}. \quad (25)$$

Assuming the initial state is thermal, we use the following fit function to extract the temperature, $T = (\beta k_B)^{-1}$:

$$N_{0n}^{\text{fit}} = \frac{\exp(-k_n^2 \sigma_{\text{PSF}}^2)}{\exp(\hbar \omega_{0n} \beta) - 1}. \quad (26)$$

Note that we also introduce the effect of the imaging system by multiplying the modes with a Gaussian point spread function with width σ_{PSF} which corresponds to a convolution with a Gaussian point-spread function in the real space [27]. To extract the KG mass, we fit the KG dispersion relation, $\omega_{0n}^{\text{fit}} = \sqrt{c^2 k_n^2 + M^2 c^4 / \hbar^2}$, to the calculated ω_{0n} .

The single-particle tunnelling rate, J is related to the KG mass via the following equation:

$$J = \frac{M^2 c^3}{8 \hbar \sqrt{m g_{1D} n_{1D}}}. \quad (27)$$

The correlation length, as mentioned in the main text, is then given by $l_C = \sqrt{\hbar / (4mJ)}$ [27].

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