Fundamental noise in matter interferometers

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These lecture notes discuss two effects which contribute to the reduction of the interference fringe contrast in matter interferometers. The first effect is the shot noise arising from a finite number of atoms used in experiments. Focusing on a single shot measurement we provide explicit calculations of the full distribution functions of the fringe contrast for the interference of either the coherent or the number states of atoms. Another mechanism of the suppression of the amplitude of interference fringes discussed in these lecture notes is the quantum and thermal fluctuations of the order parameter in low dimensional condensates. We summarize recent theoretical and experimental studies demonstrating that suppression of the interference fringe contrast and its shot to shot variations can be used to study correlation functions within individual condensates. We also discuss full distribution functions of the fringe amplitudes for one and two dimensional condensates and review their connection to high order correlation functions. We point out intriguing mathematical connections between the distribution functions of interference fringe amplitudes and several other problems in field theory, systems of correlated electrons, and statistical physics.

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I. INTRODUCTION

A. Interference experiments with cold atoms

From the earliest days of quantum mechanics its probabilistic nature was the cause of many surprises and controversies [1]. Perhaps the most unusual manifestation of the quantum uncertainty is a quantum noise: measurements performed on identical quantum mechanical systems can produce results which are different from one experimental run to another. At the level of a single particle quantum mechanics, the quantum noise is no longer a research topic but is discussed in undergraduate physics textbooks [2]. However, the situation is different when we talk about quantum mechanics of many-body systems. One can ask seemingly simple questions to which there is no obvious answer: does it still make sense to talk about quantum noise when discussing measurements on many-body quantum states? How does the quantum noise manifest itself? Can one use this noise to extract nontrivial information about the system?

The idea of the quantum noise analysis of many-particle systems is common to many areas of condensed matter physics [3, 4, 5, 6] and quantum optics [7, 8, 9, 10, 11, 12]. In the field of ultracold atoms it has been successfully employed in a variety of recent experiments [13, 14, 15, 16, 17, 18, 19] with many more theoretical proposals awaiting their turn [20, 21, 22, 23, 24, 25, 26]. These lecture notes address a very specific experimental probe of the cold atomic ensembles – interference experiments. Our discussion focuses on a variety of interesting and important phenomena which originate from the fundamental quantum and/or thermal noise of cold atoms condensates and can be studied in interference experiments. Although focused on the specific type of experiments, the general methodology discussed in these lecture notes can be extended to a variety of other measurements on systems of cold atoms.

Interference experiments constitute an important part of the modern toolbox for studying ultracold atoms. Original experiments used large three dimensional Bose-Einstein condensates (BEC) to demonstrate macroscopic coherence [27]. More recently interference experiments have been done with one and two dimensional condensates [28, 29, 30, 31, 32, 33] and demonstrated the important role of fluctuations in low dimensional systems. Matter interferometers using cold atoms [29, 30, 34, 35, 36, 37, 38, 39, 40, 41, 12, 43, 44, 45, 46, 47, 48, 49, 50] have been considered for applications in accelerometry, gravitometry, search for quantum gravity, and many other areas (for a review see Ref. [51]). Interference experiments have been used to measure the condensate formation [52, 53] as well as the critical properties of the BEC transition [54]. What is common to most interference experiments is the focus on the phase of interference patterns. Suppression of the fringe contrast is considered to be a spurious effect caused by noise and fluctuations. On the contrary, these lecture notes focus on understanding physical phenomena that underlie the imperfect visibility of interference fringes. As we discuss below, suppression of the fringe visibility comes from fundamental physical phenomena, such as the noise intrinsic to performing a classical measurement on a quantum mechanical wave function (shot noise) or classical and quantum fluctuations of the order parameter.

In these lecture notes we discuss how one can use analysis of the contrast of interference fringes to learn about fluctuations of the order parameter [32, 33, 55, 56, 57, 58]. We will demonstrate that important information is contained not only in the average contrast but also in its shot to shot variations. For example, when we discuss fluctuating condensates, we will show that high moments of the interference fringe amplitudes contain information about high order correlation functions and thus provide valuable information about the system.

The basic scheme of interference experiments is shown in Fig. 1. Originally two condensates are located at distance $d$ away from each other. At some point they are allowed to expand ballistically [59] until sizes of the clouds become much larger than the original separation between the clouds $d$. After the expansion the density is measured by shining a laser beam through the cloud. Interference leads to the appearance of the density modulation at a wave vector $Q = m d/\hbar t$ (see Fig. 1 and discussion in section 11). When the two condensates are coherent, the position of interference fringes is determined by the relative phase between the two clouds. Surprisingly the interference pattern will be observed even for two independent condensates which do not have a well defined relative phase (see e.g. Fig. 2). To beginning readers it may seem confusing that we can observe interference in the absence of coherence between the two clouds. Or even more confusing, we discuss interference patterns when both clouds are number states and phases of individual condensates are not well defined. Several theoretical frameworks have been introduced to understand the origin of interference patterns in the absence of phase coherence [60, 61, 62, 63, 64]. In this paper we explain
FIG. 1: Schematic view of the interference experiments with ultracold atoms. Two condensates, originally separated by the distance $d$, are released from the traps and expand until the clouds overlap. The imaging beam measures the density of the atoms after the expansion. Quantum interference leads to the periodically modulated density projected on the screen. Projected density image is taken from the actual experimental data of Hadzibabic et al. [32].

FIG. 2: Interference pattern from the first set of experiments with sodium atoms. Taken from Ref. [27].

the origin of interference fringes using the language of correlation functions and point out connections to Hanbury Brown and Twiss (HBT) experiments [2] in optics. This section provides a simple heuristic picture and a more formal discussion is left to section II.

What one measures in experiments is the density profile after the expansion $\rho(r)$. Interference pattern appears as the density modulation $\rho(r) = \rho_Q e^{-iQr} + \text{c.c.} + \text{const}$. The absolute value of $\rho_Q$ determines the amplitude of interference fringes and its phase defines the position of the fringe maxima and minima. Schematically we can write $\rho_Q \sim e^{i\phi_1 - i\phi_2}$, where $\phi_{1,2}$ are phases of the two condensates before the expansion (see discussion in section II). In the absence of coherence $\langle e^{i\phi_1 - i\phi_2} \rangle = 0$, which implies that $\langle \rho_Q \rangle = 0$. Vanishing of the average, however, does not mean that interference fringes are absent in each individual shot. In the present case it only shows that the phase of interference fringes is random in each shot. We remind the readers that taking an expectation value in quantum mechanics implies averaging over many measurements. On the other hand we can focus on the amplitude of interference fringes and accept the fact that we can not predict their phase. Then we need to consider the quantity which does not vanish after averaging over the unpredictable phase difference. One such quantity is given by the density-density correlation function $\rho(r)\rho(r') = |\rho_Q|^2(e^{iQ(r-r')} + \text{c.c.}) + \text{other terms}$. The right hand side of the last equation does not vanish when we average over random phases $\phi_{1,2}$, and we find finite expectation value

$$\langle \rho(r)\rho(r') \rangle = 2|\rho_Q|^2 \cos(Q(r-r')) + \text{const}. \quad (1)$$

What this correlation function tells us is that in a single shot we can not predict whether at a given point $r$ we will have a minimum or a maximum of the density modulation. However what we can say is that if there is a maximum at point $r$, it will be followed by another maximum a distance $2\pi/Q$ away. While this simple argument explains the
origin of interference patterns from independent condensates, it leaves many questions unanswered. For example, it is not obvious how accurately one can represent two independent condensates using states with a well defined but unknown phase difference. Also it is not clear how to generalize this analysis to elongated condensates, when we need to go beyond the single mode approximation and include phase fluctuations within individual systems. These lecture notes will present a uniform approach for addressing these and many other questions.

When the focus of interference experiments is on measuring the phase, one usually averages interference patterns obtained in several shots. The result is easy to interpret: an average of many experimental runs is precisely what we define as a quantum mechanical average. However in experiments with independent condensates, summing interference patterns is not appropriate. The phase of interference patterns is random from shot to shot and adding individual images washes out interference fringes completely (for a nice experimental demonstration see Ref. [65]). Hence, in this case one needs to focus on interference patterns obtained in individual shots. In the absence of averaging, a single shot measurement contains noise. Thus to characterize such experiments we need to provide both the average value and the shot to shot fluctuations of the fringe contrast. The most comprehensive description of the fluctuating variable comes from providing its full distribution function. Theoretical calculations of the distribution functions of the fringe contrast will be the central part of these lecture notes.

It is useful to point out the analogy between the approach discussed in this paper and the famous Hanbury Brown and Twiss experiments in optics. The original motivation for HBT experiments came from astronomy: the goal was to measure the angle between two incoherent stellar sources such as two different points on the surface of the star. Since the two sources are incoherent, this can not be done using a single detector: first order interference is absent and the measured signal is simply the sum of the two intensities (12). The insight of HBT was to use two detectors and measure the correlation function of the two intensities as a function of the relative distance between the two detectors. One finds that this correlation function is given by

\[
\langle I(r)I(r') \rangle \sim \cos (\langle k_1 - k_2 \rangle (r - r')) + \text{const},
\]

where \(k_1, k_2\) are wave vectors of photons arriving from the two points on the surface of the star, \(r, r'\) are positions of the detectors, and \(I(r), I(r')\) are the intensities measured in the two detectors (see Fig. 3). Hence, the main idea of HBT experiments is that information is contained not only in the average signal \(\langle I(r) \rangle\), but also in the noise. Such noise can be characterized by looking at higher order correlations. In astronomy, HBT experiments were used to measure several important properties of distant stars, including their angular sizes and the surface temperature [66].

There is an obvious analogy between equations (1) and (2), but there is also an important difference. HBT stellar interferometers operate in real time, and averaging over time is built into the measuring procedure. In these experiments, fluctuations of \(I(r)I(r') - \langle I(r)I(r') \rangle\), which would correspond to higher order correlations in \(I(r)\), are not easy to measure. On the other hand interference experiments with cold atoms are of a single-shot type: each measurement is destructive and gives a certain density profile \(\rho(r)\). A single image contains information not only about the two point correlation function \(\langle \rho(r)\rho(r') \rangle\), but about higher order correlations as well. The most important information for our purposes is contained in the interference pattern at wave vector \(Q\). Essentially each interference pattern constitutes a classical measurement of the quantum mechanical operator \(\rho Q\). We remind the readers that we expect the phase of \(\rho Q\) to be random, so the quantity of interest will be \(|\rho Q|^2\). By performing measurements several times we will find not only the average value of this operator, but also its higher moments. Ultimately we should be able to reconstruct the entire distribution function for \(|\rho Q|^2\). So the simplified and idealized procedure that we

![FIG. 3: Schematic view of Hanbury Brown and Twiss noise correlation experiment as an example of intensity interferometry. Detectors at positions \(r, r'\) measure the intensity of light coming from two distant incoherent sources. The ”correlator” (denoted by the box) measures the coincidence events and thus the intensity-intensity correlation function.](image-url)
analyze is the following: one performs interference experiments many times. Each experiment is analyzed by doing a Fourier transform of the density to extract $\rho_Q$. The histogram of the measured values of $|\rho_Q|^2$ will be the main subject of these lecture notes. We will demonstrate the wealth of information that can be extracted from analysis of such histograms.

As a passing note, we mention that the setup considered in Fig. 1 is not the only possible configuration for interference experiments. In another common setup one makes several copies of the same cloud using Bragg pulses, and observes an interference between them.

It is useful to put our work in the general perspective of noise analysis in physical systems. Understanding photon fluctuations is at the heart of modern quantum optics and provides a basis for creation, detection, and manipulation of non-classical states of light. The field of quantum optics has a long and fruitful tradition of using the higher-order correlation functions as well as the shot noise to characterize the quantum states of light. The notion of higher-order degree of coherence was first introduced by R. Glauber in 1963, also by Klauder and Sudarshan and by Mandel and Wolf. The knowledge of photoelectron counting distribution function reveals such non-classical features of light as antibunching, sub-Poissonian statistics and probe of violation of Bell inequalities. In particular, third-order correlations provide a test for distinguishing between quantum and hidden variable theories in a way analogous to that provided by the Greenberger-Horne-Zeilinger test of local hidden variable theories. Interference of independent laser beams was first observed in Ref. and stimulated a number of theoretical studies (for reviews see Refs. [10, 77, 78]).

In condensed matter physics, noise analysis was also suggested as a powerful approach for analyzing electron systems. It was demonstrated theoretically that in certain mesoscopic systems current fluctuations should contain more information than the average current itself. In particular, the third and higher moments contain important quantum information on interaction effects, entanglement and relaxation processes (see e.g. Refs. [79, 80]). Specific proposals exist for detecting statistics of quasi-particles [81], understanding transmission properties of small conductors [4] and observing entanglement between electrons [8]. Perhaps the most spectacular experimental success of the noise analysis in electronic systems has been the demonstration of the fractional charge of quasiparticles in the fractional quantum Hall regime. However generally the noise analysis in condensed matter systems did not become the detection tool of the same prominence as in quantum optics. The main reason for this is the excruciating difficulty of the noise measurements in solid state experiments. One often needs to measure a signal which is only a part in a million of the unwanted technical noise.

In the field of ultracold atoms experiments analyzing quantum noise are only starting. However we have already seen spectacular successes in several recent experiments. Analysis of noise correlations in the time of flight experiments was used to demonstrate fermionic pairing as well as HBT type correlations for atoms in optical lattices (see also I. Bloch’s lectures in this school). Single atom detectors have been used to demonstrate HBT noise correlations for cold atoms. Strongly interacting systems of cold atoms are expected to realize analogues of important models of condensed matter systems. Being able to study noise in such systems should provide an important new perspective on strongly correlated states of matter and have a profound effect on many areas of physics. We hope that these lecture notes will stimulate more experimental work in analyzing noise in interference experiments. The first success in this direction was the recent observation of the Berezinskii-Kosterlitz-Thouless (BKT) transition by Hadzibabic et al. in Ref. [32].

B. Fundamental sources of noise in interference experiments with matter

Two fundamental sources of fluctuations in the amplitude of interference fringes are the shot noise and the order parameter fluctuations within individual condensates.

Shot noise comes from the finite number of atoms used in the experiments. Let us discuss limiting cases first. Consider an interference experiment with one atom. Before the expansion the atom is in a perfect superposition of being between the two wells. After the expansion we get a perfectly periodic wave function $\psi(r) = 2C \cos(Qr+\phi)$ (for a more detailed discussion see section 11). The expectation value of the density operator is $\langle \rho(r) \rangle = |\psi(r)|^2 = 2|C|^2(\cos(\sqrt{|\langle \rho|Q^2 \rangle}| + 1)$. However this average value will not be measured in a single shot. A single measurement finds the atom at a single point. The expectation value of the density determines probabilities with which we can find atom at any given point, but in a single measurement we collapse a quantum mechanical wave function and observe the atom at one point only. Can one reconstruct the entire amplitude of the interference pattern $|\rho_Q| = |C|^2e^{-i\phi}$ from a single measurement? Obviously the answer is no. In the opposite case of a very large number of atoms in the same single particle state one should be able to reconstruct a complete interference pattern already from a single measurement, since the measurement of positions of many atoms performs a statistical averaging implicit in quantum mechanics. In the general case of experiments with a finite number of atoms, the question arises how well one can determine the amplitude of interference fringes from doing a single shot measurement. Formulated more accurately the problem is...
Fluctuations of the order parameter are particularly important for low dimensional systems. If the condensates are confined in one or two dimensions, then the true long-range order may not exist. Rigorous theorems forbid true long range order in two dimensional systems at finite temperature and in one dimension even at zero temperature. What this means is that low dimensional condensates can not be characterized by a single phase and we need to take into account spatial fluctuations of the order parameter. Effects of such fluctuations on interference experiments are illustrated schematically in Fig. 4. Two one dimensional clouds expand in the transverse direction. Each point along the condensates has a local interference pattern, but in the presence of phase fluctuations (either thermal or quantum), these patterns are not in phase with each other. It is natural to define the net interference amplitude from the density integrated over the axis of the system (the so called columnar density). In many experiments such integration is done by the measurement procedure itself. For example, systems such as shown in Fig. 4 originally had imaging done along the axis of the interferometer. Then the laser beam integrates the atomic densities within the imaging length. Integrating over local interference patterns which are not in phase with each other leads to a reduced contrast of the net interference fringes.

In earlier literature smearing of interference patterns by fluctuations was considered an unwanted effect. The point of view presented in these notes is quite the opposite. Suppression of the fringe contrast is an interesting effect which tells us about important phenomena in Bose condensates, such as thermal and quantum fluctuations of the order parameter. By analyzing such suppression we can extract non-trivial information about the system. In particular, it has been shown in Ref. that the scaling of the average interference signal with the observation area contains information about the two-point correlation functions within each cloud. Recent experiments by Hadzibabic et al. used this approach to observe the BKT transition in two dimensional condensates (see discussion in section IV). We also note that such experiments can be used to extract information which is difficult to obtain by other means. For example, in section IV we discuss that analysis of the high moments of the contrast tells us about high order correlations within individual clouds.

These lecture notes are organized as follows. In section II we discuss how interference fringes appear for ideal non-interacting 3D BEC’s at zero temperature. In section III we analyze the shot noise for ideal condensates. In these lecture notes ideal condensates are understood as clouds of non-interacting atoms which before the expansion occupy a single mode within each of the traps. The problem of interference of independent condensates of ideal bosons has been extensively analyzed in the literature before. In an important recent paper, Polkovnikov showed that the variance of the fringe amplitude decreases as an inverse power of $N$, with a non-universal coefficient which contains information about the state of each cloud (e.g. coherent states vs Fock states). In this paper we develop a general formalism for calculating the full distribution functions of the fringe amplitudes in interference experiments with ideal condensates with a finite number of atoms. We apply this formalism to obtain distribution functions for several experimentally relevant cases such as states with a well defined phase difference between the two clouds and Fock states of atoms. Effects of the order parameter fluctuations are
II. INTERFERENCE OF IDEAL CONDENSATES

In this section we discuss why interference fringes appear for ideal non interacting BEC’s at zero temperature. We follow Ref. [64], and introduce notations for subsequent sections. First we consider the case of two clouds with a well defined relative phase, where appearance of interference fringes can be understood at a single particle level. Then we show that almost ideal interference fringes appear even when two expanding clouds are uncorrelated, provided that the number of particles in each cloud is large.

A. Interference of condensates with a well defined relative phase

1. Basics of interference experiments. First quantized representation

To illustrate how interference fringes arise, let us start by considering a simple case of two BEC clouds with a well defined relative phase. Here we neglect interactions between atoms, so initially all atoms are assumed to be in the same single-particle state (single mode approximation). After the confining potential is removed, the single-particle state evolves with time, but many body wave function remains in the product state. The interference appears as a result of single particle wave function evolution, which can be studied in detail.

Normalized single particle wave functions for two clouds will be denoted as $\psi_1(r, t)$ and $\psi_2(r, t)$, and the initial relative phase is $\varphi$. If the total number of particles equals $N$, then the complete wave function of the system in the first quantized notations at any moment of time is given by

$$\Psi(r_1, ..., r_N, t) = \prod_{n=1}^{N} \frac{1}{\sqrt{2}} (\psi_1(r_n, t)e^{i\varphi/2} + \psi_2(r_n, t)e^{-i\varphi/2}).$$

This wave function satisfies the proper symmetry requirements for permutations of $r_i$ and $r_j$, and evolution of $\psi_1(r, t)$ and $\psi_2(r, t)$ is controlled by the single particle Schrödinger equation. Initial overlap of the states $\psi_1(r) \equiv \psi_1(r, 0)$ and $\psi_2(r) \equiv \psi_2(r, 0)$ is assumed to be negligible:

$$\int dr \psi_1^*(r) \psi_2(r) \approx 0.$$  

The expectation value of the total density corresponding to the wave function (3) is

$$\langle \rho(r, t) \rangle = \frac{N}{2} (|\psi_1(r, t)|^2 + |\psi_2(r, t)|^2 + 2\text{Re}[e^{i\varphi}\psi_1(r, t)\psi_2^*(r, t)]).$$

The expectation value of the total density displays an interference pattern due to the last term of Eq. (5). As a simple example, let us assume that $\psi_1(r, t)$ and $\psi_2(r, t)$ are initially in the Gaussian states centered at points $\pm d/2$, and their widths are $R_0 \ll d$. Then the evolution of the single particle wave functions can be simply calculated, and the result is

$$\psi_1(r, t) = \frac{1}{(\pi R_t^2)^{3/4}} e^{-\frac{(r-d/2)^2 + i\varphi}{2R_t^2}},$$

$$\psi_2(r, t) = \frac{1}{(\pi R_t^2)^{3/4}} e^{-\frac{(r+d/2)^2 + i\varphi}{2R_t^2}},$$

where the widths of the wave packets, $R_t$, at time $t$ are given by

$$R_t^2 = R_0^2 + \left(\frac{\hbar t}{mR_0}\right)^2.$$
FIG. 5: Schematic view of the interference experiment with 3D condensates.

We will be interested in the regime, when the sizes of the clouds $R_t$ are much larger than original distance between the clouds, that is

$$R_t \gg d \gg R_0.$$  \hfill (9)

In this regime, the clouds overlap strongly, and the real parts of the exponents in Eqs. (6) and (7) are responsible for the broad overall density profile. Imaginary parts in the same exponents give rise to interference effects in the last term of Eq. (5). Thus the interference part of the density is equal to

$$N \left( \frac{\pi R_t^2}{R_0^2} \right)^{3/2} e^{-\frac{2\pi^2 d^2}{R_t^2}} \cos \left( \frac{\hbar}{mR_0^2R_t^2} t + \varphi \right).$$  \hfill (10)

For sufficiently large $t$, one can substitute $R_t \approx \hbar t/mR_0$, and obtain oscillations of the density at wave vector $Q = m\bar{d}/\hbar t$, with positions of the minima and the maxima controlled by the relative phase $\varphi$. The Fourier transform of the density at wave vector $Q$ is

$$\langle \rho_Q \rangle = \langle \int d\mathbf{r} \rho(\mathbf{r}, t) e^{iQ \mathbf{r}} \rangle \approx N^2 e^{-i\varphi}.$$  \hfill (11)

Physically $Q$ can be understood as the momentum difference of the two particles which have been released from the two traps and arrive simultaneously at the detection point $\mathbf{r}$. This can be seen from the following quasiclassical argument. A particle released from the condensate one and detected at time $t$ at point $\mathbf{r}$ has momentum

$$Q_1 = m(\mathbf{r} - \mathbf{d}/2)/(\hbar t).$$  \hfill (12)

During the expansion this particle picks up a phase $Q_1(\mathbf{r} - \mathbf{d}/2)$. Analogously a particle originating from the condensate two has momentum

$$Q_2 = m(\mathbf{r} + \mathbf{d}/2)/(\hbar t)$$  \hfill (13)

and picks up a phase $Q_2(\mathbf{r} + \mathbf{d}/2)$. The interference pattern arises from the oscillating structure of the phase difference with the wave vector of oscillations

$$Q = Q_1 - Q_2 = m\bar{d}/\hbar t.$$  \hfill (14)

This simple argument shows that as long as the original sizes of the clouds are much smaller than the distance between them, after sufficient expansion one should observe oscillations of the density at the wave vector $Q$ determined by the distance between the clouds.

The interference patterns which we introduced up to this point appear as a result of the time evolution of single particle states. The many-body nature of the state comes into Eq. (5) only as a prefactor $N$. In principle, one could have done the same experiment with only one particle. The same result as Eq. (14) can be obtained in this case by doing experiments many times and averaging over individual experiments: in each particular realization the particle is observed at some random point $\mathbf{r}$ with probability $\langle \rho(\mathbf{r}, t) \rangle$. In experiments with a large number of atoms, $N$, used in each shot, each absorption image is a result of $N$ measurements of single particle wave functions. This performs the statistical averaging implicit in quantum mechanics, and leads to the density profile close to Eq. (5) in each shot.
2. Second quantized representation

To set up the stage for later we will now present the discussion leading to Eq. (11) using the second-quantized formalism. Wave function (9) in the second quantization at \( t = 0 \) can be written as

\[
|\varphi, N, t \rangle = \frac{1}{\sqrt{(2^N N!)^{1/2}}(a_1^\dagger e^{i\varphi/2} + a_2^\dagger e^{-i\varphi/2})^N |0\rangle. \tag{15}
\]

Here \( a_1^\dagger \) and \( a_2^\dagger \) are creation operators for clouds one and two:

\[
a_i^\dagger = \int dr \psi_i(r) \hat{\psi}^\dagger(r). \tag{16}
\]

\( \hat{\psi}(r) \) is the second quantized operator for the boson field, which satisfies the usual commutation relations

\[
\left[ \hat{\psi}(r'), \hat{\psi}^\dagger(r) \right] = \delta(r' - r), \quad \left[ \hat{\psi}(r'), \hat{\psi}(r) \right] = \left[ \hat{\psi}^\dagger(r'), \hat{\psi}^\dagger(r) \right] = 0. \tag{17}
\]

Operators \( a_i, a_j \) and their conjugates satisfy the canonical boson commutation relations:

\[
[a_i, a_j^\dagger] = \delta_{ij}, \quad [a_i, a_j] = [a_i^\dagger, a_j^\dagger] = 0. \tag{18}
\]

Different initial states can be simply written using the Fock basis of operators \( a_1^\dagger \) and \( a_2^\dagger \). For example, the initial state for two independent condensates with \( N_1 \) and \( N_2 \) particles in clouds 1 and 2 can be conveniently written as

\[
|N_1, N_2 \rangle = \frac{1}{\sqrt{N_1!N_2!}}(a_1^\dagger)^{N_1}(a_2^\dagger)^{N_2}|0\rangle. \tag{19}
\]

Initial state written in the Fock basis of operators \( a_i^\dagger \) and \( a_j^\dagger \) contains all information about properties of the interference amplitudes. During the free expansion the occupation numbers of states one and two do not change, and only the single particle wave functions \( \psi_1(r, t) \) and \( \psi_2(r, t) \) evolve. After the expansion, the many-body wave function at time \( t \) can be obtained from the initial state written in the Fock basis of operators \( a_1^\dagger \) and \( a_2^\dagger \) using substitutions

\[
\begin{align*}
a_1^\dagger & \to \int dr \psi_1(r, t) \hat{\psi}^\dagger(r), \tag{20} \\
a_2^\dagger & \to \int dr \psi_2(r, t) \hat{\psi}^\dagger(r). \tag{21}
\end{align*}
\]

For example, substituting Eqs. (20)-(21) into Eq. (13), the wave function (3) considered earlier is written as

\[
|\varphi, N, t \rangle = \frac{1}{\sqrt{(2^N N!)^{1/2}}} \left( \int dr \psi_1(r, t) e^{i\varphi/2} + \psi_2(r, t) e^{-i\varphi/2} \right)^N |0\rangle. \tag{22}
\]

In the long time limit (9) considered earlier, single-particle wave functions \( \psi_1(r, t), \psi_2(r, t) \) can be written as

\[
\begin{align*}
\psi_1(r, t) & = u_1(r, t) e^{iQ_1 \cdot r}, \tag{23} \\
\psi_2(r, t) & = u_2(r, t) e^{iQ_2 \cdot r}, \tag{24}
\end{align*}
\]

where \( Q_1, Q_2 \) are defined by Eqs. (12)-(13) and \( u_1(r, t), u_2(r, t) \) are slowly varying real functions, which determine the overall density profiles. Since clouds overlap strongly after the expansion, these functions are normalized according to

\[
\begin{align*}
\int u_1(r, t)^2 dr & = 1, \quad \int u_2(r, t)^2 dr = 1, \tag{25} \\
\int u_1(r, t) u_2(r, t) dr & \approx 1. \tag{26}
\end{align*}
\]

The operator, which corresponds to the amplitude of density oscillation at wave vector \( Q \) is written in the second quantized notations as

\[
\hat{\rho}_Q = \int dr \hat{\rho}(r) e^{iQ \cdot r} = \int dr \hat{\psi}^\dagger(r) \hat{\psi}(r) e^{iQ \cdot r}. \tag{27}
\]
To find out the statistical average of the amplitude of density oscillations for state \( |\varphi, N, t\rangle \), we need to evaluate the following matrix element:

\[
\langle \rho_Q \rangle = \langle \varphi, N, t | \hat{\rho}_Q | \varphi, N, t \rangle = \langle \varphi, N, t | \int dr \hat{\psi}_l(r) \hat{\psi}(r) e^{iQr} | \varphi, N, t \rangle. \tag{28}
\]

To evaluate such matrix elements, first we need to know how annihilation operator \( \hat{\psi}(r) \) acts on a state \( |\varphi, N, t\rangle \). Since \( |\varphi, N, t\rangle \) is obtained from \( |\varphi, N\rangle \) by substitutions \((20)-(21)\), it is easy to see that

\[
\langle \varphi, N, t | \hat{\psi}_l(r) \hat{\psi}(r) | \varphi, N, t \rangle = \langle \varphi, N | \left( a_1^\dagger u_1(r, t) e^{-iQr} + a_2^\dagger u_2(r, t) e^{-iQ_2r} \right) \left( a_1 u_1(r, t) e^{iQ_1r} + a_2 u_2(r, t) e^{iQ_2r} \right) | \varphi, N \rangle,
\]

where \( |\varphi, N\rangle \) is defined in Eq. \((15)\) and is written only in terms of \( a_1^\dagger \) and \( a_2^\dagger \). Integration over \( dr \) in \((28)\) can be done using normalization \((20)\), and assuming that \( u_i(r, t) \) vary at scales much larger than \( 1/Q \). Since \( Q = Q_1 - Q_2 \), evaluation of Fourier transform picks only the product of the first term in the first parentheses and of the second term in the second parentheses of the equation above. We obtain

\[
\langle \rho_Q \rangle = \langle \varphi, N | a_1^\dagger a_2^\dagger \int u_1(r, t) u_2(r, t) dr | \varphi, N \rangle = \langle \varphi, N | a_1^\dagger a_2^\dagger \int u_1(r, t) u_2(r, t) dr | \varphi, N \rangle = \langle \varphi, N | \frac{N}{2} e^{-i\varphi} | \varphi, N \rangle = \frac{N}{2} e^{-i\varphi}, \tag{29}
\]

which is the same as obtained from a single particle discussion in Eq. \((11)\).

The example above illustrates how the matrix elements of many-particle operators at time \( t \) can be evaluated using initial states written in the Fock basis of operators \( a_1^\dagger \) and \( a_2^\dagger \). In general, when one needs to evaluate an expectation value of some normal ordered combination of operators \( \hat{\psi}(r) \) and \( \hat{\psi}_l(r) \) over the final state at time \( t \), one needs to make substitutions

\[
\hat{\psi}(r) \rightarrow a_1 u_1(r, t) e^{iQ_1r} + a_2 u_2(r, t) e^{iQ_2r}, \tag{30}
\]

\[
\hat{\psi}_l(r) \rightarrow a_1^\dagger u_1(r, t) e^{-iQ_1r} + a_2^\dagger u_2(r, t) e^{-iQ_2r}, \tag{31}
\]

and evaluate matrix elements over the \( t = 0 \) state, written in the Fock basis of operators \( a_1^\dagger \) and \( a_2^\dagger \). It is important that the expression needs to be normal ordered using commutation relations \((17)\) before making substitutions \((30)-(31)\), since after substitutions \((30)-(31)\) fields \( \hat{\psi}(r), \hat{\psi}_l(r) \) do not satisfy the exact commutation relations \((17)\).

Another way of seeing this is to realize that in Heisenberg representation (see e.g. chapter 6 of Ref. \((102)\)) substitutions \((30)-(31)\) perform the time evolution of a product of boson annihilation operators

\[
\hat{\psi}(r_1, t) \ldots \hat{\psi}(r_n, t) = e^{i\hat{H}t} \hat{\psi}(r_1) \ldots \hat{\psi}(r_n) e^{-i\hat{H}t} \tag{32}
\]

only when this product acts on states of the form

\[
|\Psi_0\rangle = \frac{1}{\sqrt{N_1! N_2!}} (a_1^\dagger)^{N_1} (a_2^\dagger)^{N_2} |0\rangle. \tag{33}
\]

Thus for calculating the expectation values of the form

\[
|\Psi_0\rangle e^{i\hat{H}t} \hat{\rho}(r_1) \ldots \hat{\rho}(r_n) e^{-i\hat{H}t} |\Psi_0\rangle \tag{34}
\]

one first needs to normal order \( \hat{\rho}(r_1) \ldots \hat{\rho}(r_n) \) using commutation relations \((17)\) as

\[
\hat{\rho}(r_1) \ldots \hat{\rho}(r_n) = \sum_{m \leq n} f_m(r_1, \ldots, r_n) \hat{\psi}_l(r_1) \ldots \hat{\psi}_l(r_m) \hat{\psi}(r_m) \ldots \hat{\psi}(r_1), \tag{35}
\]

and only then use substitutions \((30)-(31)\) to evaluate matrix elements:

\[
|\Psi_0\rangle e^{i\hat{H}t} \hat{\rho}(r_1) \ldots \hat{\rho}(r_n) e^{-i\hat{H}t} |\Psi_0\rangle = \sum_{m \leq n} f_m(r_1, \ldots, r_n) \langle \Psi_0 | e^{i\hat{H}t} \hat{\psi}_l(r_1) \ldots \hat{\psi}_l(r_m) e^{-i\hat{H}t} e^{i\hat{H}t} \hat{\psi}(r_m) \ldots \hat{\psi}(r_1) e^{-i\hat{H}t} |\Psi_0\rangle = \sum_{m \leq n} f_m(r_1, \ldots, r_n) \langle \Psi_0 | \left( a_1^\dagger u_1(r_1, t) e^{-iQ_1r_1} + a_2^\dagger u_2(r_1, t) e^{-iQ_2r_1} \right) \ldots \left( a_1^\dagger u_1(r_m, t) e^{-iQ_1r_m} + a_2^\dagger u_2(r_m, t) e^{-iQ_2r_m} \right) \times (a_1 u_1(r_1, t) e^{iQ_1r_1} + a_2 u_2(r_1, t) e^{iQ_2r_1}) \ldots (a_1 u_1(r_m, t) e^{iQ_1r_m} + a_2 u_2(r_m, t) e^{iQ_2r_m}) |\Psi_0\rangle. \tag{36}
\]
B. Interference of independent clouds

The surprising phenomenon which was observed in Ref. [27] is the appearance of interference fringes in the case when condensates are completely independent. To illustrate how interference fringes appear for independent clouds, let us now discuss the case when the numbers of particles in each of the clouds, $N_1$ and $N_2$, are fixed, hence the phase difference between the two clouds is not well defined. Initial state in the Fock basis in this case is given by (19):

$$|N_1, N_2\rangle = \frac{1}{\sqrt{N_1!N_2!}} (a_1^\dagger)^{N_1} (a_2^\dagger)^{N_2}|0\rangle. \tag{37}$$

Using the formalism of second quantization explained earlier, one can evaluate $\langle \rho^Q \rangle$ by analogy to Eq. (29):

$$\langle \rho^Q \rangle = \langle N_1, N_2|a_1^\dagger a_2|N_1, N_2\rangle = 0. \tag{38}$$

However, $\langle \rho^Q \rangle = 0$ doesn’t imply that there are no interference effects for independent condensates. Indeed, $\langle \rho^Q \rangle$ gives only the statistical average over many experiments, according to the usual interpretation of expectation values of operators in quantum mechanics. Being a quantum operator, $\hat{\rho}^Q$ has non vanishing quantum fluctuations. In each particular realization of experiment, complex number $\rho^Q$ can have a nonzero value. To show this, let us evaluate $\langle |\rho^Q|^2 \rangle$, which is the density-density correlation function at wave vector $Q$:

$$\langle |\rho^Q|^2 \rangle = \langle \hat{\rho}^Q \hat{\rho}^{Q\dagger} \rangle = \langle \int d\mathbf{r} d\mathbf{r}' \hat{\psi}^\dagger(r) \hat{\psi}(r) \hat{\psi}^\dagger(r') \hat{\psi}(r') e^{i\mathbf{Q}(r-r')} \rangle = \langle \int d\mathbf{r} d\mathbf{r}' \hat{\psi}^\dagger(r) \hat{\psi}(r) \hat{\psi}^\dagger(r') \hat{\psi}(r') e^{i\mathbf{Q}(r-r')} + \int d\mathbf{r} \hat{\psi}^\dagger(r) \hat{\psi}(r) \rangle. \tag{39}$$

These matrix elements can be evaluated using the second quantization prescription of the previous section, and the result is

$$\langle |\rho^Q|^2 \rangle = \langle N_1, N_2|a_1^\dagger a_2^\dagger a_1 a_2 + a_1^\dagger a_1 + a_2^\dagger a_2|N_1, N_2\rangle = N_1 N_2 + N_1 + N_2. \tag{40}$$

In the limit of large $N_1 = N_2 = N/2$, the leading contribution to $\langle |\rho^Q|^2 \rangle$ is the same as for the state with the fixed phase. Information about the full distribution of the quantum operator $\rho^Q$ is contained in higher moments of the distribution. If one considers higher moments of the type $\langle |\rho^Q|^{2n} \rangle = \langle \rho^{Qn} \rho^{Q\dagger n} \rangle$, the leading contribution in the limit of large $N_1$ and $N_2$ will again have the form

$$\langle |\rho^Q|^{2n} \rangle = (N_1 N_2)^n \left( 1 + O\left(\frac{1}{N_1}\right) + O\left(\frac{1}{N_2}\right) \right). \tag{41}$$

Corrections which appear because of the normal ordering result in subleading terms which are denoted by $O(1/N_1) + O(1/N_2)$. The leading term implies that in the limit of large $N_1$ and $N_2$, the distribution function of $|\rho^Q|^2$ is highly peaked near the value $N_1 N_2$, with the relative width which is proportional to the inverse square root of number of particles. Any operator of the form $\rho^{Qn} \rho^{Q\dagger n}$ will have zero expectation value for $m \neq n$ similar to $\rho^Q$, which means that the phase of the complex number $\rho^Q$ is uniformly distributed from 0 to $2\pi$. The expectation value of any operator which depends on the phase of $\rho^Q$ becomes zero due to the averaging over the phase.

The physical picture which emerges from the calculations is the following [61, 62, 64]: for two independent ideal clouds in the limit of large $N$ the absolute value of interference fringe amplitude is the same as for the state with a fixed relative phase, but the position of the intensity minima fluctuates from shot to shot. The state with a fixed number of particles is a superposition of states with fixed relative phases. For example,

$$|N/2, N/2\rangle = \left( \frac{\pi N}{2} \right)^{1/4} \int_0^{2\pi} d\varphi \frac{\varphi}{2\pi} |\varphi, N\rangle. \tag{42}$$

In the limit of large $N$ the phase states are almost orthogonal, and the measurement picks some value of the relative phase. Since the relative phase is not well defined for independent clouds, in each particular experiment the positions of the minima will fluctuate from shot to shot. To distinguish independent clouds from states which have correlated relative phases, one needs to do a series of experiments and measure not only the absolute magnitude of interference fringes, but also the positions of the minima. Experiments which distinguish states with a fixed relative phase from some other many body states are already being done, and can be used i.e. to measure the temperature [103] or to study the dynamical evolution [29, 30, 45, 48, 49] of the relative phase.
III. FULL COUNTING STATISTICS OF SHOT NOISE

As has been explained in the previous section, for experiments with independent clouds the average interference amplitude depends only on the number of particles per cloud. In this section we consider not only the average interference amplitude, but also its shot to shot fluctuations due to a finite number of atoms in the clouds. We will demonstrate that while the average value of $|\rho_Q|^2$ depends only on the number of particles per cloud, the full distribution function of the variable

$$R = |\rho_Q|^2$$

contains information about the states of individual clouds. Our analysis is motivated by the earlier work of Polkovnikov [99], who showed that the variance of the fringe amplitude decreases as the inverse power of the number of particles per cloud, $N$, with a non-universal coefficient which contains information about the state of the clouds. Experimental observation of the effects discussed in this section requires systems with a small number of atoms. This may be realized with micro-BEC’s on chips [29, 30, 45, 46, 47, 48, 49, 50].

The shot noise for finite $N$ has a fundamental nature, which stems from the probabilistic nature of quantum mechanics. Distributions of $R$ obtained below correspond to the following "idealized" experimental procedure: release the confining potential and take an absorption image of the columnar density on an ideal CCD camera with 100% efficiency (photon shot noise is ignored). To obtain the amplitude of interference fringes, $\rho_Q$, extract a Fourier component of the density at wave vector $Q$ from each image separately. The results of many experiments give the histogram $W(R)$ of the values of $R = |\rho_Q|^2$. We note that the quantum observable $\hat{R}$ defined in such way is a many-body operator, and calculation of its full distribution is a non-trivial task, even when all atoms are in the same state, such as for the case of a well defined relative phase between atoms in the two wells. In this section we develop a general method to find distribution functions of $R$ analytically.

We note that in our idealized setting we find interference patterns at a well defined wave vector $Q$. We expect that the finite size of the systems in transverse direction after expansion and collisions during the initial stage of expansion broaden the peak in the Fourier space to a finite, but small range of wave vectors around $Q$ [104]. Hence a one should consider $R = |\rho_Q|^2$ as an integral over the peak in the Fourier image of the density. Also in this paper we will discuss the amplitude of the interference fringes whereas experimental papers typically discuss visibility of the interference patterns. The two quantities differ only by the trivial rescaling.

Results for independent clouds in coherent (solid) and number (dashed) states for $N_1 = N_2 = 100$ are presented in Fig. 6. One can see, that even for a relatively large number of atoms, $N = 100$, fluctuations due to shot noise are appreciable. In Fig. 7 we compare the full counting statistics of $R = |\rho_Q|^2$ for independent clouds in coherent (dashed) and number (dotted) states with $N_1 = N_2 = 20$, and for clouds with a fixed relative phase (solid) with total number of atoms $N = N_1 + N_2 = 40$. Distribution functions for the cases of (i) well defined relative phase between the clouds and (ii) fixed number of atoms in each cloud are very close. They become indistinguishable when $R$ is rescaled by its average value $\langle R \rangle$, although each of them differs considerably from the Gaussian distribution. The distribution function of $R$ is wider for coherent states compared to number states, as was suggested in Ref. [99] based on the study of the variance of the two distributions. Hence the conclusion is that coherent and number states can be easily distinguished based on the statistics of fluctuations of $R$ relative to its average value.

In principle, when the two clouds are prepared with the same relative phase over many experiments, it is possible to distinguish independent condensates in number states from states with a fixed relative phase using a set of several interference experiments: minima positions are uniformly distributed for independent clouds, while for states with a fixed relative phase positions of the interference minima are always at the same points in space. However one can imagine the situation when the clouds are prepared in state with a fixed relative phase but the relative phase itself is random from realization to realization. Our results show that it is practically impossible to distinguish such states from number states by looking at the distribution of the amplitude of interference fringes. Our calculations below provide additional support to the physical interpretation [61, 62, 63, 64] presented in the previous section. Fluctuations of the absolute value of the interference amplitude are the same for the cases when clouds have random relative phase and when clouds are prepared in number states so the random relative phase is "measurement induced".

The method developed in this section can be generalized to a variety of experimental situations, i.e. several independent condensates. Different squeezed states within individual condensates can be considered, and measurement of full counting statistics of shot noise can be used as an experimental probe to distinguish between different correlated states.

As noted earlier, we will be interested in the full distribution function of the positive definite quantum observable

$$\hat{R} = |\hat{\rho}_Q|^2 = \hat{\rho}_Q \hat{\rho}_Q$$

defined by Eq. (39). To calculate its full distribution function, $W(R)$, one needs to know expressions for higher moments $\langle \hat{R}^n \rangle$. After that, one has to solve the "problem of moments", i.e. to recover the distribution function on
FIG. 6: Rescaled distribution functions of $R = |\rho_Q|^2$ for independent clouds in the coherent states (solid) or in the states with a well defined numbers of atoms (dashed). Here $N_1 = N_2 = 100$.

FIG. 7: Distribution functions of $R = |\rho_Q|^2$ for independent clouds in the coherent (dashed) and in the number (dotted) states with $N_1 = N_2 = 20$. Solid line is a distribution function of $R$ for clouds with a fixed relative phase with total number $N = N_1 + N_2 = 40$. Distribution functions for states with a fixed relative phase and with fixed numbers are very close, and become indistinguishable when $R$ is normalized by its average value $\langle R \rangle$.

the $(0, \infty)$ interval using all moments. In general, this procedure is numerically hard and unstable, unless higher moments have a certain analytical form. If the expression for higher moments is known analytically, then one can sometimes avoid the "problem of moments" by calculating the so-called characteristic function, $\chi(\lambda)$, which is the Laplace transform of $W(R)$:

$$\chi(\lambda) = \int_0^\infty e^{-\lambda R} W(R) dR = \int_0^\infty \sum_{i=0}^\infty \frac{(-\lambda R)^n}{n!} W(R) dR = \sum_{i=0}^\infty \frac{(-\lambda)^n \langle \hat{R}^n \rangle}{n!}$$

(45)

If $\chi(\lambda)$ can be calculated analytically, then $W(R)$ can be recovered by the inverse Laplace transform. In our case it is more practical to calculate not the characteristic function, but the analog of the Hankel transformation [105, 106] of $W(R)$, given by

$$Z(i\lambda) = \sum_{n=0}^{n=\infty} \frac{(i\lambda)^{2n}}{(n!)^2} \langle \hat{R}^n \rangle.$$  

(46)

Using the expansion of the zeroth order Bessel function, one can write

$$Z(i\lambda) = \int_0^\infty W(R) J_0(2\lambda \sqrt{R}) dR.$$  

(47)

(48)

The inversion of the transformation can be found using the orthogonality condition for the zeroth order Bessel functions...
By the end of this section we will provide analytical expressions for $Z(i\lambda)$ for certain cases (see Eqs. (62), (65), (72)), from which $W(R)$ can be obtained by simple numerical integration according to Eq. (49).

To proceed we note that $\hat{\rho}_{\mathbf{Q}}$ and $\hat{\rho}_{-\mathbf{Q}}$ commute with each other:

$$[\hat{\rho}_{\mathbf{Q}}, \hat{\rho}_{-\mathbf{Q}}] = 0.$$ (50)

Operators $\hat{\rho}_{\mathbf{Q}}$ are understood as in Eq. (27), without the projection on single particle states $\psi_1(\mathbf{r}, t), \psi_2(\mathbf{r}, t)$ as in Eqs. (40)-(41). The latter substitutions can be only done after normal ordering (see discussion in section II A 2). Hence we find

$$Z(i\lambda) = \langle \int_0^{2\pi} \frac{d\varphi}{2\pi} e^{i\lambda(\hat{\rho}_{\mathbf{Q}}e^{i\varphi} + \hat{\rho}_{-\mathbf{Q}}e^{-i\varphi})} \rangle.$$ (51)

Indeed, after expanding the exponent and integration, only even degrees of $i\lambda$ survive, and non vanishing terms are exactly what is needed for Eq. (10).

The normal ordering of Eq. (51) can be done using the following identity:

$$e^{\int f(r)\hat{\psi}^\dagger(r)\hat{\psi}(r) dr} = \langle e^{\int (f(r)-1)\hat{\psi}^\dagger(r)\hat{\psi}(r) dr} \rangle.$$ (52)

Here we have assumed that operators $\hat{\psi}^\dagger(r), \hat{\psi}(r)$, have the canonical commutation relations given by Eq. (17). Eq. (52) is a generalization of the simpler identity [107]:

$$e^{\lambda a^\dagger a} = \langle e^{(\lambda-1)a^\dagger a} \rangle.$$ (53)

for operators which obey the commutation relations

$$[a, a^\dagger] = 1.$$ (54)

The normal ordering signs $: :$ mean that all creation operators should be put to the left of annihilation operators in Taylor expansion of expressions being ordered. To illustrate the meaning of Eq. (53), let us consider the expansions of left and right side up to $\lambda^2$. The left hand side is

$$1 + \lambda a^\dagger a + \frac{\lambda^2}{2} a^\dagger aa^\dagger a + O(\lambda^3) = 1 + (\lambda + \frac{\lambda^2}{2})a^\dagger a + \frac{\lambda^2}{2} a^\dagger aa^\dagger a + O(\lambda^3).$$ (55)

The right hand side is

$$1 + (\lambda - 1) : a^\dagger a : + \frac{(\lambda - 1)^2}{2} : a^\dagger aa^\dagger a : + O(\lambda^3) =$$

$$1 + (\lambda + \frac{\lambda^2}{2} + O(\lambda^3)) : a^\dagger a : + \frac{\lambda^2}{2} + O(\lambda^3) : a^\dagger aa^\dagger a : + O(\lambda^3) =$$

$$1 + (\lambda + \frac{\lambda^2}{2})a^\dagger a + \frac{\lambda^2}{2} a^\dagger aa^\dagger a + O(\lambda^3).$$

Eq. (53) holds not only up to $\lambda^2$, but to all orders in $\lambda$ and plays an important role in quantum optics.

Using the definition of $\rho_{\mathbf{Q}}$ given by Eq. (27), one can apply Eq. (52) with $f(r) = 2i\lambda \cos(Qr + \varphi)$ and rewrite $Z(i\lambda)$ in Eq. (51) as

$$Z(i\lambda) = \langle \int_0^{2\pi} \frac{d\varphi}{2\pi} e^{\int 2i\lambda \cos(Qr + \varphi) - 1)\hat{\psi}^\dagger(r)\hat{\psi}(r) dr} \rangle.$$ (56)
A. Interference of two independent coherent condensates

Let us first explain how to evaluate $Z(i\lambda)$ for independent clouds in coherent states \[5\] of operators $a_1$ and $a_2$ with eigenvalues $\sqrt{N_1}e^{i\psi_1}$ and $\sqrt{N_2}e^{i\psi_2}$. Since coherent states form a complete basis, any initial state can be expanded in this basis, and thus the problem of calculating of $W(R)$ is essentially solved for arbitrary initial states. Coherent states are convenient, since they are the eigenstates of the annihilation operator, and the annihilation operator acts on them as a $c$-number. Hence after making substitutions \[30]-\[31\] into the normal ordered expression, one can substitute operators $a_1, a_1^\dagger$ by numbers $\sqrt{N_1}e^{i\psi_1}, \sqrt{N_2}e^{-i\psi_2}$. Since the normal ordered expression is obtained by the normal ordering of the Taylor expansion of Eq. \[56\], one needs to collect the Taylor series back. For coherent states, the whole procedure is equivalent to removing the normal ordering signs in Eq. \[56\], making substitutions \[30]-\[31\] and treating operators $a_1, a_1^\dagger$ as numbers $\sqrt{N_1}e^{i\psi_1}, \sqrt{N_2}e^{-i\psi_2}$. Thus we obtain

$$Z(i\lambda; \sqrt{N_1}e^{i\psi_1}, \sqrt{N_2}e^{i\psi_2}) = \int_0^{2\pi} \frac{d\phi}{2\pi} e^{\frac{i2\lambda}{2\pi} e^{i(\psi_1+\psi_2)}(N_1u_1(r,t)^2+N_2u_2(r,t)^2+2\sqrt{N_1N_2}\cos(\psi_1+\psi_2)e^{i\lambda}u_1(r,t)u_2(r,t))}d\phi.$$  \tag{57}

Similar to section II A 2 we assume the that normalized functions $u_1(r,t)$ and $u_2(r,t)$ strongly overlap and vary at scales much larger than $1/Q$, which is equivalent to

$$\int e^{i\lambda R}u_1(r,t)u_2(r,t)d\phi = \delta_{\lambda 0}.$$  \tag{58}

Then integration over $d\phi$ in the exponent of Eq. \[57\] can be done using the following equations:

$$\int \left( e^{2i\lambda \cos(Qr+\varphi)} - 1 \right) (N_1u_1(r,t)^2+N_2u_2(r,t)^2) d\phi = \sum_{m=1}^{\infty} \frac{(i\lambda)^{2m}}{(2m)!} \int (N_1u_1(r,t)^2+N_2u_1(r,t)^2)(e^{i(Qr+\varphi)} + e^{-i(Qr+\varphi)})^{2m} d\phi = \sum_{m=1}^{\infty} (N_1+N_2) \frac{(i\lambda)^{2m}}{(2m)!} \cdot \frac{m!m!}{m!m!} = (J_0(2\lambda) - 1)(N_1+N_2); \tag{59}$$

$$2\sqrt{N_1N_2} \int \left( e^{2i\lambda \cos(Qr+\varphi)} - 1 \right) \cos(\psi_1-\psi_2+Qr)u_1(r,t)u_2(r,t) d\phi = \sqrt{N_1N_2} \sum_{m=0}^{\infty} \frac{(i\lambda)^{2m+1}}{(2m+1)!} \int \left( e^{i(Qr+\varphi)} + e^{-i(Qr+\varphi)} \right)^{2m+1} \left( e^{i(\psi_1-\psi_2+Qr)} + e^{-i(\psi_1-\psi_2+Qr)} \right) u_1(r,t)u_2(r,t) d\phi = 2i\sqrt{N_1N_2}J_1(2\lambda) \cos(\psi_1-\psi_2-\varphi). \tag{60}$$

Substituting \[59] and \[60] into \[57\], and doing the integral over $\varphi$, we finally obtain the central result of this section:

$$Z(i\lambda; \sqrt{N_1}e^{i\psi_1}, \sqrt{N_2}e^{i\psi_2}) = \int_0^{2\pi} \frac{d\phi}{2\pi} e^{(J_0(2\lambda)-1)(N_1+N_2) + 2i\sqrt{N_1N_2}J_1(2\lambda) \cos(\psi_1-\psi_2-\varphi)} = e^{(J_0(2\lambda)-1)(N_1+N_2)J_0 \left(2\sqrt{N_1N_2}J_1(2\lambda)\right)}. \tag{61}$$

B. Interference of independent clouds in number states

Let us now explain how to calculate $Z'(i\lambda, N_1, N_2)$ for the Fock states with the number of particles equal to $N_1$ and $N_2$. First, we need to expand the Fock states $|N_1, N_2\rangle$ using the coherent states basis. Since the basis of coherent states is overcomplete \[108\], there are many ways to do a decomposition. For our purposes it is convenient to use

$$|N_1, N_2\rangle = (a_1^\dagger)^{N_1}(a_2^\dagger)^{N_2} |0\rangle = \sqrt{N_1N_2}e^{\alpha N_1-N_1e^2} \int_0^{2\pi} \int_0^{2\pi} d\varphi_1 R d\varphi_2 R (2\pi)^2 e^{-iN_1\varphi_1 R-iN_2\varphi_2 R} |\alpha e^{i\varphi_1 R}, \alpha e^{i\varphi_2 R}\rangle,$$  \tag{63}

where $\alpha$ is an arbitrary real positive number. Coherent states are given by

$$|\alpha e^{i\varphi_1 R}, \alpha e^{i\varphi_2 R}\rangle = e^{-\alpha^2 + \alpha e^{i\varphi_1 R}a_1^\dagger + \alpha e^{i\varphi_2 R}a_2^\dagger} |0\rangle,$$  \tag{64}

$$= e^{-\alpha^2 + \alpha e^{i\varphi_1 R}a_1^\dagger + \alpha e^{i\varphi_2 R}a_2^\dagger} |0\rangle.$$
and the overlap between them equals
\[ \langle \alpha e^{i\varphi_{1L}}, \alpha e^{i\varphi_{2L}} | \alpha e^{i\varphi_{1R}}, \alpha e^{i\varphi_{2R}} \rangle = e^{\alpha^2(e^{i\varphi_{1R} - i\varphi_{1L}} + e^{i\varphi_{2R} - i\varphi_{2L}}) - 2}. \] (65)

One can also expand the bra state \( \langle N_1, N_2 \rangle \) similarly to (63) by introducing integration variables \( \varphi_{1L}, \varphi_{2L} \). For any given values of \( \varphi, \varphi_{1L}, \varphi_{2L}, \varphi_{1R}, \) and \( \varphi_{2R} \), matrix elements between coherent states can be evaluated as in previous section, and by simple modification of Eq. (61) we obtain
\[ Z^f(i\lambda, N_1, N_2) = \int \int \int d\varphi_{1L} d\varphi_{2L} d\varphi_{1R} d\varphi_{2R} \frac{(2\pi)^5}{N_1! N_2!} \alpha^{-2N_1 - 2N_2} e^{iN_1(\varphi_{1L} - \varphi_{1R}) + iN_2(\varphi_{2L} - \varphi_{2R})} \times \langle \alpha e^{i\varphi_{1L}}, \alpha e^{i\varphi_{2L}} | \alpha e^{i\varphi_{1R}}, \alpha e^{i\varphi_{2R}} \rangle \times e^{(J_0(2\lambda) - 1)\alpha^2(e^{i\varphi_{1L} - i\varphi_{1L}} + e^{i\varphi_{2R} - i\varphi_{2L}})} e^{i\alpha^2 J_1(2\lambda)(e^{i(\varphi_{1L} + \varphi_{2L})} - e^{-i(\varphi_{1L} + \varphi_{2L})})}, \] (66)

One can now substitute Eq. (65) into the equation above, introduce variables \( \varphi_1 = \varphi_{1L} - \varphi_{1R} \) and \( \varphi_2 = \varphi_{2L} - \varphi_{2R} \), and integrate over \( \varphi, \varphi_{1R}, \varphi_{2R} \). Multiple cancellations occur, and eventually we obtain
\[ Z^f(i\lambda, N_1, N_2) = \int \frac{d\varphi_1 d\varphi_2}{(2\pi)^2} N_1! N_2! \alpha^{-2N_1 - 2N_2} e^{iN_1 \varphi_1 + iN_2 \varphi_2} e^{J_0(2\lambda)\alpha^2(e^{i\varphi_1} + e^{-i\varphi_2})} J_0(2\alpha^2 J_1(2\lambda)e^{-i(\varphi_1 + \varphi_2)/2}). \] (68)

Both integrations in the equation above can be done in a closed form for arbitrary positive integer \( N_1 \) and \( N_2 \) using hypergeometric functions. Here we will present the results only for \( N_1 = N_2 = N \). One needs to expand the last exponent and \( J_0(2\alpha^2 J_1(2\lambda)e^{-i(\varphi_1 + \varphi_2)/2}) \) using Taylor series. After integration over \( d\varphi_1 \) and \( d\varphi_2 \) dependence on \( \alpha \) disappears, and we obtain the final result for Fock states:
\[ Z^f(i\lambda, N, N) = \sum_{k=0}^{N} (N!)^2 J_0(2\lambda) \frac{2k}{(N - k)!} (iJ_1(2\lambda))^{2(N-k)} = 2 F_1 \left( \frac{-n}{1}, \frac{-J_0(2\lambda)}{J_1(2\lambda)} \right) (iJ_1(2\lambda))^{2N}, \] (69)

where \( 2 F_1 (a, b; c; x) \) in a hypergeometric function defined by
\[ 2 F_1 (a, b; c; x) = 1 + \frac{ab x}{c!} + \frac{a(a+1)b(b+1)x^2}{c(c+1)2!} + ... \] (70)

C. Clouds with a well defined relative phase

Let us now consider the case of clouds with a fixed relative phase, when the initial state \( |\varphi_0, N \rangle \) at \( t = 0 \) is given by Eq. (13). This state can be expanded using the coherent states basis as
\[ |\varphi_0, N \rangle = \frac{1}{(2\pi)^{1/2}} \langle 0 | e^{i\varphi_0/2} + e^{-i\varphi_0/2} \rangle^N |0 \rangle = \sqrt{N!} (\sqrt{2\alpha})^{-N} e^{\alpha^2} \int_0^{2\pi} \frac{d\varphi_R}{2\pi} e^{-iN\varphi_R} \langle \alpha e^{i\varphi_L + i\varphi_0/2}, \alpha e^{i\varphi_R - i\varphi_0/2} \rangle. \]

A similar expansion can be written for bra- vector \( \langle \varphi_0, N \mid \) using the phase variable \( \varphi_L \). Coherent states and their overlaps are given by Eqs. (63), (64), and one obtains an expression for the generating function \( Z(i\lambda, N) \) similar to Eq. (66):
\[ Z(i\lambda, N) = \int \int \int \frac{d\varphi_{1L} d\varphi_{2L}}{(2\pi)^3} N! (\sqrt{2\alpha})^{-2N} e^{2\alpha^2} \langle \alpha e^{i\varphi_L + i\varphi_0/2}, \alpha e^{i\varphi_R - i\varphi_0/2} \rangle \times \times e^{2(J_0(2\lambda) - 1)\alpha^2(e^{i\varphi_R - i\varphi_L})} e^{i\alpha^2 J_1(2\lambda)(e^{i(\varphi_0 + \varphi_R - \varphi_L)} + e^{-i(\varphi_0 + \varphi_R - \varphi_L)})}. \]

The integrand depends only on the difference \( \Delta \varphi = \varphi_L - \varphi_R \), and the integral over \( \varphi \) can be done analytically. Dependence of \( Z(i\lambda, N) \) on \( \varphi_0 \) drops out, as expected:
\[ Z(i\lambda, N) = \int \int \frac{d\Delta \varphi}{2\pi} N! \alpha^{-2N} e^{iN\Delta \varphi} e^{2J_0(2\lambda)\alpha^2 e^{-i\Delta \varphi}} J_0(2\alpha^2 J_1(2\lambda)e^{-i\Delta \varphi}). \] (71)

Expanding the last exponent and \( J_0(2\alpha^2 J_1(2\lambda)e^{-i\Delta \varphi}) \) in the expression above and integrating over \( \Delta \varphi \), we obtain the final expression for even \( N \):
\[ Z(i\lambda, N) = N! 2^{-N} \sum_{k=0}^{N/2} \frac{(iJ_1(2\lambda))^{2k} (2J_0(2\lambda))^{N-2k}}{(k!)^2} \frac{N!}{(N - 2k)!} = 2^{-N} (iJ_1(2\lambda))^N C_N^N \left( \frac{J_0(2\lambda)}{iJ_1(2\lambda)} \right), \] (72)

where \( C_n^a(x) \) is a Gegenbauer polynomial (106).
IV. INTERFERENCE OF FLUCTUATING LOW-DIMENSIONAL GASES

As we discussed in previous sections, for macroscopic three dimensional Bose-Einstein Condensates the long-range phase coherence manifests itself in the nearly perfect interference fringes between two independent condensates [27]. For low dimensional Bose gases, the situation is different, since phase fluctuations are very effective in destroying the long-range order. In one dimension, long-range coherence is prohibited even at zero temperature [92], while in two dimensions any nonzero temperature destroys long-range order [91]. In addition, the Berezinskii-Kosterlitz-Thouless (BKT) phase transition occurs [85, 86], which separates the low temperature phase with power-law correlations from the high temperature phase with short-range correlations. Phase fluctuations reduce the average visibility of the interference fringes, and result in the shot to shot fluctuations of the visibility.

In this section, we discuss how measurements of interference fringes can reveal information about spatial correlations within individual condensates. The typical experimental setups are shown in Figs. 4 and 8. They correspond to the so-called open boundary conditions (OBC). Essentially the OBC mean that the imaged area is cut out of a larger system. As a theoretical model one can also consider a one-dimensional condensate with periodic boundary conditions (PBC), which corresponds to interference experiments with two coaxial rings lying in two parallel $xy$ planes. While this model is somewhat artificial from the experimental point of view (see however Ref. [109]), it allows a very elegant theoretical analysis, hence we will discuss it in these lecture notes as well.

The confining potential is highly anisotropic, and after it is switched off, the clouds predominantly expand in the transverse direction, while no significant expansion occurs in the axial (for 1D gases) or in-plane (for 2D gases) directions. For low-dimensional gases the phase of the condensate doesn’t have a long-range order due to quantum or thermal fluctuations. Locally the phase determines the positions of the minima of the absorption intensity, and fluctuations of the phase lead to fluctuations of the interference fringe positions along the condensates, as shown in Fig. 4. Fluctuations of the fringe positions contain information about the original phase fluctuations present in the system, which are preserved during expansion.

To extract information about fringe position fluctuations for the 1D case, we will integrate the intensity along the axes of the clouds. Fluctuations of the relative phase result in fluctuations of the minima positions for different $x$. For each $y$, the image can be integrated along the $x$ direction to give the integrated fringe amplitude $\rho_Q(L)$ (see Fig. 3). Note that the integrated fringe amplitude depends on the integration length $L$. One experimental image can be
used to extract information for different values of $L$. Many images are still required to obtain distribution functions for each $L$. For 2D gases, the setup is analogous and is shown in Fig. 3. Here part of the integration is performed by the imaging beam itself. The size of the integration area along the direction of the imaging beam can be controlled by applying magnetic field gradients, so that only a specified section of the cloud is resonant with the probe light.

The operator which corresponds to the fringe amplitude $\rho Q(L)$, illustrated in Fig. 3, is the same as $\rho Q$ defined by Eq. (27), where the integration along $x$-dimension is limited to the section of length $L$. Let us first consider the expectation value $\langle |\rho Q(L)|^2 \rangle$ (expectation values of operators which depend on the phase of $\rho Q(L)$ vanish, similar to 3D case, since two clouds are assumed to be independent). One has to use modified formulas (30)-(31), where operators $a_1^\dagger, a_i$ are now allowed to have $x$-dependence. In the limit when the number of particles in the section of size $L$ is large, the average value of $\langle |\rho Q(L)|^2 \rangle$ is given by

$$\langle |\rho Q(L)|^2 \rangle = \int_0^L \int_0^L dx_1 dx_1' \langle a_1^\dagger(x_1) a_1^\dagger(x_1') a_1(x_1') a_1(x_1) \rangle.$$  \hspace{1cm} (73)

Note that in Eq. (73) we used the normal ordered form of the operators which means that we neglect the shot noise considered in section II. This is justified for long condensates as we discuss below. From now on, we will concentrate on the case when independent clouds are identical and have the same density of particles with equal interaction strengths. Then

$$\langle |\rho Q(L)|^2 \rangle = \int_0^L \int_0^L dx_1 dx_1' \langle a_1^\dagger(x_1) a(x_1') \rangle^2.$$ \hspace{1cm} (74)

To gain intuition into the physical meaning of the average amplitude of interference fringes, we address two limiting cases. First, consider the situation when $\langle a_1^\dagger(x_1) a(x_0) \rangle$ decays exponentially with distance and the correlation length is given by $\xi << L$. Then Eq. (74) implies $\langle |\rho Q(L)| \rangle \propto \sqrt{L/\xi}$, which has a simple physical interpretation. Since the phase is only coherent over a length $\xi$, the system is effectively equivalent to a series of $L/\xi$ pairs of independent condensates. Each pair contributes interference fringes with a constant amplitude proportional to $\xi$ and a random phase. The total amplitude $\rho Q(L)$ is the result of adding $L/\xi$ independent vectors of constant length $\xi$ and random direction. Adding random uncorrelated vectors gives a zero average except for a typical square root fluctuation. Thus scaling of the absolute value of the net interference amplitude is $\sqrt{L/\xi}$. This observation is similar in spirit to that made in Ref. 55 of interference between 30 independent condensates in a chain. Fringes can be seen, though their average amplitude is suppressed by a factor of $\sqrt{30}$ compared to the interference between two condensates. Now consider the opposite limit of perfect condensates, for which $\langle a_1^\dagger(x_1) a(x_0) \rangle$ is constant. In this case Eq. (74) implies that $\langle |\rho Q(L)| \rangle \propto L$. Pictorially this is the result of adding vectors which are all aligned, resulting in a fringe amplitude absolute value of which scales as the total size of the system.

Methods developed in this section for analyzing $\langle |\rho Q(L)|^2 \rangle$ can be applied to condensates with either uniform and non uniform densities. For simplicity, we concentrate on the case when $L$ is much smaller than the size of the clouds, so the change in the atomic density along the clouds can be ignored. In this case correlation functions for 1D gases are described by the Luttinger liquid theory 110, 111. For OBC at zero temperature two-point correlation functions are given by

$$\langle a_1^\dagger(x) a(y) \rangle \sim \rho \left( \xi_h / |x - y| \right)^{1/2K}.$$ \hspace{1cm} (75)

Here $\rho$ is the particle density, $\xi_h$ is the healing length, which also serves as the short range cutoff, and $K$ is the so-called Luttinger parameter, which characterizes the strength of interactions. For bosons with a repulsive short-range potential, $K$ ranges between 1 and $\infty$, with $K = 1$ corresponding to strong interactions, or “impenetrable” bosons, while $K \rightarrow \infty$ for weakly interacting bosons. Substituting Eq. (75) into Eq. (74) and assuming that $L \gg \xi_h$, we obtain

$$\langle |\rho Q(L)|^2 \rangle = \tilde{C} \rho^2 L^2 \left( \frac{\xi_h}{L} \right)^{1/K},$$ \hspace{1cm} (76)

where $\tilde{C}$ is a constant of order unity. We see that the amplitude of the interference fringes $\langle |\rho Q(L)| \rangle$ scales as a non trivial power of the imaging length. In the non interacting limit ($K \rightarrow \infty$) the scaling is linear $\langle |\rho Q(L)| \rangle \sim L$, as expected for a fully coherent system. Interestingly, $\langle |\rho Q(L)| \rangle \sim \sqrt{L}$ appears in the hard core limit ($K = 1$), as in systems with short range correlations which were discussed above.

One may be concerned that Eq. (76) gives only the long distance asymptotic behavior of the correlation functions, and does not describe the short distance behavior. From Eq. (74) one finds that the contribution of the short
distance part of the correlation functions to $\langle |\rho_\mathcal{Q}(L)|^2 \rangle$ scales as $L$. In the physically relevant case of $K > 1$ and in the limit of large $L$ this contribution is smaller than Eq. (76). We note that in principle one can use the exact Bethe ansatz solution of the Lieb-Liniger model to obtain correlation functions valid at all distances. Another contribution which has been neglected is the shot noise. The shot noise contribution to $\langle |\rho_\mathcal{Q}(L)|^2 \rangle$ comes from the normal ordering of operators, and equals

$$\int_0^L \int_0^L dx_1 dx'_1 \langle a_1^{\dagger}(x_1)a_2(x_1)a_2(x'_1)a_1(x'_1) \rangle - \int_0^L \int_0^L dx_1 dx'_1 \langle a_1^{\dagger}(x_1)a_2(x'_1)a_1(x'_1)a_2(x_1) \rangle = \int_0^L \int_0^L dx_1 dx'_1 \delta(x_1 - x'_1)\langle a_1^{\dagger}(x_1)a_1(x'_1) \rangle = n_{1D}L. \quad (77)$$

In the limit of large $L$ and $K > 1$ this is again a subleading contribution and can be neglected.

For 2D, one can use similar approach to describe the contrast distribution at finite temperature below the BKT transition. We note that we assume that the temperature is small enough such that 2D gas is in a quasicondensate regime, when only phase fluctuations are present. In this case, correlation functions are given by

$$\langle \rho^{(2)}(\mathcal{Q}) \rangle = \left( \frac{\xi_h}{\eta} \right)^{\eta(T)}, \quad (78)$$

where $\eta(T) = mT/(2\pi\hbar^2\rho_s(T))$ depends on the temperature and the superfluid density $\rho_s(T)$. The BKT transition takes place at the universal value $\eta_c(T_c) = 1/4$. To keep connection to the 1D case, we introduce

$$K = 1/(2\eta(T)), \quad (79)$$

and restrict our attention to $K > K_c = 2$. For temperatures above the BKT transition Eq. (78) doesn’t hold, and correlations decay exponentially. This means that the integrated interference amplitude will only increase as the square root of the integration area.

Fig. 8 illustrates experiments performed with 2D gases to identify the BKT transition. Two independent 2D condensates are confined in transverse directions using an optical lattice potential. After the optical potential is switched off and clouds expand, the density is imaged on a CCD camera. When temperature is small, interference of atom density to obtain an interference pattern is a classical measurement on a quantum mechanical wave function. The process of the measurement itself introduces an intrinsic quantum mechanical noise. Said differently, from shot to shot we will not get precisely the same value of the amplitude of interference fringes. Expressions for $\langle |\rho_\mathcal{Q}(L)|^2 \rangle$ which we derived in Eqs. (73)-(76) correspond to averaging over many shots. For example, data points

A. Interference amplitudes: from high moments to full distribution functions

Measuring atom density to obtain an interference pattern is a classical measurement on a quantum mechanical wave function. The process of the measurement itself introduces an intrinsic quantum mechanical noise. Said differently, from shot to shot we will not get precisely the same value of the amplitude of interference fringes. Expressions for $\langle |\rho_\mathcal{Q}(L)|^2 \rangle$ which we derived in Eqs. (73)-(76) correspond to averaging over many shots. For example, data points...
FIG. 9: Emergence of quasi-long-range order in a 2D gas. a, Examples of average integrated interference contrasts $\langle \tilde{C}^2 \rangle$ are shown for low (blue circles) and high (red squares) temperatures; $L_x$ is the integration length along $x$-direction. The lines are fits to the data using the power-law functions $1/(L_x)^{2\alpha}$. b, Exponent $\alpha$ as a function of the central contrast $c_0$. Central contrast $c_0$ serves as a "thermometer", such that smaller values of $c_0$ correspond to higher temperatures. Dashed lines indicate theoretically expected values of $\alpha$ above and below the BKT transition in a uniform system. Taken from Ref. [32].

from Fig. 9, they correspond to averaging over approximately one hundred measurements [33]. However each individual shot will give the value of $|\rho_Q(L)|^2$ which may be different from its average value.

An interesting question to consider is how this amplitude fluctuates from one experimental run to another. To address this question we need to consider higher moments of the operator $|\rho_Q(L)|^2$. Generalizing the argument which lead to Eq. (74), we obtain

$$\langle |\rho_Q(L)|^{2n} \rangle = \int_0^L \ldots \int_0^L dx_1 \ldots dx_n dx'_1 \ldots dx'_n |a^\dagger(x_1) \ldots a^\dagger(x_n)a(x'_1) \ldots a(x'_n)|^2.$$  (80)

In Eq. (80) we used a normal ordered correlation function similar to Eq. (74). One can calculate corrections due to normal ordering for higher moments of $|\rho_Q(L)|^2$, and show that in the limit of large $L$ and $K > 1$ they can be neglected.

From Eq. (74) we observe that $\langle |\rho_Q(L)|^2 \rangle$ contains information about two-point correlation functions of individual clouds. Eq. (80) shows that higher moments of $|\rho_Q(L)|^2$ contain information about higher order correlation functions. The full distribution of the fluctuating variable $|\rho_Q(L)|^2$ contains information about all high order correlation functions.

In the Luttinger liquid theory fluctuations of the phase are described by the Gaussian action. For Gaussian actions higher order correlation functions are simply related to two-point correlation functions (see e.g. Ref. [116]):

$$\langle a^\dagger(x_1) \ldots a^\dagger(x_n)a(x'_1) \ldots a(x'_n) \rangle = \prod_{i=1}^n \langle a^\dagger(x_i) a(x'_j) \rangle / \prod_{i<j} \langle a^\dagger(x_i) a(x_j) \rangle \prod_{i<j} \langle a^\dagger(x'_i) a(x'_j) \rangle.$$  (81)

Using this formula together with Eq. (75), the higher moments of fringe amplitudes can be written as

$$\langle |\rho_Q(L)|^{2n} \rangle = A_0^{2n} Z_{2n}, \text{ where } A_0 = \sqrt{C \rho^2 \xi_{th}^{1/K} L^{2-1/K}}.$$  (82)
$C$ is a constant of the order of unity and for OBC in 1D
\[
Z_{2n} = \int_0^1 \cdots \int_0^1 du_1 \cdots dv_n \left| \prod_{i<j} |u_i - u_j| \prod_{i<j} |v_i - v_j| \right|^\frac{n}{2} = \int_0^1 \cdots \int_0^1 du_1 \cdots dv_n e^{\frac{n}{2} \left( \sum_{i<j} G(u_i, u_j) + \sum_{i<j} G(v_i, v_j) - \sum_{i<j} G(u_i, v_j) \right)}.
\]

Here for OBC $G(x, y)$ is given by
\[
G(x, y) = \log |x - y|.
\]

Integrals similar to Eq. (83) appeared in the literature before [120, 121], but they are not easy to compute.

The Gaussian model possesses a powerful conformal symmetry [111, 116], which dictates the form of the correlation functions for periodic boundary conditions or nonzero temperatures. For PBC with circumference of the condensates equal to the imaging length, the change in the correlation functions leads to [111]
\[
G_{\text{per}}(x, y) = \frac{1}{\pi} \sin \pi |x - y|.
\]

For nonzero temperature, $Z_{2n}$ depends on $K$ and the thermal length $\xi_T = h v_s/(k_B T)$, where $v_s$ is the sound velocity:
\[
G(x, y, \xi_T) = \log \left( \frac{\xi_T}{\pi L} \sinh \frac{\pi |x - y| L}{\xi_T} \right).
\]

The analysis in this section (except for section IV C) doesn’t depend on the particular form of $G(x, y)$. The only general restriction is
\[
G(x, y) = G(y, x).
\]

Eq. (83) is also valid for 2D case below the BKT transition temperature, with $u_i$ and $v_i$ being 2D variables on a rectangle with $G(\vec{x}, \vec{y}) = \log |\vec{x} - \vec{y}|$, and with a properly redefined $A_0$.

In what follows, we will be interested in distribution functions $W(\alpha)$ of the variable $\alpha = |\rho_Q(L)|^2/A_0^2$ and of its normalized version $\tilde{\alpha} = |\rho_Q(L)|^2/(\rho_Q(L)^2)$. By the definition of higher moments
\[
Z_{2n} = \int_0^\infty W(\alpha) \alpha^n d\alpha, \quad \text{and} \quad \frac{Z_{2n}}{Z_2^n} = \int_0^\infty \tilde{W}(\tilde{\alpha}) \tilde{\alpha}^n d\tilde{\alpha}.
\]

The general problem we consider now is how to construct distribution functions $W(\alpha)$ given by Eq. (83) with $Z_{2n}$ given by Eq. (83), $G(x, y)$ can be an arbitrary symmetric function of $x$ and $y$, and not necessarily the function of their difference. This allows us to study systems with non-uniform density in external traps, where one needs to use modified correlation functions [111]. In section IV B we will show that $W(\alpha)$ is connected to the partition functions of various Sine-Gordon models and Coulomb gases, and methods developed in this section provide a new non perturbative tool for calculating partition functions of such models.

One can think of $Z_{2n}$ in Eq. (83) as partition functions of a classical two-component gas of fictitious charged "particles" in a microcanonical ensemble, with $K$ being the "temperature". We first briefly comment on the limiting cases $K \gg 1$, and $K \to 1$. For $K \gg 1$, expansion of $Z_{2n}$ gives $Z_{2n} \approx 1$ in the zeroth order. This means that $\tilde{W}(\tilde{\alpha}) \approx \delta(\tilde{\alpha} - 1)$, i.e. a very narrow function peaked near its average value. Higher order terms in the expansion give $\tilde{W}(\tilde{\alpha})$ a small width of the order of $1/K$, and are studied in detail in Appendix A. For 1D gases at zero temperature with $K \to 1$ or nonzero temperatures with $\xi_T K/L \ll 1$ the distribution function $\tilde{W}(\tilde{\alpha})$ is Poissonian irrespective of OBC or PBC: $\tilde{W}(\tilde{\alpha}) = e^{-\tilde{\alpha}}$. To demonstrate this we need to verify that $Z_{2n}/Z_2^n = \int_0^\infty \tilde{\alpha}^n e^{-\tilde{\alpha}} d\tilde{\alpha} = n!$. This can be shown using the classical gas analogy: as $K \to 1$, $Z_2 = \int |x - y|^{1/K} dxdy$ starts to diverge for $x \to y$, and the main contribution to $Z_{2n}$ comes from "molecular" states of the two-component gas, i.e. from the parts of the configuration space in which each "particle" has a "particle" of the opposite charge in its neighborhood: $n!$ is just the number of ways to form such pairs. In this language $K \gg 1$ corresponds to the "plasma" phase of the classical charged gas, and the evolution of the distribution function can be understood as a formation of "molecules" out of the "plasma" phase as the "temperature" (i.e. the Luttinger parameter $K$) is lowered. For finite $\xi_T$, the main contribution to $Z_{2n}$ comes from distances $\sim \xi_T K/L$, and if this parameter is much smaller than 1, then again "molecular" contributions dominate.

There is a simple physical interpretation to $W(\alpha)$ for both $K \gg 1$ and $K \to 1$. When bosons do not interact and $K = \infty$, there should be no phase fluctuations within individual condensates. Hence in each experiment we should
find a perfect interference pattern although positions of the density minima are unpredictable (see discussion in section 11). Alternatively in the regime of strong interactions when $K \to 1$ we can think of the net interference as a result of adding many random uncorrelated two dimensional vectors (see discussion after Eq. (73)). Earlier we used the fact that for a random walk the net displacement is proportional to the square root of the number of steps. But we also know that for 2D random walks the distribution function of the square of the net displacement is Poissonian, which is what we find for $\tilde{W}(\tilde{\alpha})$.

As $K$ varies from 1 to $\infty$, the distribution function $\tilde{W}(\tilde{\alpha})$ should evolve from being a very broad Poissonian function to a narrow delta function. The evolution of the distribution in the intermediate regimes will be studied in detail below. In Appendix A we develop a systematic expansion of $Z_{2n}$ in powers of $1/K$ for large $K$ which works for different dimensions and boundary conditions, and investigate the distribution functions in this limit. The rest of this section is organized as follows. In section IV D we discuss the connection of our problem to the Sine-Gordon models, which describe various physical problems, ranging from flux line in superconductors \cite{122} to string theory \cite{123}. In section IV C we obtain full distributions for the 1D case with PBC using the exact solution of the boundary Sine-Gordon model on a circle \cite{56}. In section IV D we present a novel non-perturbative solution which is applicable for any value of $K$ and for various dimensions and boundary conditions. We discuss the connection of the distribution functions of fringe visibilities to the statistics of random surfaces \cite{58}, and prove that for 1D case with periodic boundary conditions in the limit of large $K$ the distribution of fringe visibilities is given by the Gumbel distribution, one of the extreme value statistical distributions \cite{124}.

B. Connection of the fringe visibility distribution functions to the partition functions of Sine-Gordon models

To illustrate the connection of the fringe visibility distribution functions to the partition functions of Sine-Gordon models we start from a 1D system with periodic boundary conditions. Shortly we will demonstrate that this example provides a unique possibility to make an excursion into very different subjects of physics and mathematics. As shown earlier, for a one-dimensional ring condensate with PBC and circumference equal to the imaging length, the correlation functions lead to

$$G^{per}(x, y) = \log\left(\frac{1}{\pi} \sin(\pi(x - y))\right).$$

(89)

This function describes a two-component gas of particles interacting via a 1D Coulomb potential, which is a logarithmic in the interparticle separation. For PBC the distance is a chord function $\sin(\pi|x - y|)/\pi$. It is common to define the grand partition function of such Coulomb gas as

$$Z_2(K, g) = \sum_{n=0}^{\infty} \frac{g^{2n}}{(n!)^2} Z_2^{per}.$$

(90)

where $Z_2^{per}$ is given by Eq. (83) with $G(x, y)$ substituted by $G^{per}(x, y)$.

Several different physical problems can be related to the partition function given by Eq. (90). Here we mention only a few examples: a) the anisotropic Kondo model \cite{125}; b) the quantum impurity model in one dimensional interacting electron systems introduced in Ref. 104, which has been extensively studied in the context of edge states in Quantum Hall systems \cite{126}; c) the background-independent string theory and the model of strings attached to a D-brane originally introduced in Ref. \cite{123} (see Ref. \cite{126} for a recent review); d) Calogero-Sutherland model \cite{128}, which has numerous application as an effective model; e) flux line pinning in superconductors \cite{122}; f) quantum tunneling in the presence of dissipation within Caldeira-Leggett model \cite{129}; g) interference of two one-dimensional condensates \cite{55, 56, 58}.

We now briefly describe how the expression given by Eq. (90) appears as a partition function in physical systems. We consider the imaginary time action

$$S_{per}[\beta] = \frac{1}{2} \int_{-\infty}^{\infty} dx \int_{0}^{1} d\tau \left( \partial_{\tau} \phi \right)^2 + \left( \partial_x \phi \right)^2 + 2g \int_{0}^{1} d\tau \cos[\beta \phi(x = 0, \tau) + 2\pi pr]$$

(91)

known as the boundary Sine-Gordon model \cite{130}. The quantum field $\phi(x, \tau)$ in Eq. (91) is defined on an infinite line in $x$ direction and is assumed to be periodic along $\tau$ direction: $\phi(x, \tau) = \phi(x, \tau + 1)$. The interaction term is present only at $x = 0$. A typical physical system which is described by action (91) is an interacting 1D electron liquid scattered by an impurity \cite{106}. In this case the cos-term describes backscattering of electrons within the Luttinger liquid formalism. In Eq. (91) we also added a $p$-dependent phase winding term. For a quantum impurity problem this is somewhat reminiscent of having a finite voltage \cite{131}, while in interference experiments this term corresponds...
Partition function of Eq. (91) is defined as \[ Z = \int \mathcal{D}\phi e^{-S_{\text{pert}}[\phi]}. \] Here the microcanonical partition functions \( Z_{2n}^{(p)} \) are represented by the following integrals

\[
Z_{2n}^{(p)} = \int_0^{2\pi} \cdots \int_0^{2\pi} \prod_{i=1}^n \frac{du_i}{2\pi} \frac{dv_i}{2\pi} e^{i2\pi \sum (u_i - v_i)} \left| \prod_{l<j} 2\sin \left( \frac{u_l - u_j}{2} \right) \prod_{l<k} 2\sin \left( \frac{v_l - v_k}{2} \right) \right|^p. 
\]

\( Z_{2n}^{(p)} \) for \( p = 0 \) is related to \( Z_{2n}^{\text{per}} \) defined above by

\[
Z_{2n}^{(0)} = Z_{2n}^{\text{per}} (2\pi)^{-n/K}. 
\]

Then

\[
Z_0(K, g) = \sum_{n=0}^{\infty} \frac{g^{2n}}{(n!)^2} Z_{2n}^{\text{per}} = \sum_{n=0}^{\infty} \frac{g^{2n}}{(n!)^2} \langle \alpha^n \rangle = \sum_{n=0}^{\infty} \frac{g^{2n}}{(n!)^2} \int_0^{\infty} W(\alpha) \alpha^n d\alpha = \int_0^{\infty} W(\alpha) J_0(2g\sqrt{\alpha}) d\alpha. 
\]

\( Z_0(K, g) \) is essentially a Hankel transformation of \( W(\alpha) \), and inverting Eq. (71) we can express the probability \( W(\alpha) \) through the partition function \( Z_0(K, g) \). Noting that \( I_0(ix) = J_0(x) \) and using the completeness relation for Bessel functions, \( \int_0^{\infty} J_0(\lambda x) J_0(\lambda y) |x| |y| d\lambda = \delta(|x| - |y|) \), we obtain

\[
W(\alpha) = 2 \int_0^{\infty} Z_0(K, ig) J_0(2g\sqrt{\alpha}) g dg. 
\]

It is important that the last equation has the partition function at imaginary value of the coupling constant. This should be understood as an analytic continuation of \( Z_0(K, g) \). For PBC in 1D, partition functions \( Z_p(K, g) \) can be evaluated using the exact solution of the boundary Sine-Gordon model with periodic boundary conditions, and this approach will be presented in section IV.C.

For open boundary conditions one can also write the grand canonical partition function

\[
Z(g) = \sum_{n=0}^{\infty} \frac{g^{2n}}{(n!)^2} Z_{2n}, 
\]

where \( Z_{2n} \) are given by Eqs. (83), (84). We can express Eq. (99) as a partition function of a certain Sine-Gordon model:

\[
Z(g) = \frac{\int \mathcal{D}\phi e^{-S(g)}}{\int \mathcal{D}\phi e^{-S(0)}}. 
\]

where

\[
S(g) = \pi K \int_{-\infty}^{\infty} dx \int_{-\infty}^{\infty} d\tau [ (\partial_x \phi)^2 + (\partial_x \phi)^2 ] + 2g \int_0^1 d\tau \cos[2\pi \phi(x = 0, \tau)]. 
\]
Note different limits of the $\tau$ integration in the first and the second terms. One can see that Eqs. (99) and (100) define the same $Z(g)$ by observing that $-\log((x - y)/2\pi)$ is a free propagator of the Gaussian action on a plane.

So far we established the relation between the distribution functions of the amplitude of interference fringes of 1D condensates and the boundary Sine-Gordon models (91), (101). Generalization of this argument to the case of 2D condensates is straightforward. Higher moments of the interference amplitude are given by integrals of type (83), but each $u$ and $v$ is now a two-dimensional coordinate. Distribution function of fringe amplitudes is then related to the partition function of the bulk Sine-Gordon model

$$S_{2D}(g) = \pi K \int_{-\infty}^{\infty} dx \int_{-\infty}^{\infty} d\tau ((\partial_x \phi)^2 + (\partial_\tau \phi)^2) + 2g \int_{0}^{1} d\tau dx \cos[2\pi \phi(x, \tau)]. \quad (102)$$

Expanding the partition function corresponding to action (102) in powers of $g$, one finds the expression identical to Eq. (97), but with $W(\alpha)$ corresponding to interference of 2D condensates.

When we describe interference of systems with OBC, we use Eqs. (101) and (102), in which the field $\phi(x, \tau)$ is defined on a whole plane, and is not periodic in $\tau$. In both these cases the interaction is present only in some part of the system, so translational invariance is lost in both $x$ and $\tau$ dimensions. Thus to calculate $W(\alpha)$ using Eq. (98) one needs to calculate partition functions of inhomogeneous Sine-Gordon models. Exact solution of Sine-Gordon model is available only for periodic boundary conditions, so to treat open boundary conditions one needs to develop alternative methods. Conversely, if one has a solution for $W(\alpha)$, this provides a tool for calculating partition functions of inhomogeneous Sine-Gordon models and Coulomb gases using Eq. (97). In section IV D we present a novel mapping of $W(\alpha)$ for arbitrary $G(x, y)$ to the statistical properties of random surfaces, which provides a new tool for calculating partition functions of a wide class of Sine-Gordon models and Coulomb gases. In general, partition functions with inhomogeneous $g(x, \tau)$ can be evaluated using this mapping as well. In this case the function $g(x, \tau)$ will appear in the integrand of Eq. (102). We point out that the mapping which we use is not related to the existence of the exact solution of Sine-Gordon models, but relies only on the structure of the correlation functions in the absence of interactions. We also note that a suitable extension of our method can be used to compute correlation functions of Coulomb-gas models in equilibrium and non-equilibrium situations.

C. Distribution functions for 1D gas with periodic boundary conditions

There are three natural ways to compute partition function $Z_0(K, g)$ for complex $g$, which can be used to construct distribution functions for 1D gas with PBC using Eq. (88). The first one is related to the theory of symmetric polynomials and is described in Appendix [3]. The second one relies on the integrability of the quantum impurity model defined by the action (91), and makes use of the thermodynamic Bethe ansatz (TBA). This approach is discussed in Appendix [C]. These two approaches have been presented in Refs. [120], [132]. They are difficult to implement directly and are included for completeness. The third approach described below in section IV C 1 leads to the most transparent answer. Although it is intrinsically related to the first two, it has a broader applicability and will be studied in detail.

At imaginary $g$ the theory (91) is apparently non-Hermitian, but as we will see later the model belongs to a special class of non-Hermitian field theories, which have real spectrum. We describe an interesting connection with $PT$-symmetric quantum mechanics and conformal field theories with negative central charges. Here we also make close contact with the impurity problem in a Luttinger liquid [100]. This connection is based on the following observation: when we consider Taylor expansion of the quantum impurity partition function in powers of $g$, we need to take multipoint correlation functions at the same $x$ but at different times. To calculate higher moments of the interference amplitude signal we need to use correlation functions for equal time, but at different points in space. The two are the same because of the relativistic invariance of the Luttinger liquid [111], [116].

1. Mapping to integrable structure of CFT and singular anharmonic oscillator

To introduce our method we need to make several formal remarks regarding one dimensional conformal field theories. In a recent series of papers Bazhanov, Lukyanov, and Zamolodchikov explored an integrable structure of conformal field theories focusing on connections to solvable problems on lattices [133]. The key ingredients of the solvability of lattice models are the so-called transfer matrix operators $T(\lambda)$. These operators contain information about all integrals of motion as well as excitation spectra of the system. Transfer matrices are defined as a function of the so-called spectral parameter $\lambda$ (in the continuum limit $\lambda$ corresponds to rapidity) and commute for different values of $\lambda$. The latter property is a direct manifestation of the existence of an infinite number of commuting integrals of motion. In
his studies of the 8-vertex model, Baxter [134] introduced the operator $Q(\lambda)$ which helps to find eigenvalues of $T$. Operators $T$ and $Q_\pm$ satisfy a set of commutation relations, in particular [133]
\[
T(\lambda)Q_\pm(\lambda) = Q_\pm(q\lambda) + Q_\pm(q^{-1}\lambda),
\]
where $q = \exp(i\pi/2K)$. So $T$ matrices can be obtained explicitly when one knows the $Q$ operators.

By construction, these operators act in the representation space of Virasoro algebra which can be constructed from the Fock space of free bosonic operators $a_{\pm n}$ defined such that
\[
a_n|p\rangle = 0, \quad \text{for} \quad n > 0, \quad P|p\rangle = p|p\rangle, \quad [a_n, a_m] = \frac{2\pi}{K} \delta_{n+m,0}.
\]
Here $|p\rangle$ denotes the vacuum vector and $p$ is the zero mode of the corresponding bosonic field $\phi(u)$:
\[
\phi(u) = iQ + iPu + \sum_{n\neq 0} \frac{a_n}{n} e^{inu}.
\]

Operators $A_\pm(\lambda) = Q_\pm(\lambda)\lambda^{\mp 4p/K}$ act in the representation space of Virasoro algebra, which can be constructed from the Fock space of bosonic operators $a_{\pm n}$ satisfying $a_n|p\rangle = 0, \, (n > 0)$. The Fock vacuum state $|p\rangle$ is an eigen state of the momentum operator, $P|p\rangle = p|p\rangle$. For $p = N/2$ ($N = 0, 1, 2, ...$) the vacuum eigenvalues of the operator $A_\pm(\lambda)$, $A_\pm(\lambda)|p\rangle = A_\pm^{(vac)}(\lambda)|p\rangle$, are given by (below we consider only the quantities with the + subscript which correspond to the positive $p$)
\[
A^{(vac)}(\lambda) = Z_p(K, -ig),
\]
where
\[
\lambda = \frac{x}{\pi} \sin \left(\frac{\pi}{2K}\right).
\]

At this point it is not clear what we gained by connecting the analytically continued partition function of the impurity problem $Z_p(K, ig)$ to the expectation value of the operator $A^{(vac)}(\lambda)$. As we discuss below, a considerable number of important results have been derived for $A^{(vac)}(\lambda)$. We will be able to make use of these results to obtain the distribution functions of the fringe amplitudes. The function $A^{(vac)}$ has known large-$\lambda$ asymptotics [133]
\[
\log(A^{(vac)}(\lambda)) \sim M(K) \left(-\lambda^2\right)^{\frac{2K}{2K-1}},
\]
where the constant $M(K)$ is given by
\[
M(K) = \frac{\sqrt{T}\Gamma(\frac{1}{2K-1})[\Gamma(\frac{2K}{2K-1})]}{\cos(\frac{\pi}{2K})\Gamma(\frac{K}{2K-1})}.
\]

The function $A^{(vac)}(\lambda)$ is entire function for $K > 1$ and is completely determined by its zeros $\lambda_k, \, k = 0, 1, ...$. Therefore $A^{(vac)}(\lambda)$ can be represented by the convergent product
\[
A^{(vac)}(\lambda) = \prod_{k=0}^{\infty} \left(1 - \frac{\lambda^2}{\lambda_k^2}\right), \quad A^{(vac)}(0) = 1.
\]

On the basis of analysis of a certain class of exactly solvable model, corresponding to the integrable perturbation of the conformal field theory, it was conjectured in [132] that the so-called $Y$-system and related $T$ system (where $Y = e^{\epsilon r}$ and $e_r$ are the Bethe-ansatz energies parametrized by $r$, the nodes of the Dynkin diagrams (see TBA-section) satisfy the same functional equations and possess the same analytical structure and asymptotics as the spectral determinant of the one-dimensional anharmonic oscillator. Further, the same functional equations, analytical properties [110] and asymptotics [109] are satisfied for the vacuum eigenvalues of $Q$-operator for special values of $p$ and the latter are given by the spectral determinant of the following Schrödinger equation
\[
(-\partial_x^2 + x^{2\alpha})\Psi(x) = E\Psi(x).
\]
The spectral determinant is defined as
\[
D(E) = \prod_{n=1}^{\infty} \left(1 - \frac{E}{E_n}\right).
\]
Soon after, in Ref. [136], this conjecture has been extended to all values of $p$:

$$A^{(\text{vac})}(\lambda, p) = D(\rho \lambda^2),$$  \hspace{1cm} (113)

where now $D(E)$ is the spectral determinant of the Schrödinger equation

$$-\partial_x^2 \Psi(x) + \left(x^{2K - 2} + \frac{l(l+1)}{x^2}\right) \Psi(x) = E \Psi(x),$$  \hspace{1cm} (114)

with $l = 4pK - 1/2$. Here $\rho = (4K)^{2-1/K}[\Gamma(1-1/(2K))]^2$. In relation to the interference problem, $p = (md/\hbar) \tan \theta$, where $m$ is the atom’s mass, $d$ is the separation between the two condensates and $t$ is the time when the measurement was done after the free expansion started.

For some values of parameters the Eq. (114) can be solved exactly:

a) $K = 1$. In this case, corresponding to a singular harmonic oscillator $E_n = 4n + 2l - 1$, $n = 1, 2, \ldots$ and

$$D(E, l) = \frac{\Gamma(\frac{3}{4} + \frac{1}{2}) e^{-CE}}{\Gamma(\frac{1}{4} + \frac{1}{2} - \frac{K}{4})},$$  \hspace{1cm} (115)

where $C$ is nonuniversal renormalization constant.

b) The case $K \to \infty$ is recovered by the rigid well potential for which the eigen energies are given by the zeroes of the Bessel function, and therefore

$$D(E, l) = \Gamma(l + 3/2)(\sqrt{E}/2)^{-l-1/2} J_{l+1/2}(\sqrt{E}).$$  \hspace{1cm} (116)

c) For $K = 3/4$ and $l = 0$ the result is expressed in terms of Airy function.

For generic values of $K$ and $l$ the Schrödinger equation can be solved numerically with a very good precision with subsequent computation of the spectral determinant. Alternatively, for $n \geq 5-10$ the spectrum of the equation (114) is very well approximated by the standard WKB expression (see e.g. [137])

$$E_n = \gamma(K)(n - \gamma(K)) \frac{2K}{n - \gamma(K)},$$  \hspace{1cm} (117)

where $n = 1, 2, \ldots$ Here $\gamma(K)$ is the Maslov index. For $1/2 < K < \infty$, $\gamma(K) = \frac{1}{4} - \frac{l}{2}$, for $K = \infty$, $\gamma(K) = -l/2$ and for $0 < K < \frac{1}{2}$, $\gamma(K) = \frac{4K - 2l - 1}{8K}$. The function $\gamma(K)$ in Eq. (117) reads

$$\gamma(K) = \left[\frac{2\sqrt{\pi} \Gamma(\frac{1}{2} + \frac{1}{4K-2})}{\Gamma(1 + \frac{1}{4K-2})}\right]^{2K-1}. $$  \hspace{1cm} (118)

In principle, the function $\gamma(K)$ can be a smooth function interpolating between limiting values given above and can be considered as a noninteger Maslov index [138]. This interpolation allows to (approximately) evaluate the partition function and the distribution function in many cases. In the limiting cases $K \to 1$ and $K \to \infty$ the WKB approximation gives the exact spectrum. We point out however, that using the approximate WKB spectrum carelessly can result in non-physical results, e.g. negative values of the distribution function $W(\alpha)$, and thus have to be used with care. WE emphasize that the solution of the ODE (114) for several potentials gives us already analytically continued function $Z_p(K, ig)$. It seems that this approach to finding the solution is easier and more elegant than the solution of TBA equations with subsequent analytical continuation. Moreover a variety of approximate methods are available for solving Schrödinger-like 1D equations.

Note that Eq. (114) has an interesting duality symmetry which generalizes the Coulomb-harmonic oscillator duality and which allows to relate the $K > 1/2$ and $0 < K < 1/2$ sectors. Making the transformation $x \to y^{1/\alpha}$, $\Psi(x) \to y^\alpha \phi(y)$ with $\lambda = -(1/2) + 1/(2A)$, and $A = 2K$ and then rescaling $y \to \sqrt{\alpha}y$ with $\alpha = (-4K^2/E)^{2K}$ we obtain an equation of the same form with parameters $K', l', E'$ given by

$$4K' = \frac{1}{K}, \hspace{1cm} l' = 2p - 1/2, \hspace{1cm} E' = \frac{1}{4K^2} \left(\frac{4K^2}{E}\right)^{2K}. $$  \hspace{1cm} (119)

The point $K = 1/2$ is a self-dual point of this transformation. Presumably this symmetry is the origin of the Seiberg-Witten type duality in the impurity problem observed in Ref. [139].
2. PT-symmetric quantum mechanics

The surprising link between the ordinary differential equation (114) and the exactly solvable problems related to conformal field theories with negative central charge and/or two-color Coulomb gas on a circle is not limited to the examples considered above. There is a deep connection between these problems and a non Hermitian PT-symmetric quantum mechanics. The latter has been formulated and studied by C. Bender in various contexts [140]. This story goes back to the (unpublished) work of Bessis and Zinn-Justin where, inspired by studies of Lee-Yang model, they conjectured that the 1D Schrödinger equation with potential $V(x) = (ix)^3$ has a real spectrum. Later in Ref. [141] this statement was generalized for arbitrary power-law potentials, $V(x) = (ix)^n$ for real $n$. It was also conjectured that the Schrödinger equation is invariant with respect to combined action of the PT symmetry: $\mathcal{P}$ (parity inversion): $x \rightarrow -x, p \rightarrow -p$; $\mathcal{T}$: $x \rightarrow x, p \rightarrow -p, i \rightarrow -i$.

As we have seen, the partition function of the boundary Sine-Gordon model (as well as of the other related models) is real for the imaginary value of the coupling constant. This is of course not a generic property of field theory models. Therefore the boundary sine-Gordon model with imaginary coupling also belongs to the class of $\mathcal{PT}$-symmetric systems. It is then not surprising why the ODE (114) appears in such a theory.

This intriguing relation between ordinary differential equations and various integrable models has been recently extended to a large class of models (for recent review and references see [142]) by making a close link to the $\mathcal{PT}$-symmetric systems. These relationships have been used [143] for the so-called circular brane model which in certain limit describes the so-called Ambegaokar-Eckern-Schön model [144]. It is therefore reasonable to expect other new interesting applications of these findings.

3. Analysis of distribution functions

Using exact analytic expressions for $Z_p(K, ig)$ at $K = 1$ after the integral transform we obtain the Poissonian distribution for $W_p(\alpha)$. The case $K \rightarrow \infty$ is recovered easily as well. One sees that when $p = 0$, the distribution function is a delta function in the leading order (corrections to this result will be considered in Appendix A). As $p$ increases, $W_p(\alpha)$ rapidly broadens. In the limit of large $K$, the function $W_p(\alpha)$ depends only on the product $Kp$ and takes a simple form:

$$W_p(\alpha) \approx \begin{cases} 
4Kp(1 - \alpha)^{4Kp - 1}, & \alpha < 1 \\
0, & \alpha > 1.
\end{cases} \quad (120)$$

This function is peaked at $\alpha = 1$ for $Kp < 1/4$, it becomes a step function exactly at $Kp = 1/4$, and for $Kp > 1/4$, $W_p(\alpha)$ is a monotonically decreasing function of $\alpha$. When the product $Kp$ becomes large $Kp \gtrsim 1$ the function $W_p$ becomes Poissonian: $W_p(\alpha) \approx 4Kp \exp(-4Kp\alpha)$. In general, the tendency of broadening of the distribution function remains true for all values of $K$.

The distribution given in (120) is a particular representative of class of extreme-value distributions called Weibull distributions. At $p = 0$ the limit of $K \rightarrow \infty$ it will be analytically proven in section IV D 2 that the normalized amplitude $\tilde{\alpha}$ is characterized by the universal Gumbel distribution, for which

$$\tilde{W}(\tilde{\alpha}) = Ke^{K(\tilde{\alpha} - 1 - \gamma) - e^{K(\tilde{\alpha} - 1 - \gamma)}}, \quad (121)$$

where $\gamma \approx 0.577$ is the Euler gamma-constant. The Gumbel distribution, which also belongs to a class of extreme value statistics frequently appears in various situations (see e.g. books [124] for historic introduction), including number theory [145], $1/f$ noise [101], Kardar-Parisi-Zhang growth [146], free Bose gas [147], etc. Probably the whole distribution function in our interference context can in general be characterized as a certain extreme-value statistics.

D. Non perturbative solution for the general case

In this section we present a novel approach for calculating $W(\alpha)$, which is based on a mapping to the statistics of random surfaces [55]. We use this method to evaluate $W(\alpha)$ numerically for a variety of situations. We point out that our method is not related to the existence of the exact solution of Sine-Gordon models, but relies only on the structure of the multi-point correlation functions in the absence of interactions. As has been discussed in section IV B, $W(\alpha)$ is connected to the partition functions of Sine-Gordon models and the partition functions of Coulomb gases. Our mapping provides a new non perturbative tool for calculating partition functions of such models. We expect that there should be numerous applications of our new method to other physical problems which can be related to Sine-Gordon...
and Coulomb gas models. As a concrete application of our approach we prove analytically that for periodic boundary conditions in the limit of large $K$ the distribution of fringe visibilities is given by the Gumbel distribution, one of the extreme value statistical distributions [124].

1. Mapping to the statistics of random surfaces

We start by observing that $G(x, y)$ is real and symmetric. Hence it can be diagonalized on $(0, 1)$ by solving the eigenvalue equations

$$
\int_0^1 G(x, y) \Psi_f(y) dy = G(f) \Psi_f(x).
$$

Here $f$ is an integer index, which goes from 1 to $\infty$. $\Psi_f(x)$ can be chosen to be real and normalized according to

$$
\int_0^1 \Psi_f(x) \Psi_k(x) dx = \delta(f, k).
$$

Then, $G(x, y)$ is given by

$$
G(x, y) = \sum_{f=1}^{f=\infty} G(f) \Psi_f(x) \Psi_f(y).
$$

Decomposition given by Eqs. (122)-(124) is similar to diagonalization of a symmetric matrix by finding its eigenvectors and eigenvalues. Now we can write $Z_{2n}$ from Eq. (83) as

$$
Z_{2n} = \int_0^1 \cdots \int_0^1 du_1 \cdots du_n dv_1 \cdots dv_n e^{\frac{i}{\hbar} \left( \sum_{i<j} G(u_i, u_j) + \sum_{i<j} G(v_i, v_j) - \sum_{i,j} G(u_i, v_j) \right)} = \\
\int_0^1 \cdots \int_0^1 du_1 \cdots dv_n e^{\sum_{f} G_{f}(\xi) \left[ \frac{1}{2} \sum_{i=1}^{f} \Psi_f(u_i)^2 + \sum_{i=1}^{f} \Psi_f(v_i)^2 - \sum_{i=1}^{f} \Psi_f(u_i)^2 - \sum_{i=1}^{f} \Psi_f(v_i)^2 + \left( \sum_{i} \Psi_f(u_i) \right) \left( \sum_{i} \Psi_f(v_i) \right) \right]} = \\
\int_0^1 \cdots \int_0^1 du_1 \cdots dv_n e^{\sum_{f} G_{f}(\xi) \left[ \frac{1}{2} \sum_{i=1}^{f} \Psi_f(u_i) \Psi_f(v_i) - \sum_{i=1}^{f} \Psi_f(u_i) + \sum_{i=1}^{f} \Psi_f(v_i) \right]}.
$$

Square of the first $i-$sum in the last line above can be decoupled by introducing Hubbard-Stratonovich integrations over auxiliary variables $t_f$ :

$$
Z_{2n} = \int_0^1 \cdots \int_0^1 du_1 \cdots dv_n \prod_{f=1}^{f=\infty} \int_{-\infty}^{\infty} dt_f \frac{e^{-t_f^2/2} e^{\sum_{i=1}^{f} t_f \sqrt{G_{f}(\xi) \left( \Psi_f(u_i) - \Psi_f(v_i) \right)} - \frac{G_{f}(\xi)}{\hbar} \left( \Psi_f(u_i)^2 + \Psi_f(v_i)^2 \right)}}{\sqrt{2\pi}}.
$$

Now we can simply integrate over $du_1, ..., dv_n$, since all integrals over $u-$variables are the same (integrals over $v-$variables are also identical):

$$
Z_{2n} = \left( \prod_{f=1}^{f=\infty} \int_{-\infty}^{\infty} e^{-t_f^2/2} dt_f \right) \frac{\sqrt{2\pi}}{\sqrt{2\pi}} g(t_f)^n g(-t_f)^n,
$$

where

$$
g(t_f) = \int_0^1 dx e^{\sum_{i} t_f \sqrt{G_{f}(\xi) \left( \Psi_f(x) - \frac{G_{f}(\xi)}{\hbar} \Psi_f(x) \right)}}.
$$

If all eigenvalues $G(f)$ are negative, then

$$
g(-t_f) = g(t_f)^* \quad g(t_f) g(-t_f) = |g(t_f)|^2.
$$
From comparison of Eqs. (88) and (127) we obtain the central result of this section

$$W(\alpha) = \prod_{f=1}^{f=\infty} \int_{f}^{\infty} e^{-\frac{t_f^2}{2\pi}} dt_f \delta \left(\alpha - |g(\{t_f\})|^2\right).$$  

(129)

Eq. (129) can be used to simulate distributions $W(\alpha)$ using Monte-Carlo approach. First, one needs to solve integral Eqs. (122)-(124) numerically to obtain eigen functions and eigen vectors. Then one needs to choose random numbers typically $\alpha$ given by Eq. (128). Fringe amplitude $\alpha$ is always kept within 1% from its expected value. For most of the presented results, all eigenvalues $G(f)$ are negative, and Eq. (129) can be directly applied. Special care should be taken for 1D case with nonzero temperature, where one of the eigenvalues $G(1)$ can be positive. This situation can be handled by subtracting sufficiently large positive constant $C$ from $G(x, y, \frac{E}{\hbar^2})$, which makes all eigenvalues negative. According to (83), this leads to rescaling of $\alpha$ by a factor $e^{-C}$, which can be easily taken into account.

In Fig. 10 we show scaled distribution of contrast $\tilde{W}(\alpha)$ at $T = 0$ for 1D gases with open boundary conditions, shown for Luttinger parameters $K = 2$ (dashed), $K = 3$ (dotted), and $K = 5$ (solid). The inset shows a comparison between open (solid) and periodic (dashed) boundary conditions for $K = 5$. The figure is taken from Ref. [58].

![Graph](image)

FIG. 10: Scaled distribution functions of the normalized interference amplitude $\tilde{W}(\alpha)$ at $T = 0$ for 1D gases with open boundary conditions, shown for Luttinger parameters $K = 2$ (dashed), $K = 3$ (dotted), and $K = 5$ (solid). The inset shows a comparison between open (solid) and periodic (dashed) boundary conditions for $K = 5$. The figure is taken from Ref. [58].

Result (129) can be interpreted as the mapping of our problem to the statistics of random surfaces (strings in 1D), subject to classical noise. In this mapping, $\Psi_f(x)$ are the eigen modes of the surface vibrations, $t_f$ are the fluctuating mode amplitudes, and $|G(f)|$ is the noise power. This mapping holds as long as all eigenvalues $G(f)$ are negative, as discussed earlier. Infinite dimensional integral over $\{t_f\}$ variables can be understood as an averaging over fluctuations of the surface. For particular realization of noise variables $\{t_f\}$, complex valued surface coordinate at point $x$ is given by

$$h(x; \{t_f\}) = \sum_f t_f \sqrt{\frac{G(f)}{K}} \Psi_f(x) - \frac{G(f)}{2K} \Psi_f(x)^2.$$

(130)

$g(\{t_f\})$ is a number defined for each realization of a random surface $\{t_f\}$, given by Eq. (128). Fringe amplitude $\alpha$ for
FIG. 11: Scaled distribution functions of the normalized interference amplitude \( \tilde{W}(\tilde{\alpha}) \) for a 1D Bose gas with open boundary conditions at nonzero temperature and \( K = 5 \). Different curves correspond to ratios \( K\xi_T/L = \infty \) (solid), \( K\xi_T/L = 1 \) (dotted), and \( K\xi_T/L = 0.25 \) (dashed). \( \xi_T \) is the thermal correlation length, \( K \) is the Luttinger parameter and \( L \) is the imaging length. The figure is taken from Ref. [58].

FIG. 12: Scaled distribution functions of the normalized interference amplitude \( \tilde{W}(\tilde{\alpha}) \) for a two dimensional Bose gas with the aspect ratio of the imaging area equal to unity and open boundary conditions. Temperature is below the Berezinskii-Kosterliz-Thouless (BKT) transition temperature. Different curves correspond to \( T = T_{BKT} \) (BKT transition point, solid), \( T = (2/3)T_{BKT} \) (dashed line) and \( T = (2/5)T_{BKT} \) (dotted line). Above the BKT transition temperature the distribution function is Poissonian. The figure is taken from Ref. [58].

each realization of \( \{t_f\} \) variables is given by

\[
\alpha = |g(\{t_f\})|^2 = \left| \int_0^1 dx e^{h(x;\{t_f\})} \right|^2. \tag{131}
\]

2. From interference of 1D Bose liquids of weakly interacting atoms to extreme value statistics

To illustrate the power of the interpretation of the interference fringe amplitudes in terms of random surfaces, we now prove analytically that for periodic boundary conditions in 1D and in the limit of large \( K \), the normalized distribution \( \tilde{W}(\tilde{\alpha}) \) is given by the Gumbel distribution \([121]\), one of the extreme value statistical distributions. We note that to prove such result using \( 1/K \) expansion described in Appendix [A] one needs to go to the infinite order of the perturbation theory, so essentially our result is non-perturbative in \( K \).

The Gumbel function

\[
P_G(x) = e^{-(x+\gamma)-e^{-(x+\gamma)}} \tag{132}
\]
plays the same role in the extreme value statistics \([124]\) as the usual Gaussian distribution plays in the statistics of the average value. According to the central limit theorem, the average value of \( N \) numbers taken from the same ensemble in the limit of large \( N \) is distributed according to Gaussian function. One can prove similar theorems for
the distribution of the extreme, i.e. largest value of \( N \gg 1 \) numbers taken from the same ensemble. Gumbel function \( G_{\alpha}(x, y) \) is one of the universal functions describing ensembles which decay faster than any algebraic function at infinity. The Gumbel distribution often appears in applied mathematics, including problems in finance or climate studies, where it is used to describe rare events such as stock market crashes or earthquakes.

For the periodic case corresponding to Eq. (131), the eigenvalue problem \((122), (124)\) can be solved analytically, and the noise has \( 1/f \) power spectrum:

\[
G_{\text{per}}(x, y) = \log \frac{1}{\pi} \sin \pi |x - y| = - \log 2\pi - \sum_{f=1}^{\infty} \frac{1}{2f} \left[ (\sqrt{2}\cos 2\pi fx)(\sqrt{2}\cos 2\pi fy) + (\sqrt{2}\sin 2\pi fx)(\sqrt{2}\sin 2\pi fy) \right].
\]  
(133)

Eigen modes are given by simple harmonic functions, and all of them, except one, are doubly degenerate: we will denote corresponding noise variables as \( t_{1,f}, t_{2,f} \) and eigen values as \( G(f) = -1/(2f) \). Simplification in the limit \( K \gg 1 \) stems from the fact that exponent in Eq. (131) can be expanded in Taylor series, since \( h(x; \{t_f\}) \) is small for large \( K \) according to its definition \( (130) \). In this case it can be shown that the distribution of \( \alpha \) is linearly related to roughness, or mean square fluctuation of the surface, as defined in Ref. [101]. It has been shown in Ref. [101] that \( 1/f \) noise results in the Gumbel distribution of the roughness, which has been interpreted in terms of extreme value statistics in Ref. [148].

Terms of the order of \( 1/\sqrt{K} \) vanish in the Taylor expansion of Eq. (131), since the average values of \( \cos(2\pi fx) \) and \( \sin(2\pi fx) \) on interval \((0, 1)\) are equal to 0. To order \( 1/K \), we obtain

\[
\alpha \approx 1 + \frac{1}{K} \log 2\pi - \sum_{f=1}^{\infty} \frac{1}{2fK} (t_{1,f}^2 + t_{2,f}^2 - 2).
\]  
(134)

The constant term \( - \log 2\pi \) in \( G_{\text{per}}(x, y) \) gives a constant rescaling of the distribution function of \( \alpha \), and doesn’t show up in the normalized fringe amplitude \( \tilde{\alpha} \):

\[
\tilde{\alpha} \approx 1 - \sum_{f=1}^{\infty} \frac{1}{2fK} (t_{1,f}^2 + t_{2,f}^2 - 2).
\]  
(135)

Thus to the leading order in \( 1/K \) expansion, the distribution \( Y(x) \) of the rescaled variable

\[
x = -K(\tilde{\alpha} - 1) = \sum_{f=1}^{\infty} \frac{1}{2f} (t_{1,f}^2 + t_{2,f}^2 - 2)
\]  
(136)
equals

\[
Y(x) = \prod_{f=1}^{F=\infty} \int_{-\infty}^{\infty} dt_{1,f} dt_{2,f} e^{-\sum_{f=1}^{\infty} \frac{1}{2f} (t_{1,f}^2 + t_{2,f}^2 - 2)} \delta \left( x - \sum_{f=1}^{f=\infty} \frac{1}{2f} (t_{1,f}^2 + t_{2,f}^2 - 2) \right).
\]  
(137)

To prove Eq. \((121)\), we need to show \( Y(x) = P_G(x) \), where Gumbel function \( P_G(x) \) is given by Eq. \((132)\). Let us introduce new positive variables

\[
u_f = \frac{(t_{1,f}^2 + t_{2,f}^2)}{2f} > 0,
\]  
(138)
then

\[
Y(x) = \prod_{f=1}^{f=\infty} \int_{0}^{\infty} f du_f e^{-\sum_{f=1}^{\infty} f u_f} \delta \left( x - \sum_{f=1}^{f=\infty} (u_f - 1/f) \right).
\]  
(139)

To prove that \( Y(x) = P_G(x) \) we will calculate their Fourier transforms:

\[
Y(is) = \int_{-\infty}^{\infty} dx Y(x) e^{isx} = \prod_{f=1}^{f=\infty} \int_{0}^{\infty} f du_f e^{-fu_f} e^{iu_f - 1/f} = \prod_{f=1}^{f=\infty} \frac{f e^{-is/f}}{f - is} = e^{-\gamma s} \Gamma[1 - is],
\]  
(140)

\[
P_G(is) = \int_{-\infty}^{\infty} dx P_G(x) e^{isx} = \int_{-\infty}^{\infty} dx e^{-(x+\gamma)-e^{-(x+\gamma)}} e^{isx} = e^{-\gamma s} \Gamma[1 - is].
\]  
(141)
The proof above doesn’t illustrate the meaning of $Y(x)$ as of a distribution of extreme value. Here we will follow the method of Ref. [148] and explicitly construct the variable, extreme value of which generates Gumbel distribution. Let us impose a finite cutoff $f_{\text{max}} = N$, and at the end of calculation we will send $N$ to infinity. If one identifies

\begin{align*}
    z_1 &= u_N, \\
    z_2 &= u_{N-1} + u_N, \\
    &\vdots \\
    z_N &= u_1 + u_2 + \ldots + u_N,
\end{align*}

then $\{z_1, \ldots, z_N\}$ is an ordered set (since $u_i > 0$) of outcomes taken from Poissonian distribution, since

$$e^{-\sum_{f=1}^{N} f u_f} = e^{-\sum_{f=1}^{N} z_f},$$

and Jacobian of transformation from variables $\{u_f\}$ to $\{z_f\}$ variables is unity. Then

$$Y_N(x) = N! \int_0^\infty e^{-z_1} dz_1 \int_{z_1}^\infty e^{-z_2} dz_2 \ldots \int_{z_{N-1}}^\infty e^{-z_N} dz_N \delta \left( x - \left( z_N - \sum_{f=1}^{N} \frac{1}{f} \right) \right)$$

is nothing but the shifted distribution of the largest of $N$ numbers taken from the Poissonian distribution, and in the limit of large $N$ this distribution converges to Gumbel function.

One can understand the appearance of the Gumbel distribution by noting that for $K \gg 1$ the distribution function of the interference amplitude is dominated by rare fluctuations of the random periodic 1D string, which are spatially well localized. The Gumbel distribution was introduced precisely to describe similar rare events such as stock market crashes or earthquakes. For open boundary conditions the universal distribution for large $K$ is slightly different from Gumbel function, similar to $1/f$ noise in other systems [101]. But the main properties, like the presence of asymmetry or the asymptotic form of the tails are preserved.

\section{V. CONCLUSIONS}

\subsection{A. Summary}

When we discuss interference experiments with ultracold atoms, the conventional idea of the particle-wave duality takes a new meaning. On the one hand, these experiments probe phase coherence which is typically associated with coherent non-interacting waves. On the other hand, one can use powerful tools of atomic physics to control interactions between atoms in a wide range and to reach the regime of strong correlations. One can also prepare atomic systems in states which would be difficult if not impossible to obtain in optics, e.g. low dimensional condensates with strong thermal or quantum fluctuations. This remarkable combination places interference experiments with ultracold atoms in a unique position: they can address a problem of how the interactions, correlations, and fluctuations affect the coherent properties of matter. This question appears in many areas of physics, including high energy and condensed matter physics, nonlinear quantum optics, and quantum information. While the naive answer that interactions suppress interference turns out to be correct in most cases, the goal of these lecture notes was to demonstrate that the quantitative analysis of this suppression can provide a lot of nontrivial information about the original systems.

We discussed two effects which contribute to the reduction of the interference fringe contrast in matter interferometers. The first effect is the shot noise arising from a finite number of atoms used in a single measurement. This analysis is particularly important for interference experiments with independent condensates in which the position of interference fringes is random and averaging over many shots cannot be performed. In this case one needs to rely on single shot measurements to observe interference patterns. While interference of independent condensates has been discussed before [60, 61, 62, 63, 64, 65, 66, 67, 68, 69, 70, 71, 72], to our knowledge, we provide the first derivation of the full distribution function of the amplitude of interference fringes. Another mechanism of the suppression of the amplitude of interference fringes discussed in these lecture notes is the quantum and thermal fluctuations of the order parameter in low dimensional condensates. The motivation for this discussion comes from the observation that interference experiments between independent fluctuating condensates can be used to study correlation functions in such systems [52]. For example, one can use the scaling of the integrated amplitude of interference patterns to analyze two point correlation functions. This method has been successfully applied by Hadzibabic et al. [32] to observe the Berezinskii-Kosterlitz-Thouless transition in two dimensional condensates. One conceptual approach to understanding interference experiments with independent condensates is to consider them as analogues of the Hanbury Brown and
measurements scales as robust to the inhomogeneous density of atoms. Indeed the Poissonian distribution is related to the fast regime. It is also worth pointing out that we expect the limiting case of the Poissonian distribution to be particularly distribution function computed in the central region. One can also reach similar conclusions in the strongly interacting and at the boundary of the observation region is given by $1 - L^2 / S R_{TF}^2$, thus giving only a small correction to the distribution function computed in the central region. One can also reach similar conclusions in the strongly interacting regime. It is also worth pointing out that we expect the limiting case of the Poissonian distribution to be particularly robust to the inhomogeneous density of atoms. Indeed the Poissonian distribution is related to the fast $1 / \sqrt{x}$ decay of the one-particle correlation functions in the strongly interacting limit. This scaling is a universal feature of the Tonks-Girardeau limit of bosons and is not affected by the weak harmonic trap $[152]$. 

### B. Some experimental issues

We now comment on a few issues relevant for experimental analysis of noise in interference experiments. The amount of information contained in the experimentally measured distribution function is directly related to the number of cumulants which can be accurately extracted. This includes the second cumulant $k_2$, which corresponds to the width of the distribution; the third cumulant $k_3$, which is related to skewness, $g_1 = k_3 / k_2^2$, and describes the asymmetry of the distribution function, and so on. In general, the statistical error in determining the $n$th order cumulant after $N$ measurements scales as $\sqrt{A_n / N}$, where $A_n$ is a constant which grows with $n$ and depends on the higher moments of the distribution. For example, to experimentally distinguish the normal and Gumbel distributions it is necessary that the statistical error in skewness is at least a factor of two smaller than the mean skewness, which for the Gumbel distribution is $g_1 \approx -1.14$. Thus the minimal number of measurements required is $N_{\text{min}} \approx 24 / g_1^2 \approx 20$, where we used $A_3 \approx 6$, appropriate for the normal distribution $[149]$. In practice the required number of measurements may be higher because of the influence of other possible sources of noise. However, it is certainly experimentally feasible.

Another experimentally relevant issue is the effect of the inhomogeneous density due to the parabolic confining potential. While the approach discussed in these lecture notes can be extended to include the inhomogeneous density profile, interpretation of the experimental results is more straightforward when density variations can be neglected. We note that when the condensate density varies gradually in space, the power-law decay of the correlation functions is not strongly affected $[150]$, except that the exponent may be different in different parts of the trap (correlation function exponents typically depend on the density). We expect that qualitatively our results will not change provided that the power law decay of the correlation functions is much stronger than the change of the condensate density in the measured part of the cloud. To be more precise, the best comparison with theory can be achieved when the observation region $L$ is much smaller than the size of the condensate, determined by the effective Thomas-Fermi length $[151]$, $R_{TF} = (3 \hbar^2 / (m^2 \omega_a^2 a_{1D}))^{1/3}$ (here $N$ is the number of atoms of mass $m$ and $a_{1D}$ is the one-dimensional scattering length). As long as $L$ remains much larger than a healing length, our analysis is valid. In the regime of weakly interacting atoms, one can show that the ratio between the effective Luttinger parameter $K$ at the center of the harmonic trap and at the boundary of the observation region is given by $1 - L^2 / 8 R_{TF}^2$, thus giving only a small correction to the distribution function computed in the central region. One can also reach similar conclusions in the strongly interacting regime. It is also worth pointing out that we expect the limiting case of the Poissonian distribution to be particularly robust to the inhomogeneous density of atoms. Indeed the Poissonian distribution is related to the fast $1 / \sqrt{x}$ decay of the one-particle correlation functions in the strongly interacting limit. This scaling is a universal feature of the Tonks-Girardeau limit of bosons and is not affected by the weak harmonic trap $[152]$. 

### C. Outlook

Before concluding these lecture notes we would like to discuss questions which still need to be understood in the context of interference experiments with ultracold atoms. We also suggest an outlook for future theoretical work.

*Combining shot noise with the order parameter fluctuations.* A careful reader has undoubtedly noticed that we discussed either the shot noise or the order parameter fluctuations. At this point we are still lacking theoretical tools which would allow to include both effects simultaneously. One of the difficulties is that such analysis requires the knowledge of the correlation functions for all distances rather than
the long distance asymptotic form. Indeed, in section IV we showed that the short distance part of the correlation functions gives contribution of the same order as the shot noise. In the particular case of the interference of 2D condensates, the knowledge of the short distance behavior of the correlation functions is needed to include the effect of the vortex excitations below the BKT transition.

Stacks of independent condensates

In these lecture notes we focused on interference patterns from a single pair of condensates. However in experiments one often has a stack of several condensates (see e.g. [31]). In this case interference arises from all possible pairs, and the system provides intrinsic averaging and suppression of the noise. For a finite number of condensates self-averaging is not complete and one expects finite fluctuations of the fringe contrast. It would be useful to generalize analysis of the shot noise and order parameter fluctuations to such systems.

Dynamics of interacting atoms.

One of the advantages of the cold atoms systems is the possibility to study non equilibrium coherent dynamics of interacting systems. In particular dynamical splitting of a single condensate into a pair of condensates has been performed in experiments on microchips [30, 42, 48, 49] and stimulated theoretical work on the subject [153, 154, 155, 156, 157, 158]. Similar experiments can also be done using superlattice potentials in optical lattices which are now available in experiments [153, 160]. While analysis of fringe amplitude distribution functions presented in these lecture notes dealt exclusively with systems in the thermodynamic equilibrium, it would be interesting to generalize it to systems undergoing nonequilibrium dynamical evolution.

Interference experiments with fermions

The discussion presented in these notes was limited to the case of interference of bosons. Such experiments can also be done with fermions [161], which are available experimentally in different dimensions [162, 163, 164]. For fermions, modulation of the density can be related to fermion antibunching [20, 165, 166]. Analysis of the noise of the fringe contrast visibility for fermions would be an interesting problem too.

Generalization to other systems

We note that mapping of the Coulomb gas into the statistics of random surfaces introduced in section IV D should have applications beyond calculating the distribution functions of the interference fringe amplitudes. This is a new non perturbative tool to calculate partition functions of a variety of other systems that can be represented as Coulomb gas models. Examples include quantum impurity-related problems [100], Sine-Gordon models where interaction strength can depend on position, and many others. Our mapping is not related to the existence of the exact solution of Sine-Gordon models, but relies only on the factorable structure of the many-point correlation functions in the absence of interactions, which is a general property of a Gaussian action.

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APPENDIX A: LARGE $K$ EXPANSION

In this Appendix, we will describe a systematic ”diagrammatic technique” to calculate $Z_{2n}$ or $Z_{2n}^{pert}$ as an expansion in small parameter $1/K$. It corresponds to the ”high temperature” limit of the classical gas analogy discussed in section IV A. This expansion can be applied both in 1D or 2D, and can be used to study the limiting distribution at large $K$, which for PBC in 1D has been conjectured [56] and proven [58] to be the Gumbel distribution [124].

1. Expansion to order $(1/K)^2$.

We will start from the 1D case by expanding the exponent in Eq. (83):
The total number of terms is also \( n(n-1)/2 \), and the total number of \( u - v \) terms is \( n^2 \). The latter come with \(-1\) sign, so the total expression for \( Z_{2n} \) up to \( O(1/K^2) \) is

\[
Z_{2n} = 1 - \frac{1}{K} n I_0 + O\left(\frac{1}{K^2}\right). \tag{A3}
\]

In second order, dependence of \( Z_{2n} \) on \( G(x, y) \) comes through three different integrals:

\[
I_{1,2} = \int_0^1 \int_0^1 \int_0^1 dx dy dz dt G(x, y) G(z, t) = \left(\int_0^1 \int_0^1 dx dy G(x, y)\right)^2 = I_0^2, \tag{A4}
\]

\[
I_{2,2} = \int_0^1 \int_0^1 \int_0^1 dx dy dz G(x, y) G(y, z), \tag{A5}
\]

\[
I_{3,2} = \int_0^1 \int_0^1 dx dy G(x, y) G(x, y). \tag{A6}
\]

Pictorially, these expressions can be represented by the corresponding diagrams, shown in Fig. 13. There, horizontal solid line corresponds to \( G(x_i, x_j) \). If two ends are connected by a dashed line, then in the integral these two ends should correspond to the same variable. To write the expression for \( Z_{2n} \) in \( 1/K^2 \) order, one has to calculate the total number of terms corresponding to \( I_{1,2}, I_{2,2}, I_{3,2} \). Clearly, it will be some universal polynomial depending on \( n \), which is determined by combinatorics. Here we describe how to do this combinatorics in detail.

When parenthesis are multiplied in \( (A1) \) there are several types of expressions which appear, and come with different signs to the integral:

\[
\left( \sum_{i<j} G(u_i, u_j) + \sum_{i<j} G(v_i, v_j) - \sum_{i<j} G(u_i, v_j) \right)^2 = \sum_{i_1, j_1, i_2, j_2} G(u_{i_1}, v_{j_1}) G(u_{i_2}, v_{j_2}) + \sum_{i_1<j_1, i_2<j_2} (G(u_{i_1}, u_{j_1}) G(u_{i_2}, u_{j_2}) + G(v_{i_1}, v_{j_1}) G(v_{i_2}, v_{j_2}) + 2G(u_{i_1}, u_{j_1}) G(v_{i_2}, v_{j_2}))
\]

\[ -2 \sum_{i_1<j_1, i_2<j_2} (G(u_{i_1}, u_{j_1}) G(u_{i_2}, v_{j_2}) + G(v_{i_1}, v_{j_1}) G(u_{i_2}, v_{j_2})). \tag{A7}
\]
These are (signs are indicated)

\[ + G(u_{i1}, u_{j1})G(u_{i2}, u_{j2}), \text{ and equivalent after integration } G(v_{i1}, v_{j1})G(v_{i2}, v_{j2}); \quad (A8) \]

\[ + G(u_{i1}, v_{j1})G(u_{i2}, v_{j2}); \quad (A9) \]

\[ - G(u_{i1}, u_{j1})G(u_{i2}, v_{j2}), \text{ and equivalent after integration } - G(v_{i1}, v_{j1})G(u_{i2}, v_{j2}); \quad (A10) \]

\[ + G(u_{i1}, u_{j1})G(v_{i2}, v_{j2}); \quad (A11) \]

For terms which have \( G(u_{ik}, u_{jk}) \) or \( G(v_{ik}, v_{jk}), k = \{1, 2\} \) there is a restriction that \( ik < jk \), but there is no such restriction for \( G(u_{ik}, v_{jk}) \).

Let’s calculate the total number of \( I_{1,2}, I_{2,2} \) and \( I_{3,2} \) terms for \( u - u - u - u \) expressions of type \( \text{[AS]} \). Total number of \( I_{1,2} \) terms compatible with restrictions is

\[ \frac{1}{2!^2} n(n - 1)(n - 2)(n - 3). \quad (A12) \]

\( 1/2!^2 \) appears since pairs \( \{i1, j1\} \) and \( \{i2, j2\} \) are ordered, while \( n(n - 1)(n - 2)(n - 3) \) is the total number of ways to choose a sequence of four different not ordered numbers out of the set of size \( n \).

Total number of \( I_{2,2} \) terms compatible with restrictions is

\[ \frac{1}{2!^2} n(n - 1)(n - 2) * 4. \quad (A13) \]

Here the situation is similar to \( I_{1,2} \), but there is a "topological" prefactor of 4, which corresponds to four topologically equivalent diagrams for \( I_{2,2} \), as shown in Fig. 14a.

Finally, total number of \( I_{3,2} \) terms is

\[ \frac{1}{2!^2} n(n - 1) * 2. \quad (A14) \]

Here, 2 is again a "topological" prefactor, corresponding to two topologically equivalent diagrams shown in Fig. 14b.

Overall, \( u - u - u - u \) term of \( \text{[AS]} \) gives

\[ \frac{1}{4} n(n - 1)(n - 2)(n - 3)I_{1,2} + n(n - 1)(n - 2)I_{2,2} + \frac{1}{2} n(n - 1)I_{3,2} \quad (A15) \]

As a simple check of combinatorics one can calculate the total numbers of \( u - u - u - u \) terms of type \( \text{[AS]} \). It is

\[ \frac{1}{2!^2} n(n - 1)(n - 2)(n - 3) + \frac{1}{2!^2} n(n - 1)(n - 2) * 4 + \frac{1}{2!^2} n(n - 1) * 2 = \left( \frac{n(n - 1)}{2!} \right)^2, \quad (A16) \]

as it should be from the calculation which neglects the dashed lines, and doesn’t impose any restrictions on the terms at different horizontal rows.

Analogously, one can calculate polynomials for expressions \( \text{[A9]} - \text{[A11]} \), and results are as follows, respectively:

\[ n^2(n - 1)^2I_{1,2} + 2n^2(n - 1)I_{2,2} + n^2I_{3,2}, \quad (A17) \]

\[ - \frac{1}{2} n^2(n - 1)(n - 2)I_{1,2} - n^2(n - 1)I_{2,2}, \quad (A18) \]

\[ \frac{1}{2!^2} n^2(n - 1)^2I_{1,2}. \quad (A19) \]
Summing all the terms with corresponding prefactors from (A7), we obtain

\[ Z_{2n} = 1 - \frac{1}{K} nI_0 + \frac{1}{2!K^2} (P_{1,2}I_{1,2} + P_{2,2}I_{2,2} + P_{3,2}I_{3,2}) + O\left(\frac{1}{K^3}\right) = \]

\[ 1 - \frac{1}{K} nI_0 + \frac{1}{2!K^2} (3n(n-1)I_{1,2} - 4n(n-1)I_{2,2} + (2n-1)nI_{3,2}) + O\left(\frac{1}{K^3}\right), \]  

(A20)

where \( P_{1,2} = 3n(n-1), P_{2,2} = -4n(n-1), \) and \( P_{3,2} = (2n-1)n \) are universal polynomials of \( n \). Notice, that due to

cancellations of terms in (A8), the overall degree of polynomials in \( 1/K^2 \) order is \( m = 2 \), compared to

naively expected \( 2m = 4 \).

Using (A20), one can calculate \( Z_{2n}/Z_2^n \) up to \( O(1/K^3) \):

\[ \frac{Z_{2n}}{Z_2^n} = 1 + \frac{n(n-1)}{K^2} (I_{1,2} - 2I_{2,2} + I_{3,2}) + O\left(\frac{1}{K^3}\right) \]  

(A21)

Calculated values of \( \{I_0, I_{1,2}, I_{2,2}, I_{3,2}\} \) and \( \{I_{0,\text{per}}^{\text{per}}, I_{1,2,\text{per}}^{\text{per}}, I_{2,2,\text{per}}^{\text{per}}, I_{3,2,\text{per}}^{\text{per}}\} \) are

\[ I_0 = -\frac{3}{2} = -1.5, \quad I_{0,\text{per}} = -2\pi \approx -6.283; \]  

(A22)

\[ I_{1,2} = \frac{9}{4} \approx 2.25, \quad I_{1,2,\text{per}} = (2\pi)^2 \approx 3.141; \]  

(A23)

\[ I_{2,2} = \frac{51 - \pi^2}{18} \approx 2.28502, \quad I_{2,2,\text{per}} = (2\pi)^2 \approx 3.141; \]  

(A24)

\[ I_{3,2} = \frac{7}{2} \approx 3.5, \quad I_{3,2,\text{per}} = \pi^2 + 12(2\pi)^2 \approx 12.566. \]  

(A25)

Thus, for \( K \to \infty \), the limiting ratio of the widths of the distributions for OBC and PBC are

\[ \frac{Z_4/(Z_2)^2 - 1}{Z_4^{\text{per}}/(Z_2^{\text{per}})^2 - 1} \rightarrow \frac{I_{1,2} - 2I_{2,2} + I_{3,2}}{(I_{1,2,\text{per}}^{\text{per}} - 2I_{2,2,\text{per}}^{\text{per}} + I_{3,2,\text{per}}^{\text{per}})} \approx 1.43465. \]  

(A26)

2. General properties of \( (1/K)^m \) terms, and expansion to order \( (1/K)^3 \).

From the expansion of the previous subsection, one can formulate the general properties of the ”diagram technique” to calculate terms up to \( (1/K)^m \):

a) First, one has to draw all possible topologically inequivalent diagrams, which consist of \( m \) horizontal solid lines, with some of the ends connected by dashed lines. Each end can have at most two dashed lines coming out of it.

b) Ends which are connected by a dashed line correspond to the same variable. Diagrams for which two opposite ends of the horizontal line correspond to the same variable are excluded.

c) Expression which corresponds to a diagram is constructed the following way: if variables at the end of a given solid horizontal line are \( x \) and \( y \), then \( G(x,y) \) should be put as one of the terms in the product under the integrand. Thus the integrand consists of the product of function \( G \) of some variables \( m \) times. Diagrams for which the integrands are the same up to relabeling of the variables are considered to be identical.

d) all free variables should be integrated from 0 to 1.

Diagrams can be connected or disconnected. For example, in Fig. 13, diagram corresponding to \( I_{1,2} \) is disconnected, and diagrams corresponding to \( I_{2,2} \) and \( I_{3,2} \) are connected. The integral which corresponds to a disconnected diagram is a product of expressions, corresponding to its parts: for example, \( I_{1,2} = I_0 + I_0 \).

If the number of topologically inequivalent diagrams of order \( m \) is \( g(m) \), then the term of the order \( 1/K^m \) in \( Z_{2n} \) has the following form:

\[ \int_0^1 \ldots \int_0^1 du_1 \ldots dv_n \frac{1}{m!K^m} \left( \sum_{i<j} G(u_i, u_j) + \sum_{i<j} G(v_i, v_j) - \sum_{ij} G(u_i, v_j) \right)^m = \frac{1}{m!K^m} \sum_{r=1}^{g(m)} P_{r,m}(n) I_{r,m}, \]  

(A27)
where $P_{r,m}(n)$ are universal polynomials of $n$, which can be calculated combinatorially, as described in section A1. Polynomials $P_{r,m}(n)$ should satisfy the following requirements:

a) For positive integer $n$, the values of $P_{r,m}(n)$ are integer, and

$$P_{r,m}(0) = 0.$$  \hspace{1cm} (A28)

b) If one sets $G(x, y) = 1$, then $I_{r,m} = 1$, and lhs of (A27) can be trivially calculated. This implies

$$\sum_{r=1}^{g(m)} P_{r,m}(n) = (-n)^m.$$ \hspace{1cm} (A29)

c) Degree of each polynomial $P_{r,m}(n)$ is not larger than $m$. This has been shown above for $m = 1$ and $m = 2$, and has been checked up to $m = 5$, although we didn’t succeed in proving it directly. This conjecture is supported by the fact, that it guarantees that for any $G(x, y)$ for $K \to \infty$ distributions fit on the top of each other after proper rescaling (see next section).

One can do the combinatorial calculations similar to previous section for $m = 3$, and it takes about a day “by hand”. We will only show the results here. For $m = 3$ there are $g(3) = 8$ topologically inequivalent diagrams, which are shown in Fig. 15 out of which 5 diagrams ($I_{1,3}$ to $I_{5,3}$) are connected. Expressions corresponding to each diagram and universal polynomials are, respectively:

$$I_{1,3} = \int_0^1 \int_0^1 \int_0^1 dxdydztG(x, y)G(x, z)G(z, t), P_{1,3}(n) = -12n(n - 1)(2n - 3),$$  \hspace{1cm} (A30)

$$I_{2,3} = \int_0^1 \int_0^1 \int_0^1 dxdydztG(x, y)G(x, z)G(x, t), P_{2,3}(n) = 12n(n - 1),$$  \hspace{1cm} (A31)

$$I_{3,3} = \int_0^1 \int_0^1 \int_0^1 dxdydG(x, y)G(x, z) G(x, t), P_{3,3}(n) = -12n(n - 1),$$ \hspace{1cm} (A32)

$$I_{4,3} = \int_0^1 \int_0^1 \int_0^1 dxdydG(x, y)G(x, z) G(y, z), P_{4,3}(n) = 4n(n - 1)(2n - 1),$$ \hspace{1cm} (A33)

$$I_{5,3} = \int_0^1 \int_0^1 dxdG(x, y)^3, P_{5,3}(n) = -n, \hspace{1cm} (A34)$$

$$I_{6,3} = I_{2,2} * I_0, P_{6,3}(n) = 36n(n - 1)(n - 2), \hspace{1cm} (A35)$$

$$I_{7,3} = I_{3,2} * I_0, P_{7,3}(n) = -3n(n - 1)(2n - 3), \hspace{1cm} (A36)$$

$$I_{8,3} = I_{1,2} * I_0 = I_{0}^{3}, P_{8,3}(n) = -15n(n - 1)(n - 2). \hspace{1cm} (A37)$$

Numerically evaluated integrals for OBC and PBC are

$$I_{1,3} = -3.49399, I_{1,3}^{\text{per}} = -6.20797; \hspace{1cm} (A38)$$

$$I_{2,3} = -3.5268, I_{2,3}^{\text{per}} = -6.20797; \hspace{1cm} (A39)$$

$$I_{3,3} = -5.32704, I_{3,3}^{\text{per}} = -7.71956; \hspace{1cm} (A40)$$

$$I_{4,3} = -4.21255, I_{4,3}^{\text{per}} = -6.50848; \hspace{1cm} (A41)$$

$$I_{5,3} = -11.25, I_{5,3}^{\text{per}} = -12.5458; \hspace{1cm} (A42)$$

$$I_{6,3} = -3.42753, I_{6,3}^{\text{per}} = -6.20797; \hspace{1cm} (A43)$$

$$I_{7,3} = -5.25, I_{7,3}^{\text{per}} = -7.71956; \hspace{1cm} (A44)$$

$$I_{8,3} = -3.375, I_{8,3}^{\text{per}} = -6.20797. \hspace{1cm} (A45)$$

For $m > 3$ it becomes too cumbersome to manually calculate universal polynomials $P_{r,m}$. We wrote a program in Mathematica, which expands $m$ th term of (A11) directly, and calculates the values $\{P_{r,m}(0), ..., P_{r,m}(n)\}$ using powerful pattern recognition tools. After that, $P_{r,m}(n)$ is recovered using Newton’s formula. Results for $m = 3$ can be recovered that way. One can also check in each order that the degree of $P_{r,m}(n)$ is not larger than $m$. For each $m$
FIG. 15: Topologically inequivalent diagrams for $m = 3$. Diagrams $I_{1,3} - I_{5,3}$ are connected and $I_{6,3} - I_{8,3}$ are disconnected.

FIG. 16: Irreducible diagrams for $m = 4$.

the program needs as an input all topologically inequivalent diagrams, and currently the results have been extended to $m = 5$. For $m = 4$, the overall number of diagrams is 23, out of which 12 are irreducible, and shown in Fig. 16. For $m = 5$, the overall number of diagrams is 66, out of which 33 are irreducible. Numerical prefactors of polynomials $P_{r,m}(n)$ grow with $m$, while their overall sum has a prefactor 1, as follows from (A29). For example, one of the polynomials for $m = 5$ has a prefactor of 384. This puts stringent requirements on the errors in calculation of $I_{r,m}(n)$, so going beyond $m = 5$ will require additional numerical effort. For $m = 5$ results are reliable, since we can check the numerical accuracy of the calculation of integrals by comparison with analytically proven Gumbel distribution.

While calculation of original $Z_{2n}$ requires $2n$ – dimensional numerical integration, calculation of the $I_{r,m}$ requires at most $m$ – dimensional numerical integration. This can be seen from the fact that even for irreducible diagrams, only the parts which contain "loops" have to be integrated numerically. Horizontal lines which have a free end can be "integrated out" analytically, and the dimension of the numerical integral has at most dimension of the loop. For example, first and third horizontal bars in $I_{1,3}$ of Fig 15 can be integrated out using the following identity:

$$\int_0^1 dy \log |x - y| = -1 + \log (1 - x) - x \log (1 - x) + x \log x,$$

so one has to do only 2-dimensional integral numerically. Analogously, diagram $I_{2,3}$ requires only one dimensional numerical integration. In each order $m$ there is only one diagram which requires $m$ – dimensional integration, i.e. $I_{4,3}$ in Fig 15 for $m = 3$. All the rest require at most $(m - 1)$ – dimensional integration.
3. Properties of the $K \to \infty$ distribution

For $K \to \infty$ distribution function becomes very narrow, and it is an interesting question to investigate the limiting behavior of the distribution function. Let us consider the distribution of normalized contrast $\tilde{\alpha}$, defined by $\langle \tilde{\alpha} \rangle$. Due to normalization the following relations hold:

$$\int_0^\infty \tilde{W}(\tilde{\alpha})d\tilde{\alpha} = 1, \quad \int_0^\infty \tilde{\alpha}\tilde{W}(\tilde{\alpha})d\tilde{\alpha} = 1.$$  \hfill (A47)

For large $K$ distribution function is peaked near $\tilde{\alpha} = 1$, and its width is proportional to $1/K$. To calculate the properties of the universal function we assume that it decays relatively fast (exponential decay is enough) away from $\tilde{\alpha} = 1$ for large $K$, so we can extend integrations to infinity. If we define the fluctuation of the normalized contrast $\tilde{\beta} = \tilde{\alpha} - 1$ with distribution function $W_0(\tilde{\beta}) = W(1 + \tilde{\beta})$, then

$$-1 + \frac{Z_2}{Z_2^2} = \int_0^\infty (\tilde{\alpha} - 1)\tilde{W}(\tilde{\alpha})d\tilde{\alpha} = \int_{-1}^\infty \tilde{\beta}W_0(\tilde{\beta})d\tilde{\beta} \equiv \bar{M}_1,$$ \hfill (A48)

$$-1 + \frac{Z_4}{Z_2^2} = \int_0^\infty (\tilde{\alpha}^2 - 1)\tilde{W}(\tilde{\alpha})d\tilde{\alpha} \approx \int_{-\infty}^\infty (2\tilde{\beta} + \tilde{\beta}^2)W_0(\tilde{\beta})d\tilde{\beta} \equiv 2\bar{M}_1 + \bar{M}_2,$$ \hfill (A49)

$$-1 + \frac{Z_6}{Z_2^2} = \int_0^\infty (\tilde{\alpha}^3 - 1)\tilde{W}(\tilde{\alpha})d\tilde{\alpha} \approx \int_{-\infty}^\infty (3\tilde{\beta} + 3\tilde{\beta}^2 + \tilde{\beta}^3)W_0(\tilde{\beta})d\tilde{\beta} \equiv 3\bar{M}_1 + 3\bar{M}_2 + \bar{M}_3,$$ \hfill (A50)

$$\ldots,$$ \hfill (A51)

where $\bar{M}_1, \bar{M}_2, \bar{M}_3$ are the moments of $W_0(\tilde{\beta})$. Inverting (A48)-(A50), one can find $\bar{M}_2, \bar{M}_3, \ldots$ as series expansion in powers of $1/K$. If for large $K$ the distribution functions collapse on the top of each other after proper rescaling, as has been conjectured in \cite{50}, then the expansion of $\bar{M}_m$ in powers of $1/K$ should start from $1/K^m$. Below we will show, that this is necessarily true for all $n$, if the degrees of universal polynomials $P_{r,m}(n)$ are not larger than $m$ and $P_{r,m}(0) = 0$.

To show this, let us consider an analog of Eqs. (A48)-(A50) for the fluctuation of the unnormalized contrast $\beta = \alpha - 1$ with distribution function $W_0(\beta) = W(1 + \beta)$ (normalized contrast considered in Appendix A3 is related to $\alpha$ by simple rescaling). Then

$$Z_{2n} = \int_0^\infty \alpha^nW(\alpha)d\alpha = \int_{-1}^\infty (1 + \beta)^nW_0(\beta)d\beta \approx \int_{-\infty}^\infty \sum_{i=0}^{i=n} C^n_i \beta^i W_0(\beta)d\beta = \sum_{i=0}^{i=n} C^n_i M_i,$$ \hfill (A52)

where $M_i$ is the $i$-th moment of $W_0(\beta)$. We need to show that expansion of $M_i$ in powers of $1/K$ starts only from the terms of the order of $1/K^i$. This can be seen from expansion (A27) together with (A52). If degrees of $P_{r,m}(n)$ are not larger than $m$, then (A27) means that $1/K^m$ terms in $Z_{2n}$ grow at most as $n^m$ for large $n$. On the other hand, $C^n_i = \frac{n!}{i!(n-i)!}$ grows as $n^i$ for large $n$, and if expansion of $M_i$ in powers of $1/K$ starts before $1/K^i$, this will contradict the previous sentence.

To find the first nontrivial contribution to $M_m$, one has to go to order $1/K^m$ in the expansion. Using results for $m = 5$ calculated above, one can calculate for $K \to \infty$ limiting behavior of $K^2\bar{M}_2, \ldots, K^5\bar{M}_5$ for periodic and non-periodic cases, and compare it with the result of the Gumbel distribution

$$\tilde{W}_G(\tilde{\beta}) = Ke^{K\tilde{\beta} - \gamma - e^{K\tilde{\beta} - \gamma}},$$ \hfill (A53)

where $\gamma \approx 0.577$ is the Euler gamma-constant. One obtains:

Non-periodic : $K^2\bar{M}_2 \to 2.35991, K^3\bar{M}_3 \to -5.105577, K^4\bar{M}_4 \to 37.5258, K^5\bar{M}_5 \to -242.492, \frac{\bar{M}_2^2}{\bar{M}_2^3} \to 1.98336$ \hfill (A54)

Periodic : $K^2\bar{M}_2^{per} \to 1.64493, K^3\bar{M}_3^{per} \to -2.40411, K^4\bar{M}_4^{per} \to 14.6114, K^5\bar{M}_5^{per} \to -64.4321$ \hfill (A55)

Gumbel : $K^2\bar{M}_2^G \to 1.64493, K^3\bar{M}_3^G \to -2.40411, K^4\bar{M}_4^G \to 14.6114, K^5\bar{M}_5^G \to -64.4321, \left(\frac{\bar{M}_2^G}{\bar{M}_2^G}ight)^2 \to 1.29857$ \hfill (A56)

$$\bar{M}_2^G \equiv \frac{\bar{M}_2^G}{\bar{M}_2^G} \equiv 2.$$ \hfill (A57)
Clearly, for periodic case the agreement with the Gumbel distribution is excellent, but for non-periodic case, one can unambiguously conclude that limiting function is NOT a Gumbel function. For non-periodic case the function is more widely distributed compared to the Gumbel function, as evident from the higher moments. Here for completeness we provide numerical results for \( Z_{2n}^{\text{per}} \) up to \( m = 5 \):

\[
Z_{2n}^{\text{per}} = 1 + \frac{1.8379n}{K} - \frac{0.4112n + 2.5114n^2}{K^2} + \frac{0.1002n - 0.1548n^2 + 2.1456n^3}{K^3} - \frac{0.03382n + 1.0804n^2 - 0.7399n^3 + 1.7369n^4}{K^4} + \frac{0.01535n + 1.05941n^2 - 0.161727n^3 + 0.172302n^4 + 0.93098n^5}{K^5},
\]

\[
Z_{2n} = 1 + \frac{1.5n}{K} - \frac{0.55496n + 2.30496n^2}{K^2} + \frac{0.049657n + 0.343838n^2 + 1.481505n^3}{K^3} - \frac{0.04741n + 2.0329n^2 - 1.8735n^3 + 1.8256n^4}{K^4} + \frac{0.0106468n + 3.14868n^2 - 4.1829n^3 + 2.8980n^4 + 0.0943039n^5}{K^5}.
\]

4. \( D=2 \)

Here we will briefly consider results of the expansion in powers of \( 1/K \) for 2-dimensional case. Correlation function below the BKT transition is given by Eq. (78), and for square imaging area with unity aspect ratio, all discussions of one dimensional case carry over, with substitutions

\[
\int_0^1 du_i \rightarrow \int_0^1 \int_0^1 d^2 \vec{u}_i, \int_0^1 dv_i \rightarrow \int_0^1 \int_0^1 d^2 \vec{v}_i, G(\vec{x}, \vec{y}) = \log |\vec{x} - \vec{y}|,
\]

and \( (79) \). In 2D case, there is one extra degree of freedom which can be controlled in experiments, which is the aspect ratio of the observation region. If the aspect ratio of the observation region is very large, then the distribution function is essentially the same as in the one-dimensional case. Below we concentrate on the case with unity aspect ratio. Similarly to one-dimensional case, the integral

\[
\int_0^1 \int_0^1 dx_1 dy_1 \frac{1}{2} \log ((x_1 - x_2)^2 + (y_1 - y_2)^2)
\]

can be evaluated analytically, which somewhat simplifies the numerical evaluation. However, the dimensions of integrals grow fast with \( m \), and here we present results only up to \( m = 3 \):

\[
Z_{2n} = 1 + \frac{0.805087n}{K} - \frac{0.0740n + 0.3342n^2}{K^2} + \frac{0.015n + 0.045n^2 + 0.241n^3}{K^3} + O\left(\frac{1}{K^4}\right).
\]

The error in numerical coefficient due to integration is of the order of the last reported digit. Compared to 1D case, convergence is faster, which is consistent with our general ideology that the role of fluctuations is larger in systems with lower dimensions.

APPENDIX B: JACK POLYNOMIALS

The microcanonical partition functions \( Z_{2n}^{(p)} \) can be computed using so-called Jack polynomials. The Jack polynomials belong to a class of symmetric polynomials \([128, 171]\) and are a one-parametric generalization of Schur symmetric functions and generically can be defined as

\[
J_\lambda(x; \beta) = \sum_{\mu \leq \lambda} v_{\lambda,\mu}(\beta) m_\mu(x),
\]

where the monomial symmetric function \( m_\lambda(x) \)

\[
m_\lambda(x) = \sum_{\sigma} x_{\sigma(1)}^{\lambda_{\sigma(1)}} \cdots x_{\sigma(n)}^{\lambda_{\sigma(n)}}
\]

includes sum over all permutations \( \sigma \). Here \( v_{\lambda,\mu} \) are numerical coefficients and \( v_{\lambda,\lambda} = 1 \). The ordering in sum \( (31) \) means the ordering on the set of partitions \( \lambda = (\lambda_1, \lambda_2, \ldots, \lambda_n), \lambda_1 \geq \lambda_2 \geq \ldots \geq \lambda_n \geq 0 \) and \( \mu = (\mu_1, \mu_2, \ldots, \mu_n) \).
\[ \mu_1 \geq \mu_2 \geq \ldots \geq \mu_n \geq 0 \] of natural numbers \( \lambda \) and \( \mu \). The ordering \( \mu \leq \lambda \) here means that \( \sum_{j=1}^{k} \mu_j \leq \sum_{j=1}^{\lambda} \lambda_j \) for any \( k \leq n-1 \). The partitions can be represented using Young diagrams as follows: we put \( \lambda_1 \) boxes in the first row, \( \lambda_2 \) boxes in the second, and so on. The definition of Jack polynomials thus includes a sum over all Young diagrams for which the number of rows is smaller or equal to the number of variables \( n \). The Jack polynomials have the property of orthogonality, which allows to bring the microcanonical partition function into the following form (see [120] for more details on derivation)

\[
Z_{2n}^{(p)} = c_n \sum_{\lambda} \prod_{i=1}^{n} \frac{\Gamma[\lambda_i + \frac{2K}{2}(n-i+1)]\Gamma[p + \lambda_i + \frac{2K}{2}(n-i+1)]}{\Gamma[\lambda_i + 1 + \frac{2K}{2}(n-i)]\Gamma[p + \lambda_i + \frac{2K}{2}(n-i)]}.
\]

Here \( c_n = \Gamma[n+1]/\Gamma[1/2K] \) and sum goes over Young diagrams labeled by integers \( \lambda_1 \geq \lambda_2 \geq \ldots \lambda_n \geq 0 \).

In particular, the Jack polynomials expressions give the following results for the lowest microcanonical partition functions

\[
Z_2^{(p)} = \frac{\sin(\pi \frac{n}{K})\Gamma(1-1/K)}{\sin(\pi(\frac{1}{2K} + p)\Gamma(1 - \frac{1}{2K} + p)\Gamma(1 - \frac{1}{2K} - p))}
\]

and

\[
Z_4^{(0)} = \frac{1}{\Gamma^4(\frac{1}{2K})}\frac{2^{-1/K}\Gamma(\frac{1}{2} - \frac{1}{2K})\Gamma^2(\frac{1}{2})\Gamma^2(\frac{1}{K})}{\sqrt{\pi^2(1 - \frac{2K}{2})\Gamma^2(1 + \frac{2}{2K})}}\left[p_F\left(\frac{1}{1}, \frac{1}{K}, \left\{1 + \frac{2K}{2}, 1 + \frac{2K}{2}\right\}, 1\right)ight. \\
- \left. \sum_{n=0}^{\infty} \frac{\Gamma(\frac{1}{2} + n)^2}{\Gamma(2 + n)^2}\left[p_F\left(\frac{1}{1}, \frac{1}{2K} + n, 1 + \frac{2K}{2} + n, \{2 + n, 2 + n\}, 1\right)\right]\right],
\]

where \( p_F \) is a generalized hypergeometric function [172]. Apparently, for larger \( n \), expressions for \( Z_{2n}^{(p)} \) become more and more cumbersome and are difficult to use for real computations.

Jack polynomials appear particularly in connection to Calogero-Sutherland model, i.e. the model which describes \( n \) particles (the same \( n \) as in formulas above) on a line interacting via inverse square interaction. In the first quantized form the Hamiltonian of this model can be defined as [128]

\[
H_{CS} = -\frac{1}{2} \sum_{j=1}^{n} \frac{\partial^2}{\partial x_j^2} + \pi^2 \sum_{1 \leq i < j \leq n} \frac{g(g-1)}{\sin^2 \pi(x_i - x_j)^2}.
\]

APPENDIX C: THERMODYNAMIC BETHE ANSATZ OF THE QUANTUM IMPURITY MODEL

The model described by the partition function [122] can be solved exactly by thermodynamic Bethe ansatz (TBA) [173, 174]. This solution is non perturbative in \( g \). One should be careful about the correspondence between the perturbative expansion and the results from the non perturbative TBA calculations. The correct correspondence between \( Z(K, g)_{pert} \) and the non perturbative result of the TBA is given by

\[
Z_0(K, g)_{pert} = \frac{1}{\sqrt{2K}} \exp[\tilde{F}_{TBA}(K, g)],
\]

where \( \tilde{F}_{TBA}(K, g) \) is the free energy. It can be expressed in terms of the energies \( \epsilon(\theta) \) of elementary excitations which depend on rapidity \( \theta \) :

\[
\tilde{F}_{TBA}(K, g) = \int_{-\infty}^{\infty} \frac{d\theta}{2\pi \cosh[(2K - 1)(\theta - \alpha)]} \log(1 + \exp(\epsilon_+(\theta))).
\]

Here,

\[
\alpha = \log\left(\frac{x}{\Gamma(\frac{2K}{2})}\frac{2\sqrt{\pi}\Gamma(\frac{4K}{2} - 1)}{\Gamma(\frac{2K}{2} - 1)}\right)
\]

and the variable \( x \) is the fugacity in the perturbative expansion [94].
To clarify the issue further we briefly describe the formalism of TBA. The TBA equations are parametrized by the Dynkin diagrams corresponding to the classical algebras $A_n$, $D_n$, $E_n$ [175]. Our case under consideration is related to the diagrams $D_n$, which describes the sine-Gordon model. To each node of the diagram we associate the particle with the mass given by the Coxeter number of the corresponding diagram (which is equal to $2K - 2$) shown in Fig. (17). These are denoted by $\mu_+ = \mu_-$ (corresponding to the kink and antikink) and $\mu_j$, $j = 1...2K - 2$, (note that we consider $K$ to be of the form $n/2$, $n = 3, 4, 5, ...$) corresponding to breathers, bound states of kink and antikink. Explicitly,

$$
\mu_j = 2\mu_+ \sin\left(\frac{j\pi}{4K - 2}\right) \quad \text{for breathers} \quad j = 1...2K - 2.
$$  \hspace{1cm} (C4)

![Dynkin diagram](image)

**FIG. 17:** Dynkin diagram for the algebra $D_n$. In the context of sine-Gordon problem nodes $1...2K - 2$ correspond to breathers, whereas nodes $\pm$ correspond to kink/antikink.

Due to a great simplification of the structure of scattering matrices in integrable theory [175], the free energy of the boundary sine-Gordon problem can be expressed entirely in terms of the energies $\epsilon_+$ (see Eq. (C2) above). This can further be reduced as follows. Consider the TBA equations for the energies:

$$
\epsilon_r(\theta) = \int_{-\infty}^{\infty} \frac{d\theta'}{2\pi \cosh[2(2K - 1)(\theta - \theta')]} \sum_s N_{rs} \log(1 + \exp(\epsilon_s(\theta'))),
$$  \hspace{1cm} (C5)

where $N_{rs}$ is the incidence matrix of the corresponding Dynkin diagram. To clarify these notations consider few examples.

For $K = 3/2$ we have kink, antikink and one breather. Therefore the diagram consists of the three nodes, say $+1, 2, -$. The vertex $+$ (kink) is connected to the node 1 and the vertex 1 is connected to the vertex $-$. Vertices + and $-$ are disconnected. The incidence matrix is therefore

$$
\begin{pmatrix}
1 & 0 \\
1 & 1 \\
0 & 1 \\
\end{pmatrix}
$$

(C6)

(the number of rows is equal to the number of vertices, the number of columns is equal to the number of bonds between these vertices). Using this matrix we obtain the following set of coupled TBA equations:

$$
\epsilon_\pm(\theta) = \int_{-\infty}^{\infty} \frac{d\theta'}{2\pi \cosh[2(2(2K - 1)(\theta - \theta'))]} \log(1 + \exp(\epsilon_1(\theta'))),
$$  \hspace{1cm} (C7)

$$
\epsilon_1(\theta) = 2 \int_{-\infty}^{\infty} \frac{d\theta'}{2\pi \cosh[2(2(2K - 1)(\theta - \theta'))]} \log(1 + \exp(\epsilon_+(\theta'))),
$$  \hspace{1cm} (C8)

where we use the fact that $\epsilon_+ = \epsilon_-$. Now, from the second equation we can see that the $\epsilon_1(\theta)/2$ gives precisely the function $\tilde{F}$ from the Eq. (C2). Therefore, for the case of $K = 3/2$,

$$
Z_0\left(\frac{3}{2}, g\right)_{\text{pert}} = \frac{1}{\sqrt{3}} \exp\left[\frac{1}{2} \epsilon_1(\alpha)\right],
$$

(C9)

where $\epsilon_1$ must be determined from the set of coupled equations (C7).

To continue, consider the case of $K = 4/2$. Here we have 2 breathers, kink, and antikink. The incidence matrix is

$$
\begin{pmatrix}
1 & 0 & 0 \\
1 & 1 & 1 \\
0 & 1 & 0 \\
0 & 0 & 1 \\
\end{pmatrix}
$$

(C10)
The set of TBA equations is therefore

$$\epsilon_\pm(\theta) = \int_{-\infty}^{\infty} \frac{d\theta'}{2\pi} \frac{3}{\cosh[3(\theta - \theta')]} \log(1 + \exp(\epsilon_2(\theta'))),$$

(C11)

$$\epsilon_1(\theta) = \int_{-\infty}^{\infty} \frac{d\theta'}{2\pi} \frac{3}{\cosh[3(\theta - \theta')]} \log(1 + \exp(\epsilon_2(\theta'))),$$

(C12)

$$\epsilon_2(\theta) = \int_{-\infty}^{\infty} \frac{d\theta'}{2\pi} \frac{3}{\cosh[3(\theta - \theta')]} \log(1 + \exp(\epsilon_1(\theta'))),$$

(C13)

$$+ 2 \int_{-\infty}^{\infty} \frac{d\theta'}{2\pi} \frac{3}{\cosh[3(\theta - \theta')]} \log(1 + \exp(\epsilon_1(\theta'))),$$

from which we conclude that the $\epsilon_2/3$ gives the function $\tilde{F}$ from the Eq. (C2). Therefore, for the case of $K = 4/2$,

$$Z_0(\frac{4}{2}, g)_{\text{pert}} = \frac{1}{\sqrt{4}} \exp\left[\frac{1}{3} \epsilon_2(\alpha)\right],$$

(C14)

where $\epsilon_2$ must be determined from the set of coupled equations (C11).

For $K = 5/2$ we can find that

$$Z_0(\frac{5}{2}, g)_{\text{pert}} = \frac{1}{\sqrt{5}} \exp\left[\frac{1}{2} (\epsilon_3(\alpha) - \epsilon_1(\alpha))\right],$$

(C15)

where $\epsilon_1, 3$ must be determined from the set of corresponding coupled equations (easy to write down explicitly).

For $K = 6/2$ we have

$$Z_0(\frac{6}{2}, g)_{\text{pert}} = \frac{1}{\sqrt{6}} \exp\left[\frac{1}{2} (\epsilon_4 - \epsilon_2 + \tilde{F}(\epsilon_1))\right],$$

(C16)

and therefore it includes the function $\tilde{F}(\epsilon_1)$. This is not convenient from the point of view of numerics since it contains an additional integration.

Continuing the above computations one can observe that for the values of $K = 2k/2$, $k = 2, 3, 4, \ldots$ we will always have an additional function $\tilde{F}(\epsilon_1)$ in the expression for the perturbative function $Z_0(g, K)_{\text{pert}}$. On the other hand for the values of $K = (2k + 1)/2$, $k = 1, 2, 3, \ldots$ one can always express the result for partition function in terms of energies $\epsilon_i$. It seems that this is more convenient for numerics. The result for this case is the following:

$$Z_0(g, \frac{2n+1}{2})_{\text{pert}} = \frac{1}{\sqrt{2n + 1}} \exp\left[\frac{1}{2} \sum_{s=0}^{n-1} (-1)^s \epsilon_{(2n-1) - 2s}\right],$$

(C17)

where the set of energies must be determined from the solution of coupled integral equations.

It is useful to consider the asymptotic behavior. For a large $\theta \to \pm \infty$ values of the integration parameter, the energies of breathers may be taken in the following form (see Eq. (C4)):

$$\epsilon_s(\alpha) = \frac{\mu_s}{\mu_+} = 2 \sin\left[\frac{s\pi}{4K - 2}\right] \exp[\alpha], \quad \text{as} \quad \alpha \to \infty.$$  

(C18)

The general strategy is clear now: one can solve numerically the set of coupled integral equations for some values of $K = (2n + 1)/2$ and compute the partition function using Eq. (C17). At this point the difficulty arises: how to unambiguously define the analytic continuation of $Z_p(K, g)$ into imaginary-$g$ domain. To avoid this problem it is better to use the alternative approach described in section [V C 1].

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Readers may be concerned that the initial part of expansion may not be purely ballistic and atomic collisions may take place in expanding clouds. In the case of tight confinement such collisions are rare and have minimal effect. One can use the moment of the release using magnetic Feshbach resonances [170]. See also discussion after Eq. (43).

[167, 168, 169]. It is worth noting that the effect of collisions can be minimized by tuning the scattering length to zero at the first stages of the expansion. Effects of the interactions on the interference signal have been recently discussed in Refs. 95, 98, 103. See also discussion after Eq. (43).
