Expectation Values in the Lieb–Liniger Bose Gas

M. Kornos,1 G. Mussardo,1,2 and A. Trombettoni1

1SISSA and INFN, Sezione di Trieste, via Beirut 2/4, I-34151, Trieste, Italy
2International Centre for Theoretical Physics (ICTP), I-34151, Trieste, Italy

Taking advantage of an exact mapping between a relativistic integrable model and the Lieb–Liniger model we present a novel method to compute expectation values in the Lieb–Liniger Bose gas both at zero and finite temperature. These quantities, relevant in the physics of one-dimensional ultracold Bose gases, are expressed by a series that has a remarkable behavior of convergence. Among other results, we show the computation of the three-body expectation value at finite temperature, a quantity that rules the recombination rate of the Bose gas.

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Correlation functions are key quantities in quantum interacting systems: not only they fully encode the dynamics but also are directly related to various susceptibilities and response functions. For these reasons, there has always been an intense search to find the most efficient ways to compute them. The task is notoriously difficult, even if the system is integrable. A typical but significant example is provided by the Lieb–Liniger (LL) model [1] that describes the low-temperature properties of one-dimensional interacting Bose gases: although it can be solved through Bethe Ansatz equations, the explicit computation of its correlation functions is a long-standing problem [2 3]. The interest for the computation of correlation functions in the LL model is obviously not only theoretical. In a series of recent experimental achievements strongly interacting ultracold bosons have been confined within wave-guides by nearly one-dimensional potentials that tightly trap the particle motion in the two transverse directions while leaving it free in the third axial direction [4 6]: the coupling of these bosons with the external world can be made so weak that their behaviour is very well described by the Lieb–Liniger model. Through interference (or eventually in situ) experiments, several quantities can be detected both at zero and finite temperature: the time duration of experiments depends on the three-body recombination rate, which is proportional to the external trapping potential [9].

Over the years several theoretical quantities of the LL model have been computed by means of different techniques [10–19]. In this paper we present a compact and general way to determine the expectation values of its local operators. The method takes advantage of an exact mapping between a relativistic integrable massive model – the sinh–Gordon (ShG) – and the LL model: in a proper non-relativistic limit of the ShG model, both its S-matrix and Lagrangian coincide with those of the LL model. Since the S-matrix of an integrable relativistic model fixes the exact matrix elements of all operators of the theory (and for the ShG model these matrix elements are all known [20 21]), the correspondence between the two models opens the way to computing the corresponding quantities of the LL model in a very direct way. As shown below, this method provides a remarkable simplification of the problem. Its implementation actually requires to take into account an additional aspect of the problem: while in the ShG model the correlation functions refer to the vacuum (i.e. the state without any particles), in the LL model they relate instead to its ground state at a finite density. This aspect, however, can be successfully overcome by the Thermodynamical Bethe Ansatz (TBA) formalism developed in [22], which has the additional convenience of being applicable equally well both at zero and finite temperature. In this way we are able to compute not only the zero temperature expectation values but also their finite temperature expressions.

The LL-ShG mapping. The LL Hamiltonian for N interacting bosons of mass m in one dimension is

$$H = -\frac{\hbar^2}{2m} \sum_{i=1}^{N} \frac{\partial^2}{\partial x_i^2} + 2\lambda \sum_{i<j} \delta(x_i - x_j).$$

The corresponding non-relativistic field theory description is the quantum non-linear Schrödinger model [2], which employs the complex field $\psi$ and the Lagrangian

$$\mathcal{L} = -\frac{\hbar^2}{2m} |\nabla \psi|^2 + i \frac{\hbar}{2} \left( \psi \frac{\partial \psi}{\partial t} - \frac{\partial \psi^\dagger}{\partial t} \psi \right) - \lambda |\psi|^4. \quad (2)$$

The effective coupling constant of the LL model is given by the dimensionless parameter $\gamma = 2m\lambda/\hbar^2 n$, where $\lambda > 0$ is the coupling entering the Hamiltonian [1] while $n = N/L$ is the density of the gas ($L$ is the length of the system). Temperatures are usually expressed in units of the temperature $T_D = \hbar^2 n^2 / 2mk_B$ of the quantum degeneracy, $\tau = T/T_D$. The two-body elastic S-matrix of the LL model is [11 12]

$$S_{LL}(p, \lambda) = \frac{p - i2m\lambda/\hbar}{p + i2m\lambda/\hbar}, \quad (3)$$

where $p = \hbar \sum_i \partial / \partial x_i$ is the linear momentum operator.
where $p$ is the momentum difference of the two particles.

Consider now the ShG model in (1+1) dimensions, i.e. the integrable and relativistic invariant field theory defined by the Lagrangian

$$L_{\text{ShG}} = \frac{1}{2} \left[ \left( \frac{\partial \phi}{c \partial \tau} \right)^2 - (\nabla \phi)^2 - \frac{\mu^2}{g^2} \cosh(g \phi) \right] ,$$

(4)

where $\phi = \phi(x, t)$ is a real scalar field, $\mu$ is a mass scale and $c$ is the speed of light. The parameter $\mu$ is related to the physical (renormalized) mass $m$ by $\mu^2 = \pi \alpha e^2 / (2 \hbar^2)$, where $\alpha = \frac{\hbar c g^2}{8 \pi + c g^2}$.

The energy $E$ and the momentum $P$ of a particle can be written as $E = mc^2 \cosh \theta$, $P = mc \sinh \theta$, where $\theta$ is the rapidity. Since the ShG dynamics is ruled by an infinite energy limit of a field theory consists of expressing the real scalar coupling as a function of the rapidity. It is easy to see that taking simultaneously the non-relativistic limit $\theta, \alpha \to \infty$.

$$S_{\text{ShG}}(\theta, \alpha) = \frac{\sinh \theta - i \sin(\alpha \pi)}{\sinh \theta + i \sin(\alpha \pi)} ,$$

(5)

where $\theta$ is the rapidity difference of the two particles. It is now easy to see that taking simultaneously the non-relativistic and weak-coupling limits of the ShG model such that

$$g \to 0, \ c \to \infty, \ g c = 4 \sqrt{\lambda} / \hbar = \text{fixed} ,$$

(6)

its $S$-matrix becomes identical to the $S$-matrix of the LL model. Notice that the coupling $\lambda$ does not need to be small, i.e. with this mapping we can study the LL model at arbitrarily large values of the dimensionless coupling $\gamma$.

The mapping between the two models goes beyond the identity of their $S$-matrix: it actually extends both to their Lagrangians and TBA equations. Details will be given elsewhere, but it is simple to follow the main steps of the procedure. According to eqn [24], the non-relativistic limit of a field theory consists of expressing the real scalar field in the form

$$\phi(x, t) = \sqrt{\frac{\hbar^2}{2m}} \left( \psi(x, t) e^{-i \frac{m x^2}{2 \hbar} t} + \psi^\dagger(x, t) e^{i \frac{m x^2}{2 \hbar} t} \right) ,$$

and, when the limit $c \to \infty$ of the Lagrangian is taken, of omitting all its oscillating terms. The commutation relation $[\phi(x, t), \Pi(x', t')] = i \hbar \delta(x - x')$ implies for the non-relativistic operators $[\psi(x, t), \psi^\dagger(x', t')] = \delta(x - x')$. Furthermore, when the limit $g \to 0$ of eqn [4] is considered, the $\psi^\dagger \psi$ terms coming from the potential and kinetic parts cancel each other, while all higher terms of the series expansion of the potential, but the quartic one, vanish. Hence, the ShG Lagrangian reduces to the non-linear Schrödinger Lagrangian.

In the same way one can also show that the TBA equations of the ShG model (given for instance in [10]) reduces to the ones of the LL model, written down in [10]. In the LL model at a finite $T$ the TBA equation for the pseudo-energy $\epsilon(T, \mu)$ consists of the non-linear integral equation

$$\epsilon(T, \mu) = \frac{p^2 / 2m - \mu}{k_B T} - \varphi \circ \log \left( 1 + e^{-\epsilon} \right) ,$$

(7)

where $\mu$ is the chemical potential associated to the finite density $n$ of the gas, $\varphi(p) = -i \frac{\partial}{\partial n \log S_{\text{LL}}(p)}$ is the derivative of the phase shift and $\varphi \circ f \equiv \int_{-\infty}^{\infty} \frac{dp'}{2 \pi} \varphi(p-p') f(p')$.

The solution of this integral equation leads to the free energy and to all other thermodynamical data of the model.

**Expectation values.** At equilibrium the expectation value of an operator $\mathcal{O} = \mathcal{O}(x)$ at temperature $T$ and at finite density is given by

$$\langle \mathcal{O} \rangle = \frac{\text{Tr} \left( e^{-(H-\mu N)/(k_B T)} \mathcal{O} \right)}{\text{Tr} \left( e^{-(H-\mu N)/(k_B T)} \right)} .$$

(8)

In a relativistic integrable model the above quantity can be neatly expressed as

$$\langle \mathcal{O} \rangle = \sum_{n=0}^{\infty} \frac{1}{n!} \int_{-\infty}^{\infty} \left( \prod_{i=1}^{n} \frac{d\theta_i}{2\pi} f(\theta_i) \right) \langle \hat{\theta} | \mathcal{O}(0) | \hat{\theta} \rangle_{\text{conn}} ,$$

(9)

where $f(\theta_i) = 1/(1 + e^{\theta_i})$ and $\hat{\theta} \equiv \theta_1, \ldots, \theta_n$ for $\langle \hat{\theta} | \mathcal{O}(0) | \hat{\theta} \rangle \equiv \langle \eta_1, \ldots, \eta_1 | \mathcal{O}(0) | \eta_1, \ldots, \eta_1 \rangle$ denote the asymptotic states entering the traces in [5]. This formula employs both the pseudo-energy $\epsilon(\theta)$ and the connected diagonal form factor of the operator $\mathcal{O}$, defined as $\langle \hat{\theta} | \mathcal{O} | \hat{\theta} \rangle_{\text{conn}} = \mathcal{F} \mathcal{P} \left( \lim_{\eta_1 \to 0} \langle 0 | \hat{\theta} - \theta - i \pi + i \eta_1 | \hat{\theta} \rangle \right)$ where $\eta_1 \equiv \eta_1, \ldots, \eta_1$ and $\mathcal{F} \mathcal{P}$ in front of the expression means taking its finite part, i.e. omitting all the terms of the form $\eta_1 \eta_j$ and $1/\eta^p$ where $p$ is a positive integer.

In order to compute the expectation values of the LL model by applying eqn [9] we need: (a) to solve the integral equation [7] for $\epsilon(\theta)$; (b) to identify the relevant form factors of the ShG model; (c) to take the non-relativistic limit of both the form factors and eqn [9]. Given for granted the straightforward (numerical) solution of eqn [7], let us focus our attention on the last two points. The generic $m$-particle form factor of a local operator $\mathcal{O}$ in the ShG model can be written as

$$\mathcal{F}_m(\theta_1, \ldots, \theta_m) = \mathcal{Q}_m(x_1, \ldots, x_m) \prod_{i<j} \frac{F_{\min}(\theta_{ij})}{x_i + x_j} ,$$

(10)

where $x_i = e^{\theta_i}$ and $\mathcal{Q}_m$ are the symmetric polynomials in the $x$’s that fully characterize the operator $\mathcal{O}$. The explicit expression of $F_{\min}(\theta)$ is given in [20] but the only thing needed here is its functional equation

$$F_{\min}(i \pi + \theta) F_{\min}(\theta) = \frac{\sinh \theta}{\sinh \theta + i \sin(\pi \alpha)} .$$

We are interested in the symmetric polynomials $\mathcal{Q}_m^{(\alpha)}$ of the exponentials $\mathcal{Q}_m = e^{\alpha \phi}$ since, using their Taylor expansion, we can extract the form factors of all normal
ordered operators $\phi^k$. Their expression is [21]
\[
Q_{m}^{(q)} = [q] \left( \frac{4 \sin(\pi \alpha)}{N} \right)^{\frac{q}{2}} \det M_m(q),
\]
where $M_m(q)$ is an $(m-1) \times (m-1)$ matrix with elements $[M_m(q)]_{i,j} = \sigma^{(m)}_{i,j} |i - j + q|$. Above, $[x] \equiv \sin(x \pi \alpha) / \sin(\pi \alpha)$ while $\sigma^{(m)}_a (a = 0, 1, \ldots, m)$ are the elementary symmetric polynomials in $m$ variables.

On the basis of the results given above, we are now in the position to compute the local $k$-particle correlation function $g_k$ of the LL model defined by
\[
\langle \psi^\dagger k \psi^k \rangle = n^k g_k(\gamma, \tau),
\]
where $k$ is an integer ($k = 1, 2, 3, \ldots$). The $g_k$’s are functions of the dimensionless LL coupling $\gamma$ and of the reduced temperature $\tau$. The relation between $g_k$ in the LL model and the corresponding quantity in the ShG model in the limit (6) is given by
\[
\langle \phi^{2k} \rangle \rightarrow \left( \frac{\hbar^2}{2m} \right)^k \left( \frac{2k}{k} \right) \langle \psi^\dagger k \psi^k \rangle.
\]

Using eqn (9) and the connected form factors of the corresponding operator we arrive at the expression
\[
\langle \psi^\dagger k \psi^k \rangle = \left( \frac{2k}{k} \right)^{-1} \left( \frac{\hbar^2}{2m} \right)^{-k} \sum_{n=1}^\infty F_n,
\]
\[
F_n = \frac{1}{n!} \int_{-\infty}^{\infty} \left( \prod_{i=1}^{2n} \frac{dp_i}{2\pi} f(p_i) \right) F^{\psi^k}_{2n,\text{conn}}(p_1, \ldots, p_n),
\]
where
\[
F^{\psi^k}_{2n,\text{conn}}(\{p_i\}) = \lim_{c \to \infty, g \to 0} \left( \frac{1}{mc} \right)^n F^{\psi^k}_{2n,\text{conn}}(\{\theta_i = \frac{p_i}{mc}\})
\]
is the double limit (6) of the connected form factors. As shown below, the series (13) are nicely saturated by the first few terms for sufficiently large values of $\gamma$ ($\gamma = 0$ is a singular point of the model [1], therefore one cannot expect a priori any fast convergence nearby). A first check of eqn (13) is provided by the case $k = 1$: using (13) (with a chemical potential $\mu$ that ensures the finite density $n$) and summing up the series, one easily checks that $\langle \psi^\dagger \psi \rangle = n$ and $g_1 = 1$, as it should be for translational invariance. As shown in Fig. 1 the exact value $g_1 = 1$ (solid line) is rapidly approached by just the first terms of (13): the convergence of the series is always remarkably fast for all $\gamma \geq 1.5$, where the exact value is obtained within a 5% accuracy just using its first four terms. As a second check of (13) let us show how we can easily recover the leading order of the strong coupling (i.e. large $\gamma$) expansion of all $g_k$: since this always comes from the first non-zero integral in the series (13), we get
\[
g_k = \frac{k!}{2k} \left( \frac{\pi}{\gamma} \right)^{k(k-1)} I_k + \ldots,
\]
where
\[
I_k = \int_{-\infty}^{\infty} dk_1 \ldots \int_{-\infty}^{\infty} dk_n \prod_{i,j} (k_i - k_j)^2.
\]
This result coincides with the one obtained in [12].

The quantity $g_2$ can be exactly determined via the Hellmann–Feynman theorem [13] and its plot at $T = 0$ is shown in Fig. 1 together with our determination from eqn (13). As before, also in this case there is a fast convergent behaviour of the series. The strong coupling regime of $g_2$ can be computed by expanding in powers of $\gamma^{-1}$ all the terms in eqn (13) and for $T = 0$ we get
\[
g_2 = \frac{4 \pi^2}{3 \gamma^2} \left( 1 - \frac{6}{\gamma} + (24 - \frac{8}{5} \pi^2) \frac{1}{\gamma^2} \right) + \mathcal{O}(\gamma^{-5}),
\]
in agreement with the result of the Hellmann–Feynman theorem [14]. Expression (15) is also plotted in Fig. 1 in order to show that the determination of $g_2$ (at finite $\gamma$) obtained from the first terms of eqn (13) is closer to the exact result, because any of them contains infinitely many powers of $\gamma$. At finite temperatures the convergence of the series is also pretty good and the results are shown in Fig. 2.

As a final example, let us discuss $g_3$, a quantity known exactly at $T = 0$ [19], but only approximately at $T > 0$.
The computation of the correlators of the LL model. Then, it is easy to take their non-relativistic limit. As shown in the previous examples, it shows a nice convergent pattern to the exact values found in [19]. Figs. 3(b,c) show a significant advantage in using this method instead of employing directly the sinh–Gordon model. There is a significant advantage of non-relativistic field theory presenting a larger number of constraints (crossing invariance, for instance) which permit to pin down exactly and efficiently the matrix elements of all operators: once these quantities are known, it is then easy to take their non-relativistic limit. As shown above, this proves to be a notable simplification in the computation of the correlators of the LL model.

Conclusion. We have shown that the equilibrium expectation values for one-dimensional interacting Bose gases can be efficiently computed by using the non-relativistic limit of an integrable relativistic field theory, the sinh–Gordon model. There is a significant advantage in using this method instead of employing directly the non-relativistic Lagrangian (2). The reason is that a relativistic field theory presents a larger number of constraints (crossing invariance, for instance) which permit to pin down exactly and efficiently the matrix elements of all operators: once these quantities are known, it is then easy to take their non-relativistic limit. As shown above, this proves to be a notable simplification in the computation of the correlators of the LL model.

The method works equally well at $T = 0$ and $T \neq 0$ where the series expansion presents a remarkable convergence behaviour for finite values of $\gamma$. There is no obstruction, in principle, to compute higher form factors and further improve the result. Strong coupling expansions in $\gamma^{-1}$ can be easily derived as well but at finite $\gamma$ the form factor expansion, containing infinitely many powers of $\gamma$, is more accurate for the determination of $g_3$, as we showed comparing it with exact results. As a significant application of the method we have determined $g_3$ at finite temperature (a term which is proportional to the recombination rate of the gas). This quantity, as well as the higher $g_4$, may provide important information once the integrability of the model is broken. In the future it would be also interesting to apply this method both to two-point correlation functions and to other models.

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