Semirelativistic $N$-boson systems bound by attractive pair potentials

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Abstract
We establish bounds on the energy of a system of $N$ identical bosons bound by attractive pair potentials and obeying the semirelativistic Salpeter equation. The lower bound is provided by a ‘reduction’, with the aid of Jacobi relative coordinates, to a suitably scaled one-body Klein–Gordon problem. Complementary upper energy bounds are provided by means of a Gaussian trial function. Detailed results are presented for the exponential pair potential $V(r) = -v \exp(-r/a)$.

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1. Introduction
We consider a system of $N$ identical bosons which interact by attractive central pair potentials $V(r_{ij})$. There is no restriction in general concerning the number $d$ of spatial dimensions, but, for definiteness in the presentation, we shall assume $d = 3$. We suppose that the potential has the form $V(r) = vf(r/a)$, where $r = |r|$, $v > 0$ is the coupling parameter, $f(r/a) < 0$ is the shape of the potential and $a > 0$ is a range parameter. The Salpeter semirelativistic Hamiltonian for a system of $N$ identical bosons of mass $m$ and individual momenta $p_i$ is given by

$$H = \sum_{i=1}^{N} \sqrt{p_i^2 + m^2} + \sum_{i=1}^{N} \sum_{i<j} Vf(r_{ij}/a),$$  \hspace{1cm} (1)

where $r_{ij} = |r_i - r_j|$ and $p_i = |p_i|$. Our principal goal is to find or estimate $E$, the lowest energy of this system. The ease with which the $N$-body problem can be posed in the Salpeter theory is the main reason for our adopting the semirelativistic approach. For suitable potentials,
the Hamiltonian is bounded below and to a certain extent the spectral theory mimics the corresponding Schrödinger theory [1–3].

One path to a lower bound is to exploit the boson permutation symmetry of the exact normalized ground state \( \Psi \) to effect a ‘reduction’ generated by the equation 
\[
E = (\Psi, H \Psi) = (\Psi, \mathcal{H} \Psi),
\]
where \( \mathcal{H} \) is the Hamiltonian for a specially scaled one-body problem. A lower bound to \( E \) is then given by the bottom of the spectrum of \( \mathcal{H} \). Such reductions have a long history which is reviewed briefly in [4]. Different reductions are possible and lead to a variety of lower bounds. For example, we obtain immediately from equation (1)
\[
E = (\Psi, H \Psi) = (\Psi, H_2 \Psi),
\]
where
\[
H_2 = \frac{N}{2} \left[ \sqrt{p_1^2 + m^2} + \sqrt{p_2^2 + m^2} + (N - 1)V(r_{12}) \right].
\]
The reduction is not yet complete since \( H_2 \) is the Hamiltonian for a scaled two-body problem.

If we adopt new coordinates \( R = r_1 + r_2 \) and \( r = r_1 - r_2 \) for this problem, we find \( p_1 = P + p \) and \( p_2 = P - p \). Thus, if we write \( \sqrt{p_1^2 + m^2} = |p_1 + mk|, i = 1, 2 \), where \( k \) is a unit vector orthogonal to \( p_1 \) and \( p_2 \), then an application of the triangle inequality yields
\[
2|p + mk| = |(p - P + mk) + (p + P + mk)| \leq |p_2 + mk| + |p_1 + mk|.
\]
Thus we have the spectral inequalities
\[
H \geq H_2 \geq \mathcal{H}_{N/2} = N\sqrt{p^2 + m^2} + \gamma v f(r/a) \geq \mathcal{E}_{N/2},
\]
where \( \gamma = \frac{1}{2}N(N - 1) \) and \( \mathcal{E}_{N/2} \) is the bottom of the spectrum of \( \mathcal{H}_{N/2} \) and provides a lower bound to \( E \), the ground-state energy of \( H \). We note that this use of the triangle inequality is equivalent to setting the centre-of-mass momentum \( P \) of the two-body Hamiltonian \( H_2 \) to zero.

For boson systems there may be an energy advantage (leading to a higher lower bound) if, from the outset, Jacobi (orthogonal) relative coordinates can be accommodated by the reduction scheme. Intuitively, the use of relative coordinates allows us to remove the centre-of-mass kinetic energy of the whole system, rather than dropping the centre-of-mass energy only at the final stage, after the reduction to a scaled two-body problem such as \( H_2 \). One way of achieving this for the Salpeter problem involves squaring the reduced one-body Salpeter equation of the form \( \sqrt{p^2 + m^2} \psi = (E - V) \psi \). This allows us to invoke some remarkable expectation identities for Jacobi coordinates: these identities, which we shall refer to explicitly below, are proved in appendix A.

Since we use relative coordinates both for the upper and lower bounds, in section 2 we define the Jacobi coordinates \( \{\rho_i\}_{i=2}^N \) and note some key properties. An upper energy bound \( E_U = (\Phi, H \Phi) \) is provided by an application of a normalized Gaussian trial function \( \Phi \). This latter calculation is greatly simplified by the boson permutation symmetry of \( \Phi \) and also by the fact that this particular trial function is not only symmetric under the permutation of the relative coordinates but also factors in the form \( \Phi = \prod_{i=2}^N \phi(\rho_i) \). The upper bound is discussed in section 3. In section 4 we introduce a variant of the Jacobi system suitable for the derivation of a lower bound expressed in terms of a Klein–Gordon one-particle problem. The proof assumes that this Klein–Gordon problem has a discrete eigenvalue that is either (i) positive or (ii) monotone increasing as a function of the mass. In section 5 we consider the example of the attractive exponential potential.
2. Relative coordinates

Since boson permutation symmetry is naturally expressed in terms of the individual-particle coordinates \( \{r_i\}_{i=1}^N \), it might seem an unwelcome complication to introduce relative coordinates \( \{\rho_i\}_{i=2}^N \). Their advantage arises when we consider the reduction \( \langle H \rangle = \langle \gamma \rangle \), where \( \gamma \) is a scaled one-body problem. Jacobi relative coordinates may be defined with the aid of an orthogonal matrix \( B \) transforming old \( \{r_i\} \) to new \( \{\rho_i\} \) coordinates by \( \rho_i = B[r_i] \). The first row of \( B \), with all entries \( B_{1i} = 1/\sqrt{N} \), defines a centre-of-mass variable \( \rho_1 \), its second row defines a pair distance \( \rho_2 = (r_1 - r_2)/\sqrt{2} \), and the \( k \)-th row (\( k \geq 2 \)) first has \( k - 1 \) entries \( B_{ki} = 1/\sqrt{K(k-1)} \), the \( k \)-th entry \( B_{kk} = -\sqrt{(k-1)/k} \), and zero for all remaining entries. The momenta \( \{\pi_i\} \) conjugate to \( \{\rho_i\} \) read \( \{\pi_i\} = (B^{-1})^T[\pi] = B[\rho] \). Now, for an \( N \)-boson problem with an attractive potential \( V(r) \), let \( \Psi(\rho_2, \rho_3, \ldots, \rho_N) \) be the exact normalized ground-state eigenfunction corresponding to the lowest energy \( E \). Boson permutation symmetry is a powerful constraint that greatly reduces the complexity of this problem. Although a non-Gaussian wavefunction is not necessarily symmetric in the Jacobi coordinates, we do have the remarkable \( N \)-representability expectation expressions from appendix A: for \( i, j > 1 \)

\[
\langle \Psi, (\rho_i \cdot \rho_j) \Psi \rangle = \delta_{ij} \langle \Psi, \rho_i^2 \Psi \rangle, \quad (5)
\]

\[
\langle \Psi, (\pi_i \cdot \pi_j) \Psi \rangle = \delta_{ij} \langle \Psi, \pi_i^2 \Psi \rangle. \quad (6)
\]

3. A Gaussian upper bound

We now take advantage of the special properties of the Gaussian trial function \( \Phi \) to effect an upper reduction of \( \langle H \rangle \). In the centre-of-mass frame of the \( N \)-body system, where \( \pi_1 = 0 \), we have from the last row of the matrix \( B \)

\[
\rho_N = \frac{1}{\sqrt{N}} \pi_1 - \sqrt{\lambda} \pi_N = -\sqrt{\lambda} \pi_N, \quad \lambda = \frac{N-1}{N}. \quad (7)
\]

Meanwhile our trial wavefunction \( \Phi \) is not only a boson function but it is also symmetric and factors in the relative coordinates. We have explicitly

\[
\Phi(\rho_2, \rho_3, \ldots, \rho_N) = \prod_{i=2}^N \phi(\rho_i), \quad \phi(r) = \left( \frac{1}{\pi \sigma^2} \right)^{\frac{1}{2}} e^{-\frac{1}{2} r^2 / (\sigma^2)},
\]

where \( \sigma > 0 \) is a variational scale parameter. The boson permutation symmetry of the trial function allows us to write \( E \leq E_g = \langle \Phi, H \Phi \rangle \), where we have

\[
E_g = \langle \Phi, \left[ N \sqrt{\lambda \pi_1^2 + m^2 + \gamma V(\sqrt{2} \rho_2)]} \right] \Phi \rangle.
\]

Setting \( r = r_1 - r_2 = \sqrt{2} \rho_2 \) and \( p = \pi_2 / \sqrt{2} \), we find \( E \leq E_g = \langle \phi, \mathcal{H}_g \phi \rangle \), where

\[
\mathcal{H}_g = N \sqrt{2 \lambda p^2 + m^2 + \gamma V(r)}.
\]

We now introduce some new parameters which will allow us to express both the upper and lower energy bounds in compatible succinct forms. With \( V(r) = vf(r/a), \lambda = (N-1)/N \) and \( \gamma = \frac{1}{2} N (N-1) \), we now define

\[
\mu = ma/(2\lambda)^{\frac{3}{4}}, \quad e = E a/(2 \gamma^{\frac{1}{2}}), \quad e_g = E_g a/(2 \gamma^{\frac{1}{2}}), \quad v = \gamma^{\frac{1}{2}} (v a / 2).
\]

The Gaussian upper bound then becomes

\[
e \leq e_g = \min_{\sigma > 0} (\phi, \mathcal{H}_g \phi), \quad \text{where} \quad h_g = \sqrt{p^2 + \mu^2 + vf(r)}.
\]

(8)
We note parenthetically here that, if the same scaling is used for the reduced Hamiltonian \( \hat{\mathcal{H}}_{N/2} \) of equation (4), the result \( h_2 \) is comparable with \( h_\varepsilon \) and is given by
\[
h_2 = \sqrt{p^2/(2\lambda) + \mu^2 + \nu f(r)}, \quad \lambda = (N - 1)/N. \tag{9}
\]
We shall use this later in the paper to show that the \( N/2 \) bound is weak: we shall consider
\[
e_{2\varepsilon} = \min_{\sigma > 0} \langle \phi, h_2 \phi \rangle \geq e_{2\varepsilon}, \tag{10}
\]
where \( e_{2\varepsilon} \) is the bottom of the spectrum of \( h_2 \), and observe (in particular cases) that \( e_{2\varepsilon} < e_k \), where \( e_k \) is the Klein–Gordon lower bound that we shall derive in the next section; hence, for these problems, we know, without calculating \( e_{2\varepsilon} \), that \( e_{2\varepsilon} < e_k \).

However, our main point here is that the upper energy bound can be written in terms of an optimized expectation (with respect to a Gaussian wavefunction) of a special one-particle Salpeter problem with Hamiltonian \( h_\varepsilon \). The kinetic-energy expectation value may be expressed in terms of the modified Bessel function of the second kind \( K_1(x) \) \[7\]. Evaluating the integral, for convenience, in momentum space, yields
\[
(\phi, \sqrt{p^2 + \mu^2} \phi) = \frac{\mu}{\sqrt{2}\sigma} I(x),
\]
where \( \sigma = \sqrt{\frac{2\pi}{\mu}} \) and the integral \( I(x) \) is defined by
\[
I(x) = 4 \sqrt{\frac{\pi}{2}} \int_0^\infty \sqrt{2x + t^2} e^{-t^2} dt = \frac{2}{\sqrt{\pi}} x e^{x} K_1(x).
\]
Meanwhile, the potential-energy expectation becomes
\[
(\phi, f(r)\phi) = J \left( \frac{\sqrt{2x}}{\mu^2} \right),
\]
where the integral \( J(y) \) is given by
\[
J(y) = 4 \sqrt{\frac{\pi}{2}} \int_0^\infty e^{-t^2} f(\sqrt{y}t) t^2 dt.
\]
We now use \( s = 1/\sigma = \frac{\mu}{\sqrt{2}\pi} \) as a variational parameter and we obtain
\[
e \leq e_\varepsilon = \min_{s > 0} \left[ s I \left( \frac{4}{2 s^2} \sigma^2 \right) + v J \left( \frac{1}{s} \right) \right]. \tag{11}
\]

4. The lower energy bound

In order to effect a new reduction that incorporates \( \pi_1 = 0 \), we first choose to use \( p_N \) and \( p_{N-1} \) and we obtain the reduction
\[
E = (\Psi, H\Psi) = \left( \Psi, \left[ \frac{N}{2} \sqrt{p_N^2 + m^2} + \frac{N}{2} \sqrt{p_{N-1}^2 + m^2} + \gamma V(|r_{N-1} - r_N|) \right] \Psi \right),
\]
where \( \gamma = \frac{1}{2} N(N - 1) \). We note that the last column of \( B \) (which defines \( \pi_N \)) is given by
\[
[b, b, b, \ldots, b, -(N - 1)b], \quad \text{where} \quad b = 1/\sqrt{N(N - 1)}.
\]
Now we express the equation for \( E \) in terms of new coordinates defined by the following relations:
\[
\begin{align*}
r &= r_{N-1} - r_N = \alpha p_N - \beta p_{N-1}, & r &= |r|, \\
\begin{bmatrix} \mathcal{R} \\ r \end{bmatrix} &= \begin{bmatrix} \beta & \alpha \\ \alpha & -\beta \end{bmatrix} \begin{bmatrix} p_N \\ p_{N-1} \end{bmatrix}, & \begin{bmatrix} P \end{bmatrix} &= \frac{1}{2} \begin{bmatrix} \beta & \alpha \\ \alpha & -\beta \end{bmatrix} \begin{bmatrix} \pi_N \\ \pi_{N-1} \end{bmatrix},
\end{align*}
\]
where
\[ \alpha = \sqrt{\frac{N}{N-1}} > 1, \quad \beta = \sqrt{\frac{N-2}{N-1}} < 1, \quad \delta = \sqrt{\frac{N-2}{N}} < 1. \]

Consequently,
\[ \alpha^2 + \beta^2 = 2, \quad b^2 + \beta^2 = \frac{1}{\alpha^2} = \frac{N-1}{N}, \]
and
\[ (N-1)b = \alpha^{-1}, \quad Nb = \alpha, \quad \delta = \frac{\beta}{\alpha}, \quad 1 + \delta^2 = 2\lambda. \]

We note that because of the boson permutation symmetry, and after we set the centre-of-mass momentum variable \( \pi_1 = 0 \), we generally have that
\[ \langle \kappa(p_{N-1}) \rangle = \langle \kappa(p - \delta P) \rangle = \langle \kappa(p + \delta P) \rangle = \langle \kappa(p_N) \rangle, \]
where \( \kappa(p) \) is any appropriate kinetic energy function. The expression for the energy now has the form
\[ E = \langle \mathcal{H} \rangle, \]
where
\[ \mathcal{H} = N \sqrt{(p + \delta P)^2 + m^2 + \gamma V(r)}. \]  
(12)

The Hamiltonian \( \mathcal{H} \) is bounded below by the simple bound \( \mathcal{H}_{N/2} \) of equation (4). This result is a consequence of the inequality
\[ 2\langle |p + mk| \rangle = 2\langle (p - \delta P + mk) + (p + \delta P + mk) \rangle \leq 2\langle |p + \delta P + mk| \rangle. \]

We now consider the eigenvalue \( \mathcal{H}\psi = E\psi \), where \( \psi(r, R) \in \mathcal{D} \subset L^2(R^6) \) retains some of the boson permutation symmetry implications of the full wavefunction, namely from equation (6), \( \langle p^2 \rangle = \langle P^2 \rangle, \langle p \cdot P \rangle = 0 \), and hence
\[ \langle (p + \delta P)^2 \rangle = \langle p^2 + \delta^2 P^2 \rangle = (1 + \delta^2)\langle p^2 \rangle = 2\lambda\langle p^2 \rangle. \]

Let \( \mathcal{E} \) be the minimum of \( \langle \psi, \mathcal{H}\psi \rangle \) corresponding to normalized \( \psi \) satisfying (13) and let \( \Psi \) be the exact normalized \( N \)-body wavefunction; then
\[ E = \langle \Psi, H\Psi \rangle = \langle \psi, \mathcal{H}\psi \rangle \geq \langle \psi, \mathcal{H}\psi \rangle = \mathcal{E}. \]

The eigenvalue equation for \( \mathcal{E} \) reads
\[ N \sqrt{(p + \delta P)^2 + m^2} \psi = (\mathcal{E} - \gamma v f(r/a)) \psi. \]
(14)

Squaring these equal vectors and use of (13), and also the scaling change (to new dimensionless variables) \( \{r, p\} \rightarrow \{ra, p/a\} \), implies
\[ (\mathcal{E}a)^2 - N^2(ma)^2 \geq 4\gamma \inf_{r \in D} \psi, \left( p^2 + \frac{(va)(\mathcal{E}a)}{2} f(r) - \frac{\gamma(va)^2}{4} f^2(r) \right) \psi \].  
(15)

In contrast to our upper bound, where we worked directly with the Salpeter Hamiltonian (8), for the lower bound we are led to consider a complementary lower Klein–Gordon operator. If the minimization on the right-hand side of (15) could be effected exactly (as it can for the gravitational problem \( f(r) = -1/r \) [5]), then it would remain to extract a lower bound to \( \mathcal{E} \) from the resulting inequality. For the general problem that we are considering here, we have to find another approach. To this end we have explored what happens if we set \( \mathcal{E} = \mathcal{E}_1 \) in equation (15), and solve the resulting Klein–Gordon equation for the ground-state energy \( e \). This Klein–Gordon equation is given explicitly by
\[ ((\mathcal{E}_1a)^2 - N^2(ma)^2)\phi = 4\gamma \left( p^2 + \frac{(va)(\mathcal{E}_1a)}{2} f(r) - \frac{\gamma(va)^2}{4} f^2(r) \right) \phi. \]
(16)
This exercise is only useful if an a priori relation can be established between $\mathcal{E}_k$ and $\mathcal{E}$. More particularly, we shall now establish a relationship between $\mathcal{E}_k$ and the lower bound to $\mathcal{E}$ provided by the inequality given in equation (15). We employ a similar parameter set to that for the upper bound, that is to say $\lambda = (N-1)/N$ and $\gamma = \frac{1}{2}N(N-1)$, and also

$$\mu = ma/(2\lambda)^{1/2}, \quad e_k = \mathcal{E}_k a/(2\gamma^{1/2}), \quad \nu = \gamma^{1/2}va/2.$$  

(17)

We thus arrive at the Klein–Gordon equation in a standard form, namely

$$(p^2 + 2e\nu f(r) - (\nu f(r))^2)\phi = F(e_k)\phi = (e^2 - \mu^2)\phi.$$  

(18)

and, equivalently, in Schrödinger form

$$(p^2 + 2\epsilon_k \nu f(r) - (\nu f(r))^2)\phi = F(e_k)\phi = (\epsilon^2 - \bar{\mu}^2)\phi.$$  

(19)

The spectral function $F(e_k)$ in equation (19) describes, for each fixed coupling $\nu$, how the lowest eigenvalue of the Schrödinger operator $p^2 + 2\epsilon_k \nu f(r) - (\nu f(r))^2$ depends on the energy parameter $\epsilon_k$; the Klein–Gordon eigenvalue exists if there is a solution of the equation $F(e_k) = e^2 - m^2$; these curves are shown in figure 1 for the case of an attractive exponential potential.

Similarly, the original lower-bound inequality (15) becomes an equality in terms of the parameters (17) and also a new mass parameter $\bar{\mu} \geq \mu$ used here as a ‘slack variable’. Thus, we have

$$(p^2 + 2e\nu f(r) - (\nu f(r))^2)\phi = F(e)\phi = (e^2 - \bar{\mu}^2)\phi.$$  

(20)
In earlier work [6], on the semirelativistic and Klein–Gordon one-body problems, we have proved that, under appropriate assumptions, the Klein–Gordon energy is monotone in the mass. More specifically, in our application, we can state this result in terms of equation (20), namely, for each fixed \( \nu \), if either \( e > 0 \), or \( e \leq 0 \) and \( e > \frac{1}{2} F'(e) \), then \( e'(\bar{\mu}) > 0 \). Thus, under these conditions (which must be verified for each potential), we have from equations (19) and (20) that \( e \geq e_k \). Thus \( E_k \) provides us with a lower bound to the \( N \)-body ground-state energy \( E \). In terms of the dimensionless energy parameters, our principal results are summarized by the following inequalities:

\[
e_k(\mu, \nu) \leq e \leq e_g(\mu, \nu),
\]

where \( e_k \) is given by solving the Klein–Gordon equation (19), \( e_g \) is given by the variational equation (11) and the parameters are defined in equation (17). Thus, whenever a solution exists, we necessarily have \( F'(e) < 2e \). For example, in the case of the exponential pair potential we show this relationship explicitly in figure 1.

5. The exponential potential

We now consider the example with potential shape in dimensionless form given by \( f(r) = -e^{-r} \). The lower bound \( e_k(\mu, \nu) \) is given by solving the explicit one-particle Klein–Gordon equation

\[
(p^2 + \mu^2)\psi = (e_k + \nu e^{-r})^2 \psi
\]

and the upper bound \( e_g(\mu, \nu) \) is given by

\[
e_g = \min_{s > 0} \left[ s I \left( \frac{\mu}{2s} \right) + \nu J \left( \frac{1}{s} \right) \right],
\]

where

\[
I(x) = \frac{4}{\sqrt{\pi}} \int_0^\infty \sqrt{2x + t^2} e^{-t^2} t^2 dt = \frac{2}{\sqrt{\pi}} x e^x K_1(x),
\]

and

\[
J(y) = -\frac{4}{\sqrt{\pi}} \int_0^\infty e^{-t^2} e^{-yt} t^2 dt = \frac{y}{\pi^{\frac{3}{2}}} - \frac{1}{2} (2 + y^2)(1 - \text{erf}(y/2)) e^{y^2/4}.
\]

For the choice \( \mu = 1 \) we exhibit some numerical results in table 1 and figure 2. We note that the mass parameter \( \mu \) decreases ‘slowly’ with increasing \( N \geq 2 \) and satisfies \( ma \geq \mu \geq ma/\sqrt{2} \). In table 1 we include values of \( e_{2g} \) calculated from equation (10) for \( N \to \infty \), or \( \lambda \to 1 \): since \( e_{2g} < e_k \), we know that the Klein–Gordon bound \( e_k \) is better than the \( N/2 \) bound, that is to say \( e_2 < e_{2g} < e_k \). Moreover, we see from table 1 that for large \( N \), the Klein–Gordon lower bound is in fact much better than the \( N/2 \) bound.

The ultrarelativistic limit \( m \to 0 \), which is perhaps meaningful in the spinless Salpeter theory, has no useful counterpart expressed by the Klein–Gordon equation. The Salpeter Hamiltonian in this limit (and for the exponential example) becomes \( H = |p| - \nu e^{-r} \): this problem is not very different qualitatively from the corresponding Schrödinger problem \( H = p^2 - \nu e^{-r} \), which also has a discrete eigenvalue for every \( \nu > 0 \) not too small. By contrast, the Klein–Gordon equation (22) with \( \mu = 0 \) implies \( e_k = 0 \), and leads to the
Figure 2. Lower (L) and upper (U) energy bounds to the ground-state energy $e$ of the $N$-boson system with pair-potential shape $f(r) = -\exp(-r)$. Dimensionless units $\mu = ma/(2\lambda)^{\frac{1}{2}} = 1$, $e = Ea/(2\gamma^2)$, $\nu = \gamma^2 va/2$, are used, in which $\lambda = (N-1)/N$, and $\gamma = \frac{1}{2}N(N-1)$. The lower and upper data are given, respectively, by equations (22) and (23).

Table 1. Energy bounds $e_k \leq e \leq e_g$ for the $N$-boson system bound by an attractive pair potential with shape $f(r) = -\exp(-r)$. The parameters are given by $\mu = ma/(2\lambda)^{\frac{1}{2}} = 1$, $e = Ea/(2\gamma^2)$, $\nu = \gamma^2 va/2$, $\lambda = (N-1)/N$ and $\gamma = \frac{1}{2}N(N-1)$. The energy $e_{2g}$ is an upper bound to the $N/2$ lower bound $e_2$. The table shows $e_{2g}$ for $N \to \infty$. Since $e_{2g} < e_2$, we know that the Klein–Gordon bound $e_2$ is better than the $N/2$ bound, that is to say, for large $N$, $e_2 < e_{2g} < e_k$.

| $\nu$ | $e_{2g}$ | $e_2$ | $e_g$ |
|-------|---------|-------|-------|
| 0.6   | 0.993 207 | 0.996 825 |       |
| 0.8   | 0.962 091 | 0.980 384 | 1.002 270 |
| 1     | 0.917 724 | 0.967 625 | 0.992 897 |
| 1.1   | 0.895 169 | 0.891 979 | 0.928 149 |
| 1.5   | 0.766 126 | 0.754 224 | 0.802 883 |
| 2     | 0.573 703 | 0.580 890 | 0.642 092 |
| 2.5   | 0.352 544 | 0.380 273 | 0.454 366 |
| 3     | 0.109 627 | 0.157 791 | 0.245 264 |
| 3.5   | -0.150 549 | 0.082 440 | 0.018 650 |
| 4     | -0.424 867 | -0.609 771 | -0.476 508 |
| 5     | -1.007 410 | -0.993 110 | -0.833 505 |
| 5.67  | -0.904 766 |       |       |
| 6     | -1.015 560 |       |       |

Schrödinger problem $(p^2 - v^2 e^{-2r})\psi = 0$, which has a solution only for very particular $v$ not too small. Clearly, the Klein–Gordon lower bound is not useful for the ultrarelativistic limit.
6. Supercriticality and system size

We consider first a one-body problem with an attractive (vector) potential \( V(r) = vf(r) \) in three spatial dimensions. We contrast four theories: Dirac, Klein–Gordon, Salpeter and Schrödinger. For the singular Coulomb or gravitational case \( f(r) = -1/r \), all four theories agree that binding is obtained for arbitrarily small \( v \). Meanwhile, the three relativistic theories agree that the coupling \( v \) cannot be very large: we have for Dirac \( v < 1 \), for Klein–Gordon \( v < 1/2 \) and for Salpeter \( v < 2/\pi \). For these three relativistic theories, the discrete eigenvalues are all positive and the supercritical phenomenon that we discuss below does not arise. Nonetheless, the coupling restrictions lead to limitations on the system size, as we have discussed earlier for a static-gravity boson-star model [4].

We now set aside the Coulomb case and consider bounded short-range potentials such as the exponential potential \( f(r) = -e^{-r} \) discussed in the previous section. For this class of problems, the four theories now agree that (in three dimensions) binding is obtained only for \( v \) sufficiently large. The term ‘critical coupling’ is used for the coupling value \( v_c \), in each theory, beyond which a discrete eigenvalue appears. As \( v \) is increased further, this eigenvalue decreases. In the fully relativistic theories, another critical value \( v = v_s \) is reached at which \( E = -m \). This situation is called ‘supercriticality’ and \( v_s \) is the corresponding supercritical coupling. The relativistic Dirac and Klein–Gordon theories will not accommodate larger couplings than their particular supercritical values. By contrast, for the Salpeter and Schrödinger cases, as \( v \) is increased, the eigenvalues simply decrease without lower limit. It is now clear that if we use the Klein–Gordon eigenvalue to bound the Salpeter eigenvalue below, a limitation is introduced which was not originally present in the Salpeter problem itself, namely that the coupling should not be very large. One could perhaps argue that, just as the restriction \( E > -m \), which emerges internally and technically from the theory of the Klein–Gordon equation, is ‘understood’ in terms of another theory, namely pair creation in quantum field theory; thus, the significance of this constraint could be acquired by the Salpeter theory from the Klein–Gordon theory. All we are saying here is that, if the Salpeter theory claims to be partially relativistic, then it is natural to impose the condition that \( E > -m \), or, what is sufficient, that \( v \) not be too large. Using the Klein–Gordon theory to bound Salpeter below achieves this automatically. In terms of the \( N \)-body reductions discussed in this paper, this means that the scaled coupling \( v = \gamma^2va/2 \) should not be too large. Since \( \gamma = \frac{1}{2}N(N-1) \), this means, for a given pair interaction, that the particle number \( N \), in turn, should not be too large. Thus we arrive at the conclusion that such relativistic systems in their ground states will not become giant balls as \( N \) is increased, but will divide into smaller aggregates. For the exponential example with \( m = 1, a = \sqrt{2} \), \( v \) small and \( N \) large so that \( \lambda \approx 1 \), we have from table 1, \( v = \gamma^2va/2 < 5.7 \), which is to say, \( N < 11.4/v \). Thus we conclude that a phenomenon which seemed to be a special feature of the \(-1/r\) potential may be more widely present in relativistic many-body systems.

7. Conclusion

The principal disadvantage of the semirelativistic Salpeter equation is that it is not fully relativistic. It perfectly well accommodates the harmonic oscillator, a strong interaction not at all tolerated (as a vector potential) by the Klein–Gordon or Dirac theories. Moreover, the Salpeter Hamiltonian is non-local, and this introduces technical problems for spectral estimation. The main advantages of the Salpeter theory are (i) the ease with which the relativistic many-body problem can be formulated and (ii) the fact that the energy spectrum is characterized, and may therefore be estimated, variationally. For a wide class of attractive pair
potentials that vanish at infinity, we have shown in this paper that Jacobi relative coordinates and
the permutation symmetry of the boson states allow us to construct ground-state energy bounds
in terms of certain reduced one-body problems: the lower bound is provided by the ground-
state energy of a Klein–Gordon problem, and the upper bound is given by the expectation of a
Salpeter Hamiltonian with respect to a scale-optimized Gaussian trial function. These bounds
may well have theoretical as well as computational applications. For example, as \( N \) increases,
the effective potential coupling \( \nu \) in the scaled one-body Klein–Gordon problem (18) increases
by the factor \( \sqrt{N(N - 1)} \sim N \) whereas the kinetic energy term remains unchanged. Even
for bounded potentials, such as the exponential potential, it implies that this lower-bound
Klein–Gordon problem will, for sufficiently large \( N \), become supercritical, with \( e < -\mu \).
This may, in turn, indicate that the \( N \)-body problem will not be stable for large aggregates,
and will necessarily break up into smaller fragments. This would be the expression of a truly
relativistic many-body phenomenon with no pre-image in non-relativistic theories.

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Appendix A

For definiteness, we consider the relative momenta \( \pi_i \) and \( \pi_j \) with \( i, j > 1 \). Since \( [\pi] = B[p] \),
we have

\[
\pi_i \cdot \pi_j = \left( \sum_k B_{ik} p_k \right) \cdot \left( \sum_k B_{jk} p_k \right) = \sum_k B_{ik} B_{jk} \left( p_k^2 \right) + \sum_{k \neq l} B_{ik} B_{jl} \left( p_k \cdot p_l \right). \tag{A.1}
\]

By using the boson symmetry of the wavefunction we have

\[
\langle \pi_i \cdot \pi_j \rangle = \left( \sum_k B_{ik} B_{jk} \right) \langle p_k^2 \rangle + \left( \sum_{k \neq l} B_{ik} B_{jl} \right) \langle p_1 \cdot p_2 \rangle. \tag{A.2}
\]

But orthogonality of the rows of \( B \) to the first row tells us that

\[
0 = \sum_k B_{ik} = \left( \sum_k B_{ik} \right) \left( \sum_l B_{il} \right) = \left( \sum_k B_{ik} B_{jk} \right) + \left( \sum_{k \neq l} B_{ik} B_{jl} \right). \tag{A.3}
\]

If \( i = j \), then the orthogonality of the matrix \( B \) tells us that

\[
\sum_k B_{ik}^2 = 1 = \sum_{k \neq l} B_{ik} B_{jl}.
\]

Meanwhile, if \( i \neq j \), we know that the corresponding rows of \( B \) are orthogonal; hence, in this
case, \( \sum_k B_{ik} B_{jk} = 0 \); thus, both of the coefficients in (A.2) are zero. Hence we conclude for
all \( i, j > 1 \)

\[
\langle \pi_i \cdot \pi_j \rangle = \delta_{ij} \langle p_i^2 \rangle = \delta_{ij} \left( \langle p_i^2 \rangle - \langle p_1 \cdot p_2 \rangle \right). \tag{A.4}
\]

The proof for the relative coordinates \( \{ \rho_i \} \) is identical.
References

[1] Salpeter E E and Bethe H A 1951 Phys. Rev. 84 1232
[2] Salpeter E E 1952 Phys. Rev. 87 328
[3] Lieb E H and Loss M 1996 Analysis (New York: American Mathematical Society)
[4] Hall R L and Lucha W 2008 J. Phys. A: Math. Theor. 41 355202
[5] Hall R L and Lucha W 2006 J. Phys. A: Math. Gen. 39 11531
[6] Hall R L and Lucha W arXiv:0906.2001 [math-ph]
[7] Abramowitz M and Stegun I A (ed) 1972 Handbook of Mathematical Functions (New York: Dover)