

# Semiclassical Trace Formulas for Two Identical Particles

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Semiclassical periodic orbit theory is used in many branches of physics. However, most applications of the theory have been to systems which involve only single particle dynamics. In this work, we develop a semiclassical formalism to describe the density of states for two noninteracting particles. This includes accounting properly for the particle exchange symmetry. As concrete examples, we study two identical particles in a disk and in a cardioid. In each case, we demonstrate that the semiclassical formalism correctly reproduces the quantum densities.

## I. INTRODUCTION

Semiclassical physics has experienced a resurgence of interest, largely due to the work of Gutzwiller [1], Balian and Bloch [2] and Berry and Tabor [3]. (For recent reviews see [4,5].) These works showed that if we separate the density of states into smooth and oscillatory components, then the oscillatory part is related to the dynamics of the underlying classical system via periodic orbits. This complements the earlier work of Weyl, Wigner, Kirkwood and others who showed that the smooth component is related to the geometry of the classical phase space. Actually, the two components are related in a subtle way [6,7] since the complete geometry imparts the full dynamics and vice-versa.

Most of the theoretical work has concentrated on the single particle density of states, however, there are some notable exceptions, namely [8–10]. In [8] the focus is on the average level density and its extension to systems of identical particles. Specifically, the authors consider a system of  $N$  fermions in one dimension. Their Weyl formula for fermions works well for attractive two body interactions, but overestimates the quantum staircase function when there are repulsive two body interactions. The author of [9] develops a generalization of the canonical periodic orbit sum for the special case of  $N$  interacting spinless fermions in one dimension. It is assumed the periodic orbits are isolated and therefore it is most applicable to fully chaotic systems. The author also considers a system of noninteracting fermions and writes the many body level density as a convolution integral involving one body level densities. Finally, we mention [10] which presents an expansion of the periodic orbit sum in terms of the particle number using ideas from [8,9].

Similarly, most of the applications of semiclassical theory have been to systems which involve only single particle dynamics. Here, we mention some exceptions. The authors of [11] extend the study of scars [12] to classically chaotic few body systems of identical particles. A study of the eigenfunctions of an interacting two particle system can be found in [13]. The semiclassical approach to the helium atom, which can be understood as two inter-

acting electrons in the presence of a helium nucleus, has been studied in [14]. We also mention the novel applications of semiclassical theory to mesoscopic physics [15]. For example, orbital magnetism has been studied semiclassically for diffusive systems in [16] and for ballistic systems in [17,18]. For reviews see [19].

Ultimately, one would like to study an arbitrary number of interacting particles in any kind of potential. In the present work, we begin by exploring the structure of the trace formula for two noninteracting particles including an examination of the decomposition into bosonic and fermionic spaces. This sets the stage for the interacting  $N$ -body problem to be explored in a subsequent publication [20]. The method employed here uses the fact that the two particle density of states is the auto-convolution of the single particle density of states. Subsequently, we decompose the semiclassical two particle density of states into three distinct contributions, each of which corresponds to specific dynamical properties of the system. Of particular interest is the contribution which corresponds to two particle dynamics.

Billiards have served as prominent model systems in quantum chaos. They combine conceptual simplicity (the model of a free particle in a box) while allowing the full range of classical dynamics, from integrable to chaotic. Therefore, as initial applications of the formalism, we study two noninteracting identical particles in a disk and in a cardioid. The former problem is integrable while the second is chaotic so these two examples provide a direct test of the formalism in the two limiting cases of classical motion. In both cases, we find the semiclassical formalism does a good job of reproducing the quantum density of states.

## II. BACKGROUND THEORY

### A. Single Particle Semiclassical Theory

In this section, we review the formalism for the semiclassical decomposition of the single particle density of

states. Let  $\{\epsilon_i\}$  be the single-particle energies so that the single particle density of states is

$$\rho_1(\epsilon) = \sum_i \delta(\epsilon - \epsilon_i), \quad (1)$$

where the subscript 1 indicates that it is a single particle density. A fundamental property of the quantum density of states is that it can be exactly decomposed into an average smooth part and an oscillatory part [2]

$$\rho_1(\epsilon) = \bar{\rho}_1(\epsilon) + \tilde{\rho}_1(\epsilon). \quad (2)$$

There are various approaches for calculating these quantities [5]. For example, in systems with analytic potentials, the smooth part may be obtained from an extended Thomas-Fermi calculation which is an asymptotic expansion in powers of  $\hbar$ . In billiard systems, where the particle is confined to a spatial domain by the presence of infinitely steep potential walls, the smooth part may be obtained from the Weyl expansion. In two dimensional billiards with piecewise smooth boundaries and Dirichlet boundary conditions, the first three terms of the Weyl expansion is [21]

$$\bar{\rho}_1(\epsilon) = \left( \frac{\alpha \mathcal{A}}{4\pi} - \frac{\alpha^{1/2}}{8\pi} \frac{\mathcal{L}}{\sqrt{\epsilon}} \right) \theta(\epsilon) + \mathcal{K} \delta(\epsilon) + \dots \quad (3)$$

where  $\alpha = 2m/\hbar^2$ ,  $\mathcal{A}$  is the area,  $\mathcal{L}$  is the perimeter and

$$\mathcal{K} = \frac{1}{12\pi} \oint d\kappa(l) + \frac{1}{24\pi} \sum_i \frac{\pi^2 - \theta_i^2}{\theta_i} \quad (4)$$

is the average curvature integrated along the boundary with corrections due to corners with angles  $\theta_i$ . The oscillating part is obtained from semiclassical periodic orbit theory, and in particular the various trace formulas for  $\tilde{\rho}_1(\epsilon)$  of the form [5]

$$\tilde{\rho}_1(\epsilon) \approx \frac{1}{\pi \hbar} \sum_{\Gamma} A_{\Gamma}(\epsilon) \cos \left( \frac{1}{\hbar} S_{\Gamma}(\epsilon) - \sigma_{\Gamma} \frac{\pi}{2} \right). \quad (5)$$

$\Gamma$  denotes topologically distinct periodic orbits,  $S_{\Gamma}(\epsilon)$  is the classical action integral along the orbit  $\Gamma$ . The amplitude  $A_{\Gamma}(\epsilon)$  depends on energy, the period of the corresponding primitive orbit, the stability of the orbit, and whether it is isolated or non-isolated. The index  $\sigma_{\Gamma}$  depends on the topological properties of each orbit. For isolated orbits, it is just the Maslov index. For nonisolated orbits, there may be additional phase factors in the form of odd multiples of  $\pi/4$  which we account for, in a slight abuse of notation, by allowing  $\sigma_{\Gamma}$  to be half-integer. In the case of non-isolated orbits,  $\Gamma$  denotes distinct families of degenerate orbits. The amplitude of an isolated orbit is given by the Gutzwiller trace formula [1]

$$A_{\Gamma}(\epsilon) = \frac{T_{\gamma}(\epsilon)}{\sqrt{|\det(\tilde{M}_{\Gamma} - I)|}} \quad (6)$$

where  $T_{\gamma}(\epsilon)$  is the period of the primitive orbit  $\gamma$ , corresponding to  $\Gamma$  (*i.e.*  $\Gamma$  is an integer repetition of  $\gamma$ ) and  $\tilde{M}_{\Gamma}$  is the stability matrix of that orbit.

## B. Quantum Two Particle Density of States

Now suppose we have a system of two identical noninteracting particles. The total Hamiltonian is the sum of the single particle Hamiltonians and it follows that the energies of the composite system are just the sums of the single particle energies. The analogue of (1) is then

$$\rho_2(E) = \sum_{i,j} \delta(E - (\epsilon_i + \epsilon_j)). \quad (7)$$

A useful relation is that the two particle density of states is the autoconvolution of the single particle density of states:

$$\rho_2(E) = \int_0^E d\epsilon \rho_1(\epsilon) \rho_1(E - \epsilon) = \rho_1 * \rho_1(E), \quad (8)$$

as can be verified by direct substitution. In fact, this works even if the particles are not identical, where the full density is still the convolution of the two distinct single-particle densities. This would also apply to a single particle in a separable potential, which is mathematically equivalent. Rather than encumber the notation to explicitly allow for this possibility, we defer this discussion to Appendix A, where some formulas for nonidentical, noninteracting particles are presented.

We can decompose the two particle density of states for a system of two identical particles into a symmetric and an antisymmetric density,

$$\rho_2(E) = \rho_S(E) + \rho_A(E). \quad (9)$$

We shall use the terms symmetric/antisymmetric and bosonic/fermionic interchangeably. Each partial density may be obtained using a projection operator onto the relevant subspaces resulting in

$$\rho_{S/A}(E) = \frac{1}{2} \left( \rho_2(E) \pm \frac{1}{2} \rho_1 \left( \frac{E}{2} \right) \right). \quad (10)$$

We seek semiclassical approximations to these quantum expressions, a topic which is pursued in the following sections.

## III. SEMICLASSICAL CALCULATIONS FOR THE TWO PARTICLE SYSTEM

Decomposing the single particle density into its smooth and oscillatory components as in (2) gives a decomposition of the two particle density of states into three distinct contributions,

$$\rho_2^{\text{sc}}(E) = \bar{\rho}_1 * \bar{\rho}_1(E) + 2\bar{\rho}_1 * \tilde{\rho}_1(E) + \tilde{\rho}_1 * \tilde{\rho}_1(E). \quad (11)$$

The first term is a smooth function of energy since the convolution of two smooth functions results in a smooth function. This is followed by a cross term and finally by a purely oscillating term. The cross term is also an oscillating function. At first, this may seem incorrect since the convolution of a smooth function with an oscillating function usually yields a smooth function. As we will show, the oscillatory nature of the cross term is due to contributions from the end-points of the convolution integral. Physically, the smooth term does not depend on dynamics since it corresponds to the Weyl formula in the full two-particle space. The cross term depends only on single particle dynamics because it corresponds to the situation where one particle is stationary and the other particle is evolving dynamically on a periodic orbit. It is only the last term which contains two particle dynamics in the sense that both particles are evolving dynamically on periodic orbits. Hence, we will refer to the last term as the dynamical term.

We find a general expression for  $\tilde{\rho}_1 * \tilde{\rho}_1(E)$  by substituting a generalized trace formula for  $\tilde{\rho}_1(E)$  and then evaluating the resulting convolution integral using the method of stationary phase. Using (5), the dynamical term can be written as

$$\tilde{\rho}_1 * \tilde{\rho}_1(E) \approx \frac{1}{(\pi\hbar)^2} \sum_{\Gamma_1, \Gamma_2} \int_0^E d\epsilon A_{\Gamma_1}(\epsilon) A_{\Gamma_2}(E - \epsilon) \cos\left(\frac{1}{\hbar} S_{\Gamma_1}(\epsilon) - \sigma_{\Gamma_1} \frac{\pi}{2}\right) \cos\left(\frac{1}{\hbar} S_{\Gamma_2}(E - \epsilon) - \sigma_{\Gamma_2} \frac{\pi}{2}\right). \quad (12)$$

To evaluate this asymptotically, we should include all critical points in the integration domain. Specifically, this integral has a stationary phase point within the integration domain and finite valued endpoints. We shall show that the stationary phase point corresponds to the situation where both particles are evolving dynamically with the energy partitioned between the two particles in a prescribed way. The endpoint contributions must be evaluated at energies such that one of the particles has all of the energy while the other has no energy. However, this contradicts our assumption that both particles are evolving — this is the definition of the dynamical term. Moreover, if we were to evaluate this contribution, the result would be meaningless since it involves using the trace formula at zero energy where it is known to fail. So we shall omit the contributions from the endpoints; this is discussed more fully in section IVD and in Appendix B, as well as in reference [22].

Hence, we evaluate the integral in (12) using only the stationary phase point. To leading order, we can extend the integration limits over an infinite domain. Writing the cosine functions as complex exponentials yields four integrals; the first is

$$\begin{aligned} & \int_{-\infty}^{\infty} d\epsilon A_{\Gamma_1}(\epsilon) A_{\Gamma_2}(E - \epsilon) \exp\left(\frac{i}{\hbar} (S_{\Gamma_1}(\epsilon) + S_{\Gamma_2}(E - \epsilon))\right) \\ & \approx A_{\Gamma_1}(E_0) A_{\Gamma_2}(E - E_0) \sqrt{\frac{2\pi\hbar}{|\Upsilon(\Gamma_1, \Gamma_2, E)|}} \\ & \exp\left(\frac{i}{\hbar} (S_{\Gamma_1}(E_0) + S_{\Gamma_2}(E - E_0)) + i\nu \frac{\pi}{4}\right) \end{aligned} \quad (13)$$

where

$$\begin{aligned} \Upsilon(\Gamma_1, \Gamma_2, E) &= \left( \frac{\partial^2 S_{\Gamma_1}(\epsilon)}{\partial \epsilon^2} + \frac{\partial^2 S_{\Gamma_2}(E - \epsilon)}{\partial \epsilon^2} \right) \Big|_{E_0} \\ \nu &= \text{sign}(\Upsilon(\Gamma_1, \Gamma_2, E)). \end{aligned} \quad (14)$$

$E_0$  is determined from the stationary phase condition

$$\begin{aligned} & \left( \frac{\partial S_{\Gamma_1}(\epsilon)}{\partial \epsilon} + \frac{\partial S_{\Gamma_2}(E - \epsilon)}{\partial \epsilon} \right) \Big|_{E_0} = 0 \\ & \implies T_{\Gamma_1}(E_0) = T_{\Gamma_2}(E - E_0) \end{aligned} \quad (15)$$

where we have used the fact that the derivative of the action with respect to energy is the period.  $E_0$  is the energy of particle 1,  $E - E_0$  is the energy of particle 2 and  $E$  is the total energy of the composite system. The saddle energy  $E_0$  has a precise physical interpretation; Eq.(15) says that the energies of the two particles are partitioned so that the periods of both periodic orbits are the same. In other words, at  $E_0$ , we have orbits which are periodic in the full two particle phase space since after the period  $T$  both particles return to their initial conditions.

The next integral has the same stationary phase condition as the first integral and is its complex conjugate. The third integral is

$$\int_{-\infty}^{\infty} d\epsilon A_{\Gamma_1}(\epsilon) A_{\Gamma_2}(E - \epsilon) \exp\left(-\frac{i}{\hbar} (S_{\Gamma_1}(\epsilon) - S_{\Gamma_2}(E - \epsilon))\right) \quad (16)$$

and has no stationary phase point since setting the first derivative of the action to zero yields the stationary phase condition

$$T_{\Gamma_1}(E_0) = -T_{\Gamma_2}(E - E_0). \quad (17)$$

The trace formula only involves orbits with positive period, so we ignore this possibility. The last integral is the complex conjugate of the third and will also be ignored.

Adding the contributions from the first two integrals, we arrive at the two particle trace formula:

$$\begin{aligned} \tilde{\rho}_1 * \tilde{\rho}_1(E) &\approx \frac{2}{(2\pi\hbar)^{3/2}} \sum_{\Gamma_1, \Gamma_2} \frac{A_{\Gamma_1}(E_0) A_{\Gamma_2}(E - E_0)}{\sqrt{|\Upsilon(\Gamma_1, \Gamma_2, E)|}} \\ &\cos\left(\frac{1}{\hbar} (S_{\Gamma_1}(E_0) + S_{\Gamma_2}(E - E_0)) - (\sigma_{\Gamma_1} + \sigma_{\Gamma_2}) \frac{\pi}{2} + \nu \frac{\pi}{4}\right). \end{aligned} \quad (18)$$

This result possesses the intuitive properties that, other than factors arising from the stationary phase analysis,

the actions and Maslov indices are additive and the amplitudes are multiplicative. We note that this saddle-point analysis fails for the simplest problem in physics, the harmonic oscillator, where  $\Upsilon = 0$ . This is because the two-particle harmonic oscillator has a higher degree of symmetry than we are accounting for here. This is a nongeneric property specific to the harmonic oscillator. We also stress that we have made no assumption about the stability or structure of the orbits. They can be isolated, stable or unstable or come in families. There are also problems with coexisting isolated orbits and families, such as those of the equilateral triangle billiard [23,5].

Note that the overall  $\hbar$  dependence is not multiplicative but picks up an additional factor of  $\hbar^{1/2}$  from the stationary phase integral. For isolated orbits, the amplitudes  $A$  are independent of  $\hbar$  and the expression (18) has a  $1/\hbar^{3/2}$  prefactor as opposed to the  $1/\hbar$  in the amplitude of the single particle trace formula. The fact that the  $\hbar$  dependence is different implies that the periodic orbits of the full system come in continuous degenerate families rather than isolated trajectories, which in turn implies that there exists a continuous symmetry in the problem [24]. This is an important point which we will address in a companion paper [20]. (It was also noted in [17].) Nonetheless, it may be helpful to give a brief explanation here. Imagine the full phase space periodic orbit  $\Gamma$  consists of particle 1 on a periodic orbit  $\Gamma_1$  with energy  $E_0$  and particle 2 on a distinct periodic orbit  $\Gamma_2$  with energy  $E - E_0$ . We can define  $t = 0$  to be when particle 2 is at some prescribed point on  $\Gamma_2$ . Keeping particle 2 fixed, we can change the position of particle 1 on  $\Gamma_1$  to generate the initial condition of a distinct but congruent periodic orbit in the full phase space. Continuous time translation of the initial condition on  $\Gamma_1$  generates a continuous family of congruent periodic orbits in the full phase space. Since the time translational symmetry can be characterized by a single independent symmetry parameter, the  $\hbar$  dependence is  $\mathcal{O}(1/\sqrt{\hbar})$  stronger than for a system with isolated periodic orbits [24,5].

#### IV. TWO PARTICLE QUANTUM BILLIARDS

As an application of the formalism developed in section III, we consider the quantum billiard problem. Billiards are two dimensional enclosures that constrain the motion of a free particle. Classically, a particle has elastic collisions with the walls and depending on the geometric properties of the domain, the dynamics are either regular or chaotic. For the noninteracting problem, the two particles move independently of each other. In a billiard system, classical orbits possess simple scaling properties. For instance, the action of an orbit  $\Gamma$ ,  $S_\Gamma(\epsilon) = \sqrt{2m\epsilon}L_\Gamma$  and the period of the orbit is

$$T_\Gamma(\epsilon) = \frac{\partial S_\Gamma(\epsilon)}{\partial \epsilon} = \frac{\sqrt{2m}L_\Gamma}{2\sqrt{\epsilon}} = \frac{\hbar\sqrt{\alpha}}{2\sqrt{\epsilon}}L_\Gamma. \quad (19)$$

The parameter  $\alpha = 2m/\hbar^2$  already appeared in Eq. (3); it will recur often. For example, in all final expressions, the energy occurs with  $\alpha$ ; this is a result of the scaling property (the quantity  $\alpha E$  having the units of  $1/\text{length}^2$ .) In the theoretical development, it will be convenient to retain  $\alpha$  and use it to keep track of relative orders in the semiclassical expansions (since it contains  $\hbar$ ). However, once we have the final expressions, we are free to set it to unity for the purposes of numerical comparisons.

We also mention that for billiards, it is common to express the density of states in terms of the wave number  $k$ , where  $\epsilon = k^2/\alpha$  so that  $\rho(k) = 2k\rho(\epsilon)/\alpha$ . This is convenient since  $k$  is conjugate to the periodic orbit lengths  $L$ . Therefore, many of our results will be quoted as a function of  $k$ , although it should be stressed that all convolution integrals must be done in the energy domain. Thus, we shall write

$$\rho_2^{\text{sc}}(k) = (\bar{\rho}_1 * \bar{\rho}_1)(k) + 2(\bar{\rho}_1 * \tilde{\rho}_1)(k) + (\tilde{\rho}_1 * \tilde{\rho}_1)(k). \quad (20)$$

Here, it is understood that each of the functions in brackets is first evaluated in the energy domain and then converted to the  $k$  domain through the Jacobian relation above. This will always be the case when the argument is  $k$ , so that we will not always write brackets around the various functions. In terms of the wavenumber  $k$ , the decomposition (10) becomes

$$\rho_{S/A}(k) = \frac{1}{2} \left( \rho_2(k) \pm \frac{1}{\sqrt{2}} \rho_1 \left( \frac{k}{\sqrt{2}} \right) \right). \quad (21)$$

#### A. Smooth Term

The smooth part is defined by the convolution integral

$$\bar{\rho}_1 * \bar{\rho}_1(E) = \int_0^E d\epsilon \bar{\rho}_1(\epsilon) \bar{\rho}_1(E - \epsilon), \quad (22)$$

where  $\bar{\rho}_1$  is given by the Weyl expansion. The expansion in (3) is taken only to order  $\hbar^0$ . Hence, after expanding the integrand in (22), it is formally meaningless to include terms that are  $\mathcal{O}(1/\hbar)$  since there are corrections of the same order in  $\hbar$  that have not been calculated. Ignoring these terms and performing the necessary integrations, the smooth term is found to be

$$\bar{\rho}_1 * \bar{\rho}_1(E) \approx \frac{\alpha^2 \mathcal{A}^2}{16\pi^2} E - \frac{\alpha^{3/2} \mathcal{A} \mathcal{L}}{8\pi^2} \sqrt{E} + \frac{\alpha \mathcal{L}^2}{64\pi} + \frac{\alpha \mathcal{A} \mathcal{K}}{2\pi}. \quad (23)$$

## B. Cross Term

We next convolve  $\bar{\rho}_1$  term by term with  $\tilde{\rho}_1$ . Asymptotically, each convolution integral receives contributions from the upper and lower endpoints. However, we shall only include one of these, namely the endpoint for which the trace formula is not evaluated at zero energy. As in section III, we neglect the other endpoint for reasons explained in IV D and Ref. [22]. This is also discussed in Appendix B, where we evaluate the various integrals for the cross term exactly using isolated billiard orbits and show explicitly that an appropriate asymptotic expansion of the exact expression leads to consistent results.

After convolution, we find the area term involves the integral

$$\text{Re} \left\{ \int_0^E d\epsilon A_\Gamma(E - \epsilon) \exp \left( i\sqrt{\alpha(E - \epsilon)} L_\Gamma - i\sigma_\Gamma \frac{\pi}{2} \right) \right\}. \quad (24)$$

The lower endpoint  $\epsilon = 0$  corresponds to the physically meaningful situation while the upper endpoint is spurious in the sense mentioned above and discussed in detail below. Hence, to leading order, we can remove the amplitude factor from inside the integral, Taylor expand the argument of the exponential and extend the upper limit to infinity. This leads to

$$I_A(E) \approx \frac{\alpha A}{4\pi^2} \sum_\Gamma \frac{A_\Gamma}{T_\Gamma} \cos \left( \sqrt{\alpha E} L_\Gamma - \sigma_\Gamma \frac{\pi}{2} - \frac{\pi}{2} \right). \quad (25)$$

By similar logic, the perimeter term and curvature terms are

$$I_L(E) \approx -\frac{\sqrt{\alpha} \mathcal{L}}{8\pi^{3/2} \sqrt{\hbar}} \sum_\Gamma \frac{A_\Gamma}{\sqrt{T_\Gamma}} \cos \left( \sqrt{\alpha E} L_\Gamma - \sigma_\Gamma \frac{\pi}{2} - \frac{\pi}{4} \right) \\ I_K(E) \approx \frac{\mathcal{K}}{\pi \hbar} \sum_\Gamma A_\Gamma \cos \left( \sqrt{\alpha E} L_\Gamma - \sigma_\Gamma \frac{\pi}{2} \right). \quad (26)$$

Note that all amplitudes and periods in (25) and (26) are evaluated at the system energy  $E$ . Recall  $\alpha \propto 1/\hbar^2$  so that after convolution the sequence is an expansion in powers of  $\sqrt{\hbar}$  and not in powers of  $\hbar$  as for the original Weyl series (3). We also note that the first correction to  $I_A$  may be of the same order as  $I_K$  (as happens for the disk [25]) and should be included if this is the case. We then have  $\bar{\rho} * \tilde{\rho} \approx I_A + I_L + I_K$ .

## C. Dynamical Term

In this section, we derive a general expression for the dynamical term that is valid for any billiard problem. To this end, the first task is to determine the saddle energy from the stationary phase condition. Inserting (19) into (15) yields

$$\frac{L_{\Gamma_1}}{\sqrt{E_0}} = \frac{L_{\Gamma_2}}{\sqrt{E - E_0}} \quad (27)$$

which implies

$$\frac{E_0}{E} = \frac{L_{\Gamma_1}^2}{L_{\Gamma_1}^2 + L_{\Gamma_2}^2}, \quad \frac{E - E_0}{E} = \frac{L_{\Gamma_2}^2}{L_{\Gamma_1}^2 + L_{\Gamma_2}^2} \quad (28)$$

and

$$\Upsilon(\Gamma_1, \Gamma_2, E) = -\frac{\sqrt{2m}}{4E^{3/2}} \frac{(L_{\Gamma_1}^2 + L_{\Gamma_2}^2)^{5/2}}{L_{\Gamma_1}^2 L_{\Gamma_2}^2}. \quad (29)$$

Clearly  $\nu = -1$ . We then substitute these results into (18) to obtain the two particle trace formula for billiards

$$\tilde{\rho}_1 * \tilde{\rho}_1(E) \approx \frac{4E^{3/4}}{\sqrt{\hbar} \alpha^{1/4} (2\pi \hbar)^{3/2}} \sum_{\Gamma_1, \Gamma_2} \frac{L_{\Gamma_1} L_{\Gamma_2}}{(L_{\Gamma_1}^2 + L_{\Gamma_2}^2)^{5/4}} A_{\Gamma_1}(E_0) A_{\Gamma_2}(E - E_0) \\ \cos \left( \sqrt{\alpha E} \sqrt{L_{\Gamma_1}^2 + L_{\Gamma_2}^2} - (\sigma_{\Gamma_1} + \sigma_{\Gamma_2}) \frac{\pi}{2} - \frac{\pi}{4} \right). \quad (30)$$

If the single particle periodic orbits are not isolated, then one must make direct use of the corresponding single particle amplitudes in (5) evaluated at the appropriate energies. We will show an explicit example of this when we analyze the disk billiard. Note the amplitudes  $A_\Gamma$  typically have an energy dependence so one cannot make any general statements about the energy dependence of this term except that the greater the dimensionality of the periodic orbit families, the greater the energy prefactor. For example, for the disk, it turns out to be  $E^{1/4}$ .

If the single particle periodic orbits are isolated, the amplitudes are given by (6), which for billiards is

$$A_\Gamma(\epsilon) = \frac{\sqrt{\alpha} \hbar}{2\sqrt{\epsilon}} \frac{L_\Gamma}{\sqrt{|\det(\tilde{M}_\Gamma - I)|}}. \quad (31)$$

In this case, the Gutzwiller amplitudes are evaluated at  $E_0$  and  $E - E_0$ , so we again make use of (28). After some algebra and simplification, we find

$$\tilde{\rho}_1 * \tilde{\rho}_1(E) \approx \frac{\alpha^{3/4}}{(2\pi)^{3/2} E^{1/4}} \sum_{\Gamma_1, \Gamma_2} \frac{L_{\Gamma_1} L_{\Gamma_2} (L_{\Gamma_1}^2 + L_{\Gamma_2}^2)^{-1/4}}{\sqrt{|\det(\tilde{M}_{\Gamma_1} - I)|} \sqrt{|\det(\tilde{M}_{\Gamma_2} - I)|}} \\ \cos \left( \sqrt{\alpha E} \sqrt{L_{\Gamma_1}^2 + L_{\Gamma_2}^2} - (\sigma_{\Gamma_1} + \sigma_{\Gamma_2}) \frac{\pi}{2} - \frac{\pi}{4} \right). \quad (32)$$

Note the  $E^{-1/4}$  prefactor which implies that the amplitude decays weakly with energy. This is the same prefactor that occurs in the single particle disk problem. This is not a coincidence, but arises from the fact that in both

problems the periodic orbits come in one parameter families. Also, one must be careful to distinguish between  $L_\Gamma$ , the length of a periodic orbit and  $L_\gamma$ , the length of the corresponding primitive periodic orbit. In general  $L_\Gamma = n_\Gamma L_\gamma$  where  $n_\Gamma$  is the repetition index of that orbit.

#### D. Spurious Endpoint Contributions

As mentioned above, when confronted with convolution integrals, it is natural to analyse them asymptotically. This involves identifying the critical points and doing appropriate expansions in their neighbourhoods. In our work, these critical points are either stationary phase points or endpoints. The power of semiclassical methods is that each critical point can be given an immediate physical interpretation. For example, the stationary phase point in the dynamical term is found to be that energy such that the two particles have the same period so that the motion is periodic in the full two-particle phase space. This is intuitively reasonable. However, the same integral also has endpoints with finite valued contributions. We could do an asymptotic calculation in the vicinity of these points, but we can argue immediately that the result is spurious and not physically meaningful.

Recall the trace formulas are asymptotic in  $\hbar$  which typically also means asymptotic in energy. At the endpoints, one of the trace formulas is evaluated at small energy where it is known to be invalid. Alternatively, we can substitute for the trace formula any expression which is asymptotically equivalent to it and expect all meaningful results to be invariant to leading order. If we do this, we will find the endpoint contribution changes while the stationary phase contribution remains invariant, to leading order.

A further argument is that the structure of the endpoint contribution will be incorrect. Typically, it will be a sinusoid with an argument which does not depend on energy, but only depends on the properties of one of the orbits. Hence, it will have the same asymptotic structure as the cross term. However, we know that the cross term completely describes all such contributions and any further contribution with the same structure must be spurious.

Similarly, when we evaluate the cross term, we have two endpoint contributions. At one of these, we are evaluating the trace formula at some finite energy, which is reasonable. This endpoint corresponds to orbits which are periodic in the full phase space and in which one particle evolves on a single particle periodic orbit with all the energy, while the other remains fixed at some point in phase space with zero energy. At the other endpoint, we are evaluating the trace formula at zero energy, which is problematic. This corresponds to the contradictory situation in which the evolving particle has zero energy while

the fixed particle has all the energy. In addition, upon inspection of this endpoint contribution, we find a function which is not oscillatory in energy and therefore has the same asymptotic structure as the smooth term. However, the smooth term already completely describes the average behaviour of the two particle density of states and any further contributions with the same structure must be spurious.

These situations are further examples of a general situation described in Ref. [22] where it was shown that when integrating over the trace formula to obtain physical quantities, one should include all critical points except ones at which the trace formula is evaluated at zero energy. Such contributions should simply be ignored as spurious. In [22], the application was to thermodynamic calculations, but the principle is precisely the same. In Appendix B, we show the result of evaluating the cross term exactly for isolated orbits. An asymptotic analysis of this result leads to two terms which we can identify as coming from the two endpoints. One has the form used in this paper while the other is clearly spurious.

## V. TWO PARTICLE DISK BILLIARD

In this section, we apply our results to the problem of two identical noninteracting particles moving in a two dimensional disk billiard of radius  $R$ . Quantum mechanically, this problem is a simple extension of the one body problem. Nevertheless, the spectrum has some interesting features which we discuss below.

### A. Quantum Mechanics

For the disk billiard, a general two particle state can be written as

$$|m_1 n_1, m_2 n_2\rangle = |m_1 n_1\rangle \otimes |m_2 n_2\rangle \quad (33)$$

where the azimuthal quantum numbers  $m_1, m_2 = 0, \pm 1, \pm 2, \dots$  and the radial quantum numbers  $n_1, n_2 = 1, 2, 3, \dots$ . We shall also use a more compact notation  $|N_1, N_2\rangle = |m_1 n_1, m_2 n_2\rangle$  where  $N$  denotes a pair of integers  $(m, n)$ . We can immediately write down the wave numbers of the two particle system as

$$k_{N_1 N_2} = \sqrt{\left(\frac{Z_{N_1}}{R}\right)^2 + \left(\frac{Z_{N_2}}{R}\right)^2}, \quad (34)$$

where  $Z_N$  denotes the  $n$ th zero of the  $m$ th Bessel function  $J_m(z)$ . The set of all two particle states is given by  $\{|N_1, N_2\rangle\}$ .

The spectrum is highly degenerate. A typical state  $|N_1, N_2\rangle$  is 8-fold degenerate since we can reverse the sign of either  $m_1$  or  $m_2$  or interchange the two particles and

the resultant state has the same energy. However, if either  $m_1$  or  $m_2$  is zero or if  $N_1 = N_2$  then the state is 4-fold degenerate. If  $m_1 = m_2 = 0$  and  $N_1 \neq N_2$ , then the state is 2-fold degenerate whereas if  $m_1 = m_2 = 0$  and  $N_1 = N_2$ , then the state is nondegenerate. If the particles are in distinct states, the degenerate multiplets divide evenly between the symmetric and antisymmetric spaces. However, if the particles are in the same state,  $N_1 = N_2$ , it is somewhat less trivial. If  $N_1 = N_2$  and  $m_1 = m_2 \neq 0$ , there is a 4-fold degenerate set of states:  $|m\ n, m\ n\rangle$ ,  $|-m\ n, -m\ n\rangle$ ,  $|m\ n, -m\ n\rangle$  and  $|-m\ n, m\ n\rangle$ . The first two states belong to the symmetric space. From the second two states, we can construct one symmetric and one antisymmetric combination. (This is analogous to coupling two spin 1/2 states to construct a 3-fold symmetric  $S = 1$  state and a nondegenerate antisymmetric  $S = 0$  state.) If  $N_1 = N_2$  and  $m_1 = m_2 = 0$ , this yields the state  $|0\ n, 0\ n\rangle$ , which is singly degenerate and belongs to the symmetric space.

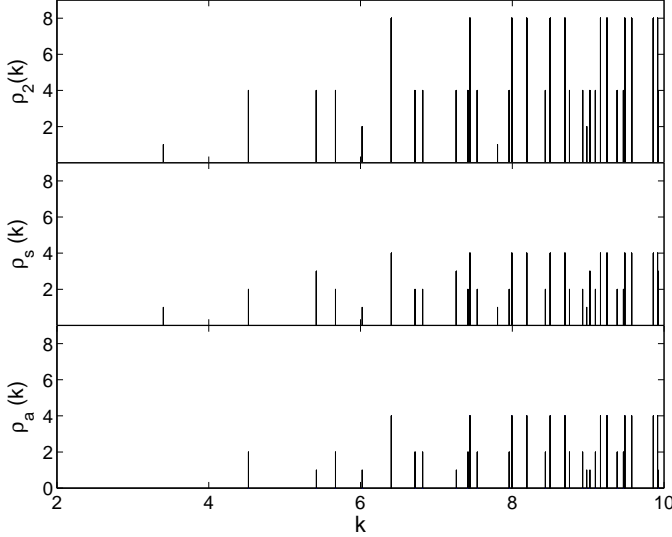


FIG. 1. (Top) The quantum density of states for two identical particles in the disk billiard. (Middle) Bosonic density of states. (Bottom) Fermionic density of states. In each case, the heights indicate the degeneracy of the state.

The quantum density of states

$$\rho_2(k) = \sum_{N_1, N_2} \delta(k - k_{N_1 N_2}) \quad (35)$$

and the corresponding symmetric and antisymmetric densities are shown in Fig. 1 as a function of the wavenumber  $k$ . Note that in this figure some of the peaks have different degeneracies in the symmetric and antisymmetric densities, as discussed above.

## B. Semiclassical Density of States

We first review the semiclassical decomposition of the single particle density of states. The smooth part of the density of states may be obtained using the general result for two dimensional billiards (3). In fact, many higher order terms have been calculated [7]. But, for our purposes, it suffices to use the first three terms as in (3) with  $\mathcal{A} = \pi R^2$ ,  $\mathcal{L} = 2\pi R$  and  $\mathcal{K} = 1/6$ .

The oscillating part of the level density can be obtained using trace formulas for systems with degenerate families of orbits. The periodic orbit families may be uniquely labelled by two integers ( $v, w$ ) where  $v$  is the number of vertices and  $w$  is the winding number around the center. The two integers must satisfy the relation  $v \geq 2w$ . The length of an orbit with vertex number  $v$  and winding number  $w$  is given by  $L_{vw} = 2vR \sin(\pi w/v)$ . With this notation, the trace formula for the oscillating part of the density of states is [26]

$$\tilde{\rho}_1(\epsilon) \approx \frac{\alpha^{3/4}}{2\sqrt{2\pi}\epsilon^{1/4}} \sum_{vw} \frac{\mathcal{D}_{vw} L_{vw}^{3/2}}{v^2} \cos\left(\sqrt{\alpha E} L_{vw} - 3v\frac{\pi}{2} + \frac{\pi}{4}\right) \quad (36)$$

where the sum goes from  $w = 1 \cdots \infty$  and  $v = 2w \cdots \infty$  and the degeneracy factor  $\mathcal{D}_{vw}$ , which accounts for negative windings, is 1 for  $v = 2w$  and 2 for  $v > 2w$ . Comparing (36) with the general form (5), we identify

$$A_{vw}(\epsilon) = \frac{\sqrt{2\pi}\alpha^{3/4}\hbar\mathcal{D}_{vw}L_{vw}^{3/2}}{4v^2\epsilon^{1/4}} \quad (37)$$

$$\sigma_{vw} = 3v - \frac{1}{2}$$

Adding the smooth and oscillating terms gives the semiclassical approximation to the single particle density of states which we denote by  $\rho_1^{\text{sc}}(\epsilon)$ .

To evaluate the semiclassical approximation to the two particle density of states, we must evaluate the smooth, cross and dynamical terms. The smooth term can be taken from Eq. (23) to be

$$\bar{\rho}_1 * \bar{\rho}_1(E) \approx \frac{\alpha^2 R^4}{16} E - \frac{\alpha^{3/2} R^3}{4} \sqrt{E} + \left(\frac{3\pi + 4}{48}\right) \alpha R^2. \quad (38)$$

The arguments of the previous section and in particular Eqs. (25) and (26) lead to the cross term

$$\bar{\rho}_1 * \tilde{\rho}_1(E) \approx \frac{\alpha^{5/4} R^2 E^{1/4}}{4\sqrt{2\pi}} \sum_{vw} \frac{\sqrt{L_{vw}} \mathcal{D}_{vw}}{v^2} \left( \cos\left(\Phi_{vw} - \frac{\pi}{2}\right) - \sqrt{\frac{\pi}{2}} \chi_{vw} \cos\left(\Phi_{vw} - \frac{\pi}{4}\right) + \left(\frac{1}{3} + \frac{R^2}{2L_{vw}^2}\right) \chi_{vw}^2 \cos \Phi_{vw} \right) \quad (39)$$

where  $\Phi_{vw} = \sqrt{\alpha E} L_{vw} - 3v\pi/2 + \pi/4$  and  $\chi_{vw} = \sqrt{L_{vw}}/(\alpha E)^{1/4} R$ . We have also included the first correction to the area term,  $I_A(E)$  which appears in the third term above [25]. The dynamical term can be obtained using (30). Noting that  $\Gamma_i$  in (30) corresponds to the pair of integers  $(v_i, w_i)$ , the result is

$$\tilde{\rho}_1 * \tilde{\rho}_1(E) \approx \frac{\alpha^{5/4} E^{1/4}}{4\sqrt{2\pi}} \sum_{v_1 w_1, v_2 w_2} \left( \prod_{i=1}^2 \frac{\mathcal{D}_i L_i^2}{v_i^2} \right) L_{12}^{-3/2} \cos \left( \sqrt{\alpha E} L_{12} - 3(v_1 + v_2) \frac{\pi}{2} + \frac{\pi}{4} \right) \quad (40)$$

where we defined  $L_i = L_{v_i w_i}$ ,  $\mathcal{D}_i = \mathcal{D}_{v_i w_i}$  and  $L_{12} = \sqrt{L_1^2 + L_2^2}$ .

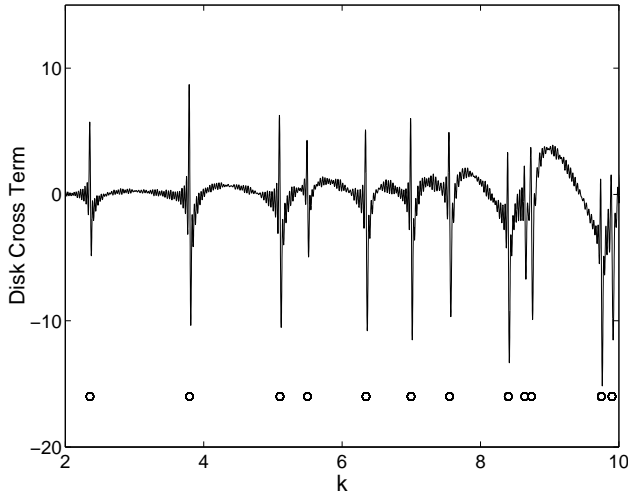


FIG. 2. The cross term (39) of the semiclassical density of states for two identical particles in the disk billiard. In this case, we truncate the sum in (39) at  $w_{\max} = 50$ ,  $v_{\max} = 100$ . The circles indicate the level sequence of the one body problem obtained from EBK quantization. Note the kinks that occur at these positions.

### C. Numerics

For numerical purposes, we take  $\alpha = \hbar = 1$  and  $R = 1$  so the single-particle energies are just the squares of the zeros of Bessel functions. Since we can only include a finite number of orbits, the periodic orbit sums must be truncated. As a representative case, we truncate the sum in (39) at  $w_{\max} = 50$ ,  $v_{\max} = 100$  (see Fig. 2) and use the same limits to truncate the quadruple sum in (40). This is a relatively small set of orbits, yet it does very well in reproducing the peaks of the quantum density of states. As an illustration, we show the first few peaks of (20) in Fig. 3. We calculated the semiclassical density of states (20) on the interval  $0 \leq k \leq 11$ . After doing so,

we found only two sets of two peaks which were not resolved. These are shown in Fig. 4. Obviously, using more orbits will produce better results, but this increases the computation time because of the quadruple sum in (40). (Although one can reduce the computational overhead by limiting the sum to orbits whose amplitude exceeds some prescribed threshold [25]).

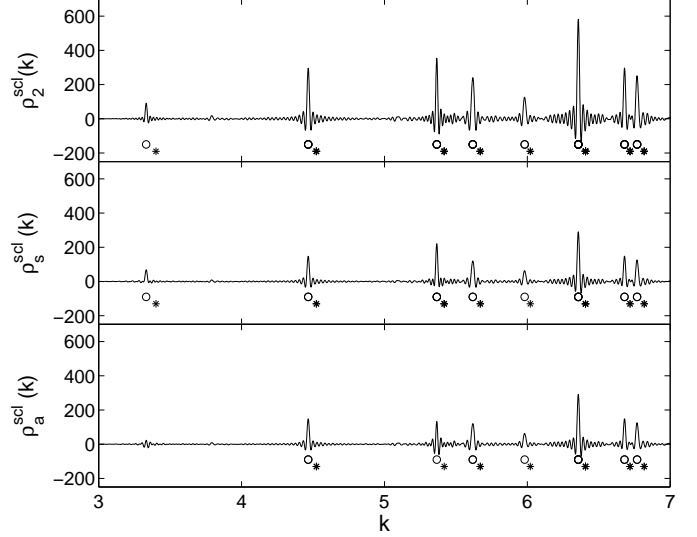


FIG. 3. (Top) The first few peaks of the semiclassical density of states (20). (Middle) Semiclassical approximation to the bosonic density of states. (Bottom) Semiclassical approximation to the fermionic density of states. In each case, the circles and stars represent the appropriate level sequences obtained from EBK quantization and quantum mechanics, respectively. Note the positions of the peaks more closely reproduce the EBK spectrum.

As an additional test, we want to determine whether (20) gives the correct degeneracies. We could do this by integrating the area under each of the peaks. However, a simpler procedure is to do a Gaussian smoothing by convolving  $\rho_2^{\text{sc}}(k)$  with an unnormalized Gaussian of variance  $\sigma$ :

$$\rho_2^{\text{sc}}(k) * G_\sigma(k) = \int_0^\infty dk' \rho_2^{\text{sc}}(k') G_\sigma(k - k') \quad (41)$$

where

$$G_\sigma(k) = \exp(-k^2/2\sigma^2). \quad (42)$$

and  $\sigma$  is the smoothing width. The reason for this is that if the variance  $\sigma$  of the Gaussian is larger than the intrinsic width of a peak in the semiclassical spectrum, then each peak acts like  $d\delta(k - k_n)$  with respect to the Gaussian. Thus, the integral in (41) becomes  $dG_\sigma(k - k_n)$  or  $d$  at  $k = k_n$ . Of course, this is invalid when the spacing between two adjacent peaks is smaller than about  $\sigma$ . Some examples are discussed in the next section.



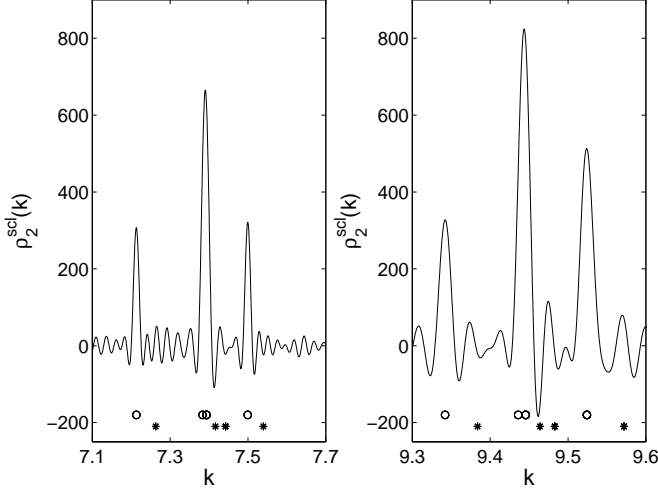


FIG. 4. Two sets of two peaks in the semiclassical density of states (20) that are not resolved. (Left) The middle peak is not resolved into the two peaks at  $k = 7.4163$  (corresponding to the quartet  $\{|0\ 1, \pm 1\ 2\rangle, |\pm 1\ 2, 0\ 1\rangle\}$ ) and  $k = 7.4423$  (corresponding to the octet  $\{|\pm 1\ 1, \pm 3\ 1\rangle, |\pm 3\ 1, \pm 1\ 1\rangle\}$ ). The corresponding EBK quartet and octet energies occur at  $k = 7.3831$  and  $k = 7.3932$  respectively. Note that  $\Delta k_{EBK} = 0.0101$  and  $\Delta k_{QM} = 0.026$  so that the spacing of the two unresolved levels is smaller in the semiclassical spectrum than in the quantum mechanical spectrum. (Right) The middle peak is not resolved into the two peaks at  $k = 9.4641$  (corresponding to the quartet  $\{|\pm 1\ 1, 0\ 3\rangle, |0\ 3, \pm 1\ 1\rangle\}$ ) and  $k = 9.4829$  (corresponding to the octet  $\{|\pm 3\ 1, \pm 1\ 2\rangle, |\pm 1\ 2, \pm 3\ 1\rangle\}$ .) The corresponding EBK quartet and octet energies occur at  $k = 9.4359$  and  $k = 9.4456$  respectively. Here,  $\Delta k_{EBK} = 0.0097$  and  $\Delta k_{QM} = 0.0188$ .

We also studied the symmetrised densities by using the expression (21) for both the quantum and semiclassical densities and convolving as above. The periodic orbit sums in the oscillating parts of the one and two body densities were truncated in the standard manner as before. The result of this numerical procedure is shown in Fig. 5. Clearly, the semiclassical approximations reproduce the correct degeneracies of the quantum spectrum as well as the approximate positions.

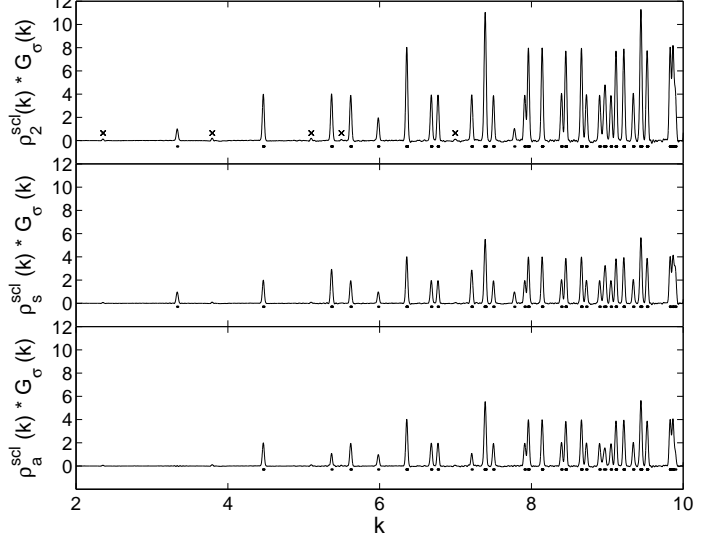


FIG. 5. (Top) The smoothed semiclassical density of states obtained from numerical convolution of (20) with (42). Note the artifacts of the single particle EBK spectrum which occur at the positions marked by an “X”. (Middle) The smoothed semiclassical bosonic density of states obtained from numerical convolution of (21) (with the + sign) and (42). (Bottom) The smoothed semiclassical fermionic density of states obtained from numerical convolution of (21) (with the - sign) and (42). In each case the sequence of dots represent the corresponding EBK spectrum and  $\sigma = 0.0125$ .

#### D. Discussion

In [26], it was noted that the trace formula replicates the single particle EBK spectrum obtained from torus quantisation more precisely than it duplicates the exact single particle quantum spectrum. After inspection of Figs. 3 and 5, we notice the same effect in the two particle spectrum. This property of the trace formulas also accounts for the unresolved peaks in the semiclassical spectrum. When the spacing of two levels of the EBK spectrum is very small, our truncated trace formulas may not resolve them, regardless of the spacing of the corresponding levels in the quantum spectrum (cf. Fig. 4).

Comparing Figs. 1 and 5, we observe generally good agreement between the quantum and semiclassical spectra. Still, there are some apparent inconsistencies, for example, the two tall peaks in Fig. 5. These are the two sets of unresolved levels in Fig. 4, in each case an octet and a quartet. The reason for the discrepancy is the level spacings are smaller than the smoothing width  $\sigma$ , in contradiction to the assumption above, so that the peak height does not equal the degeneracy. In fact, the peak heights observed are rather close to 12 since the octets and quartets are very nearly degenerate on the scale of  $\sigma$  and act almost like a 12-fold degenerate set.

It is not perfectly 12 due to the fact that the degeneracy is not perfect. However, we also observe that the integrated weight under the peak is consistent with a set of 12 energy levels. Other inconsistencies in Fig. 5 occur for the same reason. Of course, overall improvements can be made by including more orbits. As well, the artifacts of the single particle spectrum (some examples are marked by an “X” in Fig. 5) which arise from errors in the cross term, presumably decrease when corrections to the single particle trace formula are incorporated into the cross term. These preliminary numerical findings support our analytical results, which we now test in the rather different context of a chaotic billiard.

## VI. TWO PARTICLE CARDIOID BILLIARD

In this section, we study the problem of two identical noninteracting particles evolving in the cardioid billiard, which is fully chaotic [27]. Since the billiard has a reflection symmetry, all the quantum states are either even or odd (this symmetry should not be confused with the symmetric/antisymmetric symmetry due to particle exchange.) In the subsequent analysis, we will exclusively use the odd spectrum. The reason for this is to avoid the additional complication of *diffractive* orbits which strike the vertex. Classically, these orbits are undefined and are therefore not included in the standard Gutzwiller theory. Studies of diffractive effects in trace formulas can be found in [28,29]. The latter reference explores the specific application to the cardioid and shows that diffractive orbits are important in describing the even spectrum but are largely absent from the odd spectrum.

We could proceed as before by doing an explicit semiclassical analysis of each term in the decomposition of the two particle semiclassical density of states (11). However, we can simplify the analysis by removing single particle dynamics from the discussion. That is, we will focus exclusively on those quantum mechanical and semiclassical quantities that inherently describe two particle dynamics. More specifically, we compare the Fourier transform of the dynamical term

$$\tilde{F}_2^{\text{sc}}(L) = \mathcal{F}\{\tilde{\rho}_1 * \tilde{\rho}_1(k)\} \quad (43)$$

with its quantum mechanical analogue which we define to be

$$\tilde{F}_2^{\text{qm}}(L) = \mathcal{F}\{\rho_2(k) - \bar{\rho}_1 * \bar{\rho}_1(k) - 2\bar{\rho}_1 * \tilde{\rho}_1(k)\}. \quad (44)$$

The integral operator  $\mathcal{F}$  will be defined precisely below. In the semiclassical transform (43), we use (32) expressed in terms of the wavenumber. Here,  $\Gamma_1$  and  $\Gamma_2$  are periodic orbits in the fundamental domain (*i.e.* the half-cardioid.) Orbit properties are discussed in [30,29] and some representative orbits are shown in Fig. 6. The stability matrices in the denominators of the single particle

Gutzwiller amplitudes are computed using the standard prescription for the stability of free flight billiards (see, for example, [4].)

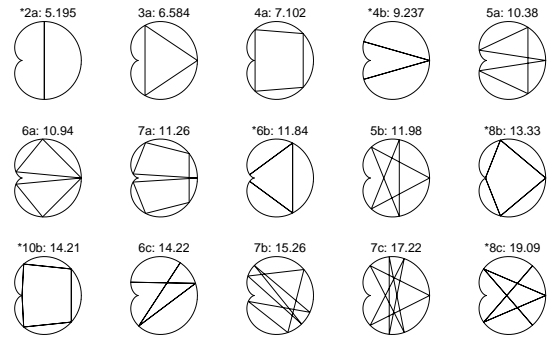


FIG. 6. Some of the shorter periodic orbits of the cardioid in the full domain. The label of each orbit includes the number of reflections and also a letter index to further distinguish it. The two orbits \*8b and \*10b reflect specularly near the cusp, contrary to appearances while the orbit 4a misses the cusp. From [29].

In the quantum mechanical analogue (44),  $\rho_2(k)$  is the quantum two particle density of states  $\rho_2(k) = \sum_I \delta(k - k_I)$  where the superindex  $I$  denotes the pair of integers  $(i, j)$  and  $k_I = \sqrt{k_i^2 + k_j^2}$ . In (44), we subtract the smooth average part and the part which contains single particle dynamics. Using  $\mathcal{A} = 3\pi/4$ ,  $\mathcal{L} = 6$ , and  $\mathcal{K} = 3/16$  in (23), the smooth term is

$$\bar{\rho}_1 * \bar{\rho}_1(E) \approx \frac{9}{256} \alpha^2 E - \frac{9}{16\pi} \alpha^{3/2} \sqrt{E} + \left( \frac{9}{16\pi} + \frac{9}{128} \right) \alpha. \quad (45)$$

The cross term is given by the general expressions (25) and (26).

### A. Numerics for the Unsymmetrised Cardioid

As before, we take  $\alpha = 1$  and use a standard sized cardioid as in [29] to obtain the single particle spectrum. In this section, we numerically compare the two particle quantum mechanics with the two particle semiclassics. We do this by making a direct comparison of the Fourier transforms in the reciprocal space of orbit lengths,  $L$ . In this space, we expect peaks at lengths which correspond to the Euclidean lengths of the *full* periodic orbits of the two particle system. For instance, if the full orbit  $\Gamma$  is comprised of particle 1 travelling on the orbit  $\Gamma_1$  and particle 2 traversing a distinct orbit  $\Gamma_2$ , we expect a peak at  $L_\Gamma = \sqrt{L_{\Gamma_1}^2 + L_{\Gamma_2}^2}$ . In the event that both particles

are on the same orbit  $\Gamma$ , we expect a peak at  $\sqrt{2}L_\Gamma$ . In this way, any peak in the two particle spectrum can be attributed to the dynamics of a particular periodic orbit of the full classical phase space.

We construct the two particle spectrum by adding the energies of the single particle spectrum. We include the first 1250 single particle energies which allows us to construct the first 766794 two particle energy levels representing all two particle energies less than  $6.8856 \times 10^3$ .

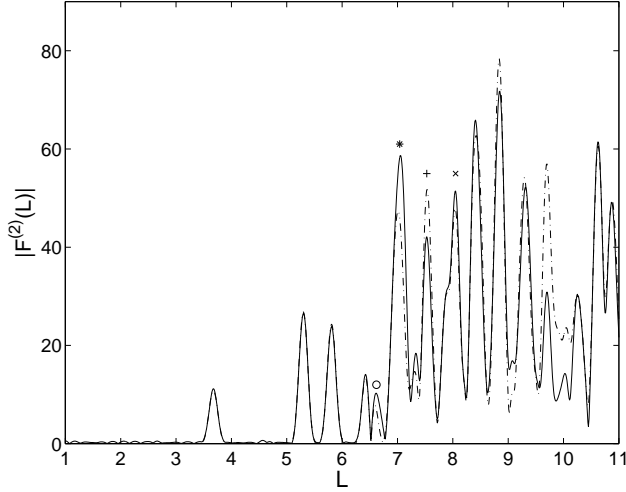


FIG. 7. The Fourier transform of the dynamical (purely oscillating) part of the two particle density of states. The solid line is the transform of the quantum mechanical two particle spectrum (44) and the dashed-dotted line is the transform of the semiclassical two particle trace formula (43). All relevant geometrical periodic orbits with length  $L_\Gamma = \sqrt{L_{\Gamma_1}^2 + L_{\Gamma_2}^2} < 11$  have been included. (Symbols described in the text.)

For a precise numerical comparison, we define the Fourier transform

$$\mathcal{F}\{f(k)\} = \int_{-\infty}^{\infty} dk w(k) e^{ikL} f(k) \quad (46)$$

as a function of the conjugate variable  $L$ . Here,  $w(k)$  is the three term Blackman-Harris window function [31]

$$w(k) = \begin{cases} \sum_{j=0}^2 a_j \cos\left(2\pi j \frac{k-k_0}{k_f-k_0}\right) & k_0 < k < k_f \\ 0 & \text{otherwise} \end{cases} \quad (47)$$

with  $(a_0, a_1, a_2) = (0.42323, -0.49755, 0.07922)$ . We choose  $k_0$  and  $k_f$  so that the window function goes smoothly to zero at the first and last eigenvalues of the two particle spectrum. Numerical integration of (43) and (44) using this definition of  $\mathcal{F}$  is displayed in Fig. 7. In the semiclassical transform, a total of 100 periodic orbits including multiple repetitions were used.

In Fig. 7, we observe good agreement between the quantum and semiclassical results for  $L < 6.5$  and  $L > 10.3$ . In the region  $6.5 < L < 10.3$ , there are appreciable discrepancies for the following reason. Recall that the amplitudes of the two particle trace formula (32) apply only to billiard systems whose single particle periodic orbits are isolated. In the single particle cardioid problem, there exist orbits which are not well isolated in phase space, in fact two geometric orbits and a diffractive orbit are sometimes very close in phase space. For example, the two geometric orbits  $4a$  and  $*10b$  together with the similar looking diffractive orbit  $4a'$  (not shown) [29]. In this event, the stationary phase approximation underlying the Gutzwiller formalism fails as does the argument in that diffractive orbits do not affect the odd spectrum. As a result, whenever a two particle orbit in the full space is comprised of one or both particles on one of these problematic single particle periodic orbits, the resulting two particle amplitude is inaccurate. (There is recent work on uniform approximations to account for such effects [32], unfortunately it seems not to apply to the cardioid which has the additional curious feature that the boundary curvature is infinite at the vertex.)

$\Lambda$	$\Gamma_1$	$\Gamma_2$	$L_\Lambda$
1	$\frac{1}{2}(*2a)$	$4a$	7.562
2	$\frac{1}{2}(*2a)$	$\frac{1}{2}(*10b)$	7.565

TABLE I. A few of the periodic orbits which conspire to give trouble around  $L=7.6$ .

We now consider some specific examples. Consider first the peak structures “o” and “\*”. In this region, the single particle trace formula is erroneous [29] and these errors propagate through to the cross term and inevitably to the quantum mechanical transform. We have also computed the cross term using quantum mechanics, that is, using  $\tilde{\rho}_1 = \rho_1 - \bar{\rho}_1$  in (44) and confirmed that these discrepancies do not arise (cf. Figs. 8 and 9). Thus, these discrepancies are due to errors in the semiclassical approximations of the cross term.

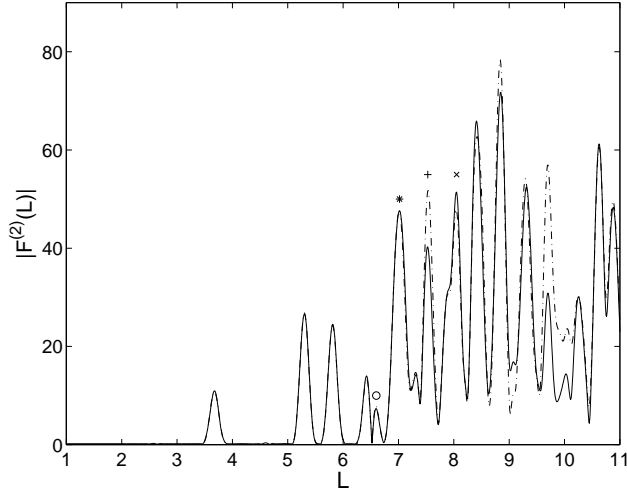


FIG. 8. Same as figure 7 except that the cross term in (44) is computed using single particle quantum mechanics. (Symbols described in the text.)

For the rest of the discussion,  $\Lambda$  refers to a particular periodic orbit family in the full phase space with each two particle orbit in this family comprised of the single particle orbits  $\Gamma_1$  and  $\Gamma_2$  and  $L_\Lambda$  are the lengths of the orbits in each family. Next, consider the peak structure at  $L \approx 7.5$  (+). There are two families of orbits,  $\Lambda_1$  and  $\Lambda_2$  that are responsible for these peaks. The underlying structures of these orbits are shown in Table I. Bearing in mind the two single particle orbits  $\Gamma_2$  are not well isolated (cf. Fig. 6), the Gutzwiller amplitude of each  $\Gamma_2$  is incorrect. Consequently, the two particle Gutzwiller amplitude will also be incorrect, as Fig. 7 demonstrates. Let us look at the next peak structure. Clearly, the quantum peak heights are underestimated at  $L \approx 8.1$  (X). We account for this by recognizing the two particle orbit structure involves the single particle orbit  $\frac{1}{2}(*8b)$  which is an orbit which passes close to the vertex. More specifically, the orbit family  $\Lambda = 3$  is composed of single particle orbits  $\Gamma_1 = \frac{1}{2}(*4b)$  and  $\Gamma_2 = \frac{1}{2}(*8b)$  and the lengths of these two particle orbits are  $L_\Lambda = 8.109$ . As a final illustration, we consider the region  $9.6 < L < 10.3$ . In this neighbourhood, the semiclassics are particularly bad. This can be

accounted for by inspection Table II.

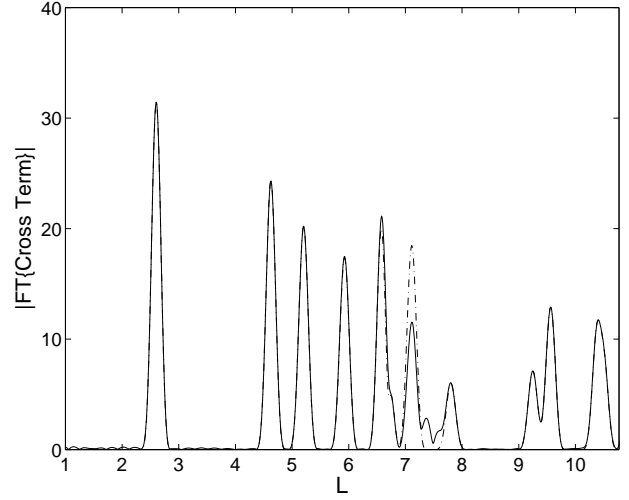


FIG. 9. The cross term for the cardioid billiard calculated using quantum mechanics (solid) and periodic orbit theory (dashed-dotted).

As the Table II and Fig. 6 show, there are many instances where both of the single particle orbits constituting the full orbit are poorly isolated. In view of this, both single particle Gutzwiller amplitudes are incorrect making the product even worse. This accounts for the gross inconsistencies in this region of the reciprocal space. The other discrepancies can be accounted for in a similar manner.

$\Lambda$	$\Gamma_1$	$\Gamma_2$	$L_\Lambda$
4	3a	4a	9.684
5	3a	$\frac{1}{2}(*10b)$	9.687
6	4a	$\frac{1}{2}(*8b)$	9.740
7	$\frac{1}{2}(*8b)$	$\frac{1}{2}(*10b)$	9.742
8	4a	4a	10.044
9	4a	$\frac{1}{2}(*10b)$	10.046
10	$\frac{1}{2}(*10b)$	$\frac{1}{2}(*10b)$	10.048

TABLE II. A few of the periodic orbits which conspire to give trouble around  $L=10$ .

## B. Symmetry Decomposition

In this section, we explore the symmetry decomposition of the two particle problem and in particular the comparison of the symmetrized two particle quantum mechanics with the corresponding two particle semiclassics. We start by defining the smooth and oscillating symmetry reduced densities of states from (21)

$$\bar{\rho}_{S/A}(k) = \frac{1}{2} \left( (\bar{\rho}_1 * \bar{\rho}_1)(k) \pm \frac{1}{\sqrt{2}} \bar{\rho}_1 \left( \frac{k}{\sqrt{2}} \right) \right) \quad (48)$$

and

$$\bar{\rho}_{S/A}^{\text{dyn}}(k) = \frac{1}{2} \left( (\tilde{\rho}_1 * \tilde{\rho}_1)(k) \pm \frac{1}{\sqrt{2}} \tilde{\rho}_1 \left( \frac{k}{\sqrt{2}} \right) \right). \quad (49)$$

While the second term in (49) is a single particle density, in a future paper [20] we will demonstrate that this term describes the physical situation in which two particles are traversing the same periodic orbit, with the same energy and are exactly half a period out of phase. It describes the effect of particle exchange on the spectrum and for this reason affects the symmetric and antisymmetric spaces differently and is based on the theory of Ref. [33]. Therefore, this second term also belongs to the two-particle dynamical term and we identify (49) as being a purely dynamical term. We want to compare it with the corresponding term in the symmetrized quantum densities of states. Hence, in analogy with the previous subsection, we compare

$$\tilde{F}_{S/A}^{\text{dyn}}(L) = \mathcal{F} \left\{ \bar{\rho}_{S/A}^{\text{dyn}}(k) \right\} \quad (50)$$

and

$$\tilde{F}_{S/A}^{\text{qm}}(L) = \mathcal{F} \left\{ \rho_{S/A}(k) - \bar{\rho}_{S/A}(k) - \bar{\rho}_1 * \tilde{\rho}_1(k) \right\}. \quad (51)$$

where  $\rho_{S/A}(k)$  is the quantum bosonic (S) or fermionic (A) density of states.

## C. Numerics for the Symmetrized Cardioid

In this section, we numerically compare the symmetrized quantum mechanics with the corresponding semiclassical quantities. In particular, we compute the transforms (50) and (51) [34]. The symmetrized quantum densities are

$$\begin{aligned} \rho_S(k) &= \sum_{i < j} \delta \left( k - \sqrt{k_i^2 + k_j^2} \right) + \sum_i \delta \left( k - \sqrt{2} k_i \right), \\ \rho_A(k) &= \sum_{i > j} \delta \left( k - \sqrt{k_i^2 + k_j^2} \right) \end{aligned} \quad (52)$$

using the same constraint on the energies as above. Of course, the sum of these symmetrized densities is the total density of states.

Before presenting our numerical results, we describe what we expect. First, all the peaks of the unsymmetrized two particle density should be present. In addition, for each periodic orbit  $\Gamma$ , there should also be peaks at lengths  $L_\Gamma/\sqrt{2}$  arising from the oscillating part of the single particle density of states. The results are shown in Fig. 10. For the two particle density term of (50), we used the same 100 two particle orbits of section VIA while in the single density term we used all single particle orbits with length  $L < 11$ . As well, we included the single particle orbit  $\frac{1}{2}(*10h)$  (not shown in Fig. 6) which has a length  $L = 10.477$ . Fig.10 displays the peak structure in the reciprocal space up to  $L = 6.75$ .

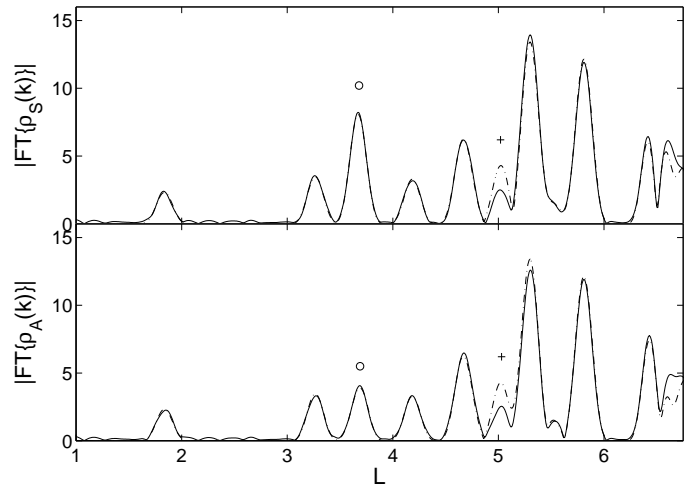


FIG. 10. The Fourier transform of the quantum and semiclassical symmetrized densities of states; the top is bosonic and the bottom is fermionic. In both cases the solid line is the transform of the quantum density of states and the dashed-dotted line is the transform of its semiclassical approximation. Peaks with symbols are described in the text.

We notice that most of the amplitude divides evenly between the symmetric and the antisymmetric densities. Nonetheless, there are exceptions such as the peak at  $L \approx 3.6$  (o). Here, both terms of (49) contribute and the difference in the sign of the second term accounts for the uneven amplitude division between the two symmetrized densities. Semiclassically, we account for the peak structure by noting that two different physical situations are responsible for the peak structure “o”. First, there is the situation in which both particles are on the orbit  $\Gamma = \frac{1}{2}(*2a)$  with no restrictions on the time phase difference between the two particles. This contribution comes from the two particle density term resulting in a peak at a length  $\sqrt{2}L_\Gamma = 3.673$  and produces identical structures in both densities. The second situation occurs when both particles are on the orbit  $\Gamma = (*2a)$  exactly half a period

out of phase. This contribution comes from the single density term at  $L_{\Gamma}/\sqrt{2} = 3.673$  and is explained more fully in [20]. Since the second contribution comes with a different sign in the two symmetries, the amplitudes are different for the symmetric and antisymmetric spaces. In this particular case, it is stronger in the symmetric density and weaker in the antisymmetric density, although the opposite may be true in other cases.

In closing, we remark that the overall agreement between the quantum and semiclassical calculations is good. The poorly reproduced peak just above  $L = 5$  (+) comes from the single density term. This is just the poorly reproduced peak of the single particle density at  $L \approx 7$  shifted down by a factor of  $\sqrt{2}$ .

## VII. CONCLUSION

Initially, we developed a semiclassical formalism to describe the two particle density of states. After deriving a trace formula describing two particle dynamics, we investigated its structure and noted intuitive properties such as the additivity of the actions and topological phase factors. As well, we briefly explained the structure of the full two particle orbits which come in degenerate families. As a first application, we wrote down a two particle trace formula for two identical particles in a billiard. The semiclassical symmetry decomposition involved formal substitution of the semiclassical quantities into the quantum mechanical expressions for the symmetrized densities. In a future paper [20], we show how these formal expressions emerge directly from the classical structures.

Following these general considerations, we studied two identical noninteracting particles in a disk and in a cardioid. In each case, we find that the formalism correctly reproduced the full and symmetrized densities of states. In the integrable problem, we found that our formalism replicates the two body EBK spectrum more precisely than the quantum spectrum, suggesting a deep connection between periodic orbit theory and EBK quantization for integrable systems. In the chaotic cardioid billiard, we note that the single particle orbits which pass close to the vertex lead to inconsistencies in the Fourier transform of the semiclassical density of states. Clearly, our formalism fails here because the Gutzwiller theory itself fails for these “semi-diffractive” orbits. For all other orbits, the two particle trace formula works very well.

The techniques employed here involve the classical phase space of each particle. In a future paper [20], we derive the same results by working in the full two particle phase space. This approach has the advantage of being more general than what we have presented here. Nonetheless, it is conceptually useful to see how the same structure emerges from these two distinct points of view. We would also like to incorporate interactions between the particles. Such a project would undoubtedly require

working in the full phase space since it is no longer true that the full density of states is the convolution of the single particle level densities. This provides an additional motivation for working out the noninteracting problem in the full phase space as a first step towards the more ambitious goal. This full phase space analysis also generalises more readily to more particles. Finally, it has the conceptual advantage that the spurious endpoint contributions discussed in IV D and Appendix B do not arise and therefore need not be explained away.

It may be argued that interacting many body systems are too complex to be accessible to the semiclassical method. However, given the intractability of the many body problem, there may well be questions which semiclassical theory can answer. In particular, we have in mind the applications of semiclassical theory to mesoscopic physics [19]. Here, our seemingly academic study of billiard systems finds physical applications in the context of nanostructures. For example, the disk billiard can serve as a realistic lowest-order approximation to the mean field of the electrons in a circular quantum dot [35]. In fact, many phenomena in ballistic mesoscopic systems can, at least qualitatively, be described by using quantum billiards with independent particles as physical models.

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## VIII. APPENDIX A: NONIDENTICAL PARTICLES

As we have mentioned, most of the discussion still applies if the two particles are not identical. Another situation is a single particle in a separable potential. For example, in two dimensions, one could have  $V(x, y) = V_a(x) + V_b(y)$  in which case, the dynamics in the  $x$  direction are completely uncoupled from the dynamics in the  $y$  direction so that the system is formally the same as if there were distinct particles executing the  $x$  and  $y$  motions. The formalism presented above follows in a natural way. The main differences are that one no longer considers the symmetrised density of states since the symmetry of particle exchange no longer exists and secondly there are two distinct cross terms so that (11) is replaced by

$$\begin{aligned} \rho_2(E) = & \bar{\rho}_{1a} * \bar{\rho}_{1b}(E) + \bar{\rho}_{1a} * \tilde{\rho}_{1b}(E) \\ & + \tilde{\rho}_{1a} * \bar{\rho}_{1b}(E) + \tilde{\rho}_{1a} * \tilde{\rho}_{1b}(E), \end{aligned} \quad (53)$$

where the indices  $a$  and  $b$  refer to the two distinct particles, while the indices 1 and 2 still refer to one or two particle densities of states.

Imagine, for example, that we have two nonidentical particles in distinct billiard enclosures. We introduce two parameters,  $\alpha_a = 2m_a/\hbar^2$  and  $\alpha_b = 2m_b/\hbar^2$ . The smooth term (23) is replaced by

$$\begin{aligned}\bar{\rho}_{1a} * \bar{\rho}_{1b}(E) &\approx \frac{\alpha_a \alpha_b \mathcal{A}^2}{16\pi^2} E \\ &- \left( \alpha_a^{1/2} + \alpha_b^{1/2} \right) \frac{\sqrt{\alpha_a \alpha_b}}{16\pi^2} \mathcal{A} \sqrt{E} \\ &+ \frac{\alpha_a^{1/2} \alpha_b^{1/2} \mathcal{L}^2}{64\pi} + \frac{(\alpha_a + \alpha_b) \mathcal{A} \mathcal{K}}{4\pi}.\end{aligned}\quad (54)$$

The cross terms each separately have the same structure as the cross term for identical particles. Obviously, they are no longer equal to each other, but functionally little has changed. It is just a question of inserting the relevant information from the different smooth and oscillating densities of states of the two particles. Following the same logic as before, we find

$$\begin{aligned}I_A(E) &\approx \frac{\alpha_a \mathcal{A}_a}{4\pi^2} \sum_{\Gamma_b} \frac{A_{\Gamma_b}}{T_{\Gamma_b}} \cos\left(\Phi_{\Gamma_b} - \frac{\pi}{2}\right), \\ I_{\mathcal{L}}(E) &\approx -\frac{\sqrt{\alpha_a} \mathcal{L}_a}{8\pi^{3/2} \sqrt{\hbar}} \sum_{\Gamma_b} \frac{A_{\Gamma_b}}{\sqrt{T_{\Gamma_b}}} \cos\left(\Phi_{\Gamma_b} - \frac{\pi}{4}\right), \\ I_{\mathcal{K}}(E) &\approx \frac{\mathcal{K}_a}{\pi \hbar} \sum_{\Gamma_b} A_{\Gamma_b} \cos(\Phi_{\Gamma_b}),\end{aligned}\quad (55)$$

where  $\Phi_{\Gamma_b} = \sqrt{\alpha_b E} L_{\Gamma_b} - \sigma_{\Gamma_b} \pi/2$ . For  $\bar{\rho}_{1b} * \bar{\rho}_{1a}(E)$ , we just interchange  $a$  and  $b$ .

The formula for  $\bar{\rho}_{1a} * \bar{\rho}_{1b}(E)$  still has the same basic structure, but should obviously use the distinct periodic orbits for particles  $a$  and  $b$ . In particular, Eqs. (12) and (18) still apply, but with two important differences. Firstly, the double sums over periodic orbits are now labelled by the distinct periodic orbits of the two particles. Secondly, the energy partition will change due to differing masses. The criterion of stationary phase will still specify that the two particles have the same period, but relations such as (27) and (28) do not apply since they assume equal masses. The generalisations are rather straight-forward to determine. For example, the saddle energies (28) are replaced by

$$\frac{E_0}{E} = \frac{m_a L_{\Gamma_a}^2}{m_a L_{\Gamma_a}^2 + m_b L_{\Gamma_b}^2}, \quad \frac{E - E_0}{E} = \frac{m_b L_{\Gamma_b}^2}{m_a L_{\Gamma_a}^2 + m_b L_{\Gamma_b}^2} \quad (56)$$

while the general dynamical expression for billiards (30) is replaced by

$$\begin{aligned}\bar{\rho}_{1a} * \bar{\rho}_{1b}(E) &\approx \frac{(2E)^{3/4} \sqrt{\alpha_a \alpha_b} \hbar}{(2\pi)^{3/2}} \\ &\sum_{\Gamma_a, \Gamma_b} \frac{L_{\Gamma_a} L_{\Gamma_b}}{(m_a L_{\Gamma_a}^2 + m_b L_{\Gamma_b}^2)^{5/4}} A_{\Gamma_a}(E_0) A_{\Gamma_b}(E - E_0) \\ &\cos\left(\sqrt{\alpha_a L_{\Gamma_a}^2 + \alpha_b L_{\Gamma_b}^2} \sqrt{E} - (\sigma_{\Gamma_a} + \sigma_{\Gamma_b}) \frac{\pi}{2} - \frac{\pi}{4}\right).\end{aligned}\quad (57)$$

In the special case of identical particles, it is simple to check that this expression reduces to (30). For lack of an immediate physical context, we do not explore this case any further.

## IX. APPENDIX B: SPURIOUS END-POINT CONTRIBUTIONS FOR THE CARDIOID

Here we evaluate the cross term integrals exactly for isolated periodic orbits. This allows us to do an asymptotic expansion to explicitly demonstrate that the additional endpoint contributions not included are spurious. We must evaluate the integral

$$\bar{\rho}_1 * \tilde{\rho}_1(E) = \int_0^E d\epsilon \bar{\rho}_1(\epsilon) \tilde{\rho}_1(E - \epsilon) \quad (58)$$

where  $\bar{\rho}_1(\epsilon)$  is given by the Weyl expansion (3) and  $\tilde{\rho}_1(\epsilon)$  for a billiard with isolated orbits is given by

$$\tilde{\rho}_1(\epsilon) \approx \frac{\alpha^{1/2}}{2\pi\sqrt{\epsilon}} \sum_{\Gamma} \frac{L_{\Gamma}}{\sqrt{|\det(\tilde{M}_{\Gamma} - I)|}} \cos\left(\sqrt{\alpha\epsilon} L_{\Gamma} - \sigma_{\Gamma} \frac{\pi}{2}\right). \quad (59)$$

This gives

$$\begin{aligned}\bar{\rho}_1 * \tilde{\rho}_1(E) &\approx \sum_{\Gamma} \frac{L_{\Gamma}}{\sqrt{|\det(\tilde{M}_{\Gamma} - I)|}} \\ &\left( \alpha^{3/2} \frac{\mathcal{A}}{8\pi^2} I_1 - \alpha \frac{\mathcal{L}}{16\pi^2} I_2 + \alpha^{1/2} \frac{\mathcal{K}}{2\pi} I_3 \right)\end{aligned}\quad (60)$$

where

$$\begin{aligned}I_1 &= \int_0^E d\epsilon \frac{1}{\sqrt{E - \epsilon}} \cos\left(\sqrt{\alpha(E - \epsilon)} L_{\Gamma} - \sigma_{\Gamma} \frac{\pi}{2}\right), \\ I_2 &= \int_0^E d\epsilon \frac{1}{\sqrt{\epsilon}} \frac{1}{\sqrt{E - \epsilon}} \cos\left(\sqrt{\alpha(E - \epsilon)} L_{\Gamma} - \sigma_{\Gamma} \frac{\pi}{2}\right), \\ I_3 &= \frac{1}{\sqrt{E}} \cos\left(\sqrt{\alpha E} L_{\Gamma} - \sigma_{\Gamma} \frac{\pi}{2}\right).\end{aligned}\quad (61)$$

If we evaluate the first two integrals exactly, we get

$$I_1 = \frac{2}{\alpha^{1/2} L_{\Gamma}} \left( \cos\left(\Phi_{\Gamma} + \phi_{\Gamma} - \frac{\pi}{2}\right) - \cos\left(\phi_{\Gamma} - \frac{\pi}{2}\right) \right) \quad (62)$$

and

$$\begin{aligned}I_2 &= \pi \cos \phi_{\Gamma} J_0(\Phi_{\Gamma}) - \pi \sin \phi_{\Gamma} \mathbf{H}_0(\Phi_{\Gamma}) \\ &\approx \sqrt{\frac{2\pi}{\sqrt{\alpha E} L_{\Gamma}}} \cos\left(\Phi_{\Gamma} + \phi_{\Gamma} - \frac{\pi}{4}\right) \\ &- \frac{2}{\sqrt{\alpha E} L_{\Gamma}} \cos(\phi_{\Gamma}) + \dots\end{aligned}\quad (63)$$

where  $\Phi_\Gamma = \sqrt{\alpha E} L_\Gamma$ ,  $\phi_\Gamma = -\sigma_\Gamma \pi/2$ ,  $J_0$  is a zero-order Bessel function and  $\mathbf{H}_0$  is a zero-order Struve function. In the second line of (63), we have used the asymptotic expansions of these two functions.

In both  $I_1$  and  $I_2$ , we note that asymptotically there are terms with two distinct structures. The first are terms which are sinusoidal in  $\sqrt{E}$  and correspond exactly to what was used as the cross term for the cardioid (*i.e.* Eqs. (25) and 26)). There are also terms which are nonsinusoidal in  $E$ . In  $I_1$ , this comes directly from the upper endpoint of the integral while in  $I_2$  it comes from the expansion of the Struve function. In each term, the nonsinusoidal terms arise from the endpoint around  $\epsilon = E$  which, as we argued in section IV D, corresponds to an unphysical situation. Therefore, keeping only the asymptotically appropriate term (*i.e.* the oscillatory one) yields the correct behaviour for the cross term.

A similar analysis would yield similar results for the spurious endpoint contributions in the cross term of the disk billiard and the dynamical term of either billiard.

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