Boundary contributions to the hypervirial theorem

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It is shown that, under certain boundary conditions, the virial theorem has to be modified. We analyze the origin of the extra term and compute it for particular examples. The Coulomb and harmonic oscillator with the point interaction have been studied in light of this generalization of the virial theorem.

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I. INTRODUCTION

The virial theorem in classical and quantum mechanics has been known for a very long time. It appears for the first time in the classical works of Clausius in 1870 [1] in the context of statistical mechanics. See also [2] for a historical account. Already in 1930, Fock [3] derived the quantum mechanical version of the theorem that relates the expectation value of the kinetic energy in an eigenstate ψ_n of the Hamiltonian to that of the Clausius virial function. More precisely,

$$2\langle \psi_n | T \psi_n \rangle = \langle \psi_n | (\mathbf{x} \cdot \nabla V) \psi_n \rangle. \tag{1}$$

The theorem was initially derived for eigenstates of nonrelativistic Hamiltonians in flat space with potentials depending only on the position, but it can be generalized in many different ways: for relativistic systems, for a particle moving in an electromagnetic field, to many-body systems, mixed states, and for compact space [4–11] to mention a few. Also, a local version of the virial theorem for systems of fermions has recently been formulated (see Refs. [12,13]).

Another kind of generalization is the hypervirial theorem [14] that works, in principle, for any operator G and can be written

$$\langle \psi_n | [H, G] \psi_n \rangle = 0.$$

The virial theorem in dimension d is recovered when

$$G = -\frac{i}{\hbar}\mathbf{x} \cdot \mathbf{p} - \frac{d}{2}.$$

In this paper we discuss the domain of validity of these theorems and, for the cases in which they fail, we show how to modify them by adding the appropriate term related to the boundary conditions of the quantum system. In this way we extend the applicability of the virial and hypervirial theorem to these systems.

In concrete terms, assume that the Hamiltonian is defined over a domain D(H). ψ_n necessarily belongs to D(H), but it may happen that $G\psi_n \notin D(H)$. In this case the hypervirial theorem has to be modified by the addition of an extra term \mathcal{A} :

$$\langle \psi_n | [H, G] \psi_n \rangle = \mathcal{A}, \tag{2}$$

with

$$\mathcal{A} \equiv \langle \psi_n | (H_* - E_n) G \psi_n \rangle,$$

where E_n is the energy eigenvalue of ψ_n and H_* is the closed extension of H to a domain that includes $G\psi_n$. Of course, the new term cancels if $H_* = H$ [i.e., when $G\psi_n \in D(H)$], but if that is not the case it is different from zero in general. We will see that \mathcal{A} can be computed as the integration of a pure derivative term and, therefore, it depends only on the value of the wave function and its derivatives at the boundary. This fact is reminiscent of the analogous objects in quantum field theory, like the chiral anomaly [15]. We would like to remark that an additional term for the virial (or hypervirial) theorem in classical mechanics may appear also when the orbits are not bounded in phase space.

This paper is organized in the following fashion: In the next section, we discuss a simple example to motivate our study and obtain the generalization of the hypervirial theorem. Then we give two examples of systems where this generalization is needed, in Sec. III we discuss the hydrogen-atom problem with point interaction in three dimensions, while Sec. IV is devoted to the isotropic harmonic oscillator. Finally, we sketch our conclusions in Sec. V.

II. GENERALIZED HYPERVIRIAL THEOREM

Before discussing the required generalization of the hypervirial theorem we would like to motivate our study by considering a very simple example.

Consider a free particle in one dimension restricted to move in $[0,\infty)$ and subject to Robin boundary conditions [i.e., $\psi'(0) + \alpha \psi(0) = 0$ with $\alpha > 0$]. In this case the free Hamiltonian has a single eigenfunction:

$$\psi_0(x) = \sqrt{2\alpha}e^{-\alpha x},$$

with eigenvalue¹

$$E_0 = -\frac{\hbar^2 \alpha^2}{2m}.$$

$$\langle \psi | T \psi \rangle = \frac{\hbar^2}{2m} [\langle \psi' | \psi' \rangle + \psi(0)^* \psi'(0)] = \frac{\hbar^2}{2m} [\langle \psi' | \psi' \rangle - \alpha | \psi(0)|^2],$$

and the second term on the right provides a negative contribution.

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¹At first sight it may seem strange that $T = p^2/(2m)$ has a negative eigenvalue. Note, however, that p is not symmetric with Robin boundary conditions and, in fact, one has

If we use (1) to compute the expectation value of the kinetic energy in this state we obtain that it vanishes, which is in contradiction with the real result

$$\langle \psi_0 | T \psi_0 \rangle = E_0. \tag{3}$$

The reason for this apparent contradiction is the fact that the domain of the Hamiltonian is not preserved by the generator of the scale transformation

$$G = -\frac{i}{\hbar}xp - \frac{1}{2},\tag{4}$$

and the virial theorem has to be modified.

To understand the origin of this modification let us consider in detail the derivation of the virial theorem. We assume a quantum system with infinite-dimensional state space and a self-adjoint Hamiltonian with domain D(H). If $G\psi_n \in D(H)$ we have

$$\langle \psi_n | [H, G] \psi_n \rangle = \langle \psi_n | HG \psi_n \rangle - \langle \psi_n | GH \psi_n \rangle = \langle H\psi_n | G\psi_n \rangle - \langle \psi_n | GH \psi_n \rangle = 0,$$
 (5)

where in the second line we have used the fact that H is self-adjoint and that ψ_n is its eigenstate.

The problems appear if $G\psi_n \notin D(H)$. In this case the expressions in the first line of (5) do not make sense unless we extend H to a larger domain that includes $G\psi_n$. In principle it can be done in an arbitrary way but, in this case, there is a well-defined prescription to obtain the required extension. First we restrict H to an appropriate dense subspace S, typically to functions whose support does not include the boundary. If we define $\hat{H} = H|_S$, the extension H_* is the adjoint of the previous restriction (i.e., $H_* = \hat{H}^+$). We will assume that the restriction is closed (which could always be achieved by taking the closure of its graph) and then we know $H_*^+ = \hat{H}$.

With the previous machinery we can rewrite (5) in a meaningful way

$$\begin{aligned} \langle \psi_n | [H_*, G] \psi_n \rangle &= \langle \psi_n | H_* G \psi_n \rangle - \langle \psi_n | G H \psi_n \rangle \\ &= \langle \psi_n | (H_* - E_n) G \psi_n \rangle \equiv \mathcal{A}, \end{aligned}$$

and obtain the generalized virial theorem as announced in the introduction. Looking at the last line of this expression one could be tempted to take the adjoint of the operator H_* and make it act from the left side of the scalar product. However, this is not possible in general because the adjoint of H_* is \hat{H} with a domain smaller than H that may not contain ψ_n . Of course, if it so happens that $\psi_n \in S$ the additional term cancels.

We would like to show now that, under certain assumptions (that we actually meet in our examples), the extra term is a boundary term; that is, it depends only on the value of the functions and their derivatives at the boundary.

First of all, we rewrite the extra term in a way that is appropriate for wave functions that are not necessarily eigenvectors of H:

$$\mathcal{A} = \langle \psi_n | H_* G \psi_n \rangle - \langle H \psi_n | G \psi_n \rangle.$$

Next, we will show that the value of \mathcal{A} does not change if we replace ψ_n by $\psi_n + \chi$ with $\chi \in S$ and $G\chi \in S$. The difference reads

$$\Delta \mathcal{A} = \langle \psi_n | H_* G \chi \rangle - \langle H \psi_n | G \chi \rangle + \langle \chi | H_* G \psi_n \rangle - \langle H \chi | G \psi_n \rangle + \langle \chi | H_* G \chi \rangle - \langle H \chi | G \chi \rangle.$$
 (6)

Now, given that $G\chi \in D(H)$, we can replace H_* by H in the first term of this expression and, if we take its adjoint, it cancels the second term. For the very same reason the last two terms cancel. Also in the third term we can take the adjoint of H_* , because χ is in the domain of $H_*^+ = \hat{H}$, and therefore it cancels the fourth term and ΔA vanishes.

Then, if as we supposed before, the elements of S are wave functions whose support does not contain the boundary, we deduce that the extra term A is invariant under deformations of ψ_n that do not affect the values of the function and its derivatives at the boundary or, in other words, the extra term depends only on the latter.

As an illustration of these results and before discussing the examples of physical interest, we will compute the new term for the simple system introduced at the beginning of this section. One immediately gets

$$\mathcal{A} = \frac{\hbar^2}{2m} \int_0^\infty \left[G\psi_0(x) \partial_x^2 \psi_0^*(x) - \psi_0^*(x) \partial_x^2 G\psi_0(x) \right] dx$$
$$= \frac{\hbar^2}{2m} \int_0^\infty \partial_x \left[G\psi_0(x) \partial_x \psi_0^*(x) - \psi_0^*(x) \partial_x G\psi_0(x) \right] dx,$$
(7)

which is a pure boundary term (for any operator G), as stated. In our particular case it amounts to

$$\mathcal{A} = \frac{\hbar^2}{4m} [3\psi_0(0)\psi_0'(0)^* - \psi_0(0)\psi_0'(0)^*]$$
$$= -\frac{\hbar^2 \alpha}{2m} \psi_0(0)^2 = -\frac{\hbar^2 \alpha^2}{m}.$$
 (8)

Now taking into account that [G,H] = 2T and plugging these expressions into the generalized hypervirial theorem (2), we get

$$2\langle\psi_0|T\psi_0\rangle = -\frac{\hbar^2\alpha^2}{m},$$

from which (3) follows.

We have seen that, if we modify appropriately the hypervirial theorem it can be applied to situations in which the action of the operator G on the energy eigenstate puts it out of the domain of the Hamiltonian. In the following sections we will apply this idea to the Coulomb problem and later on to the isotropic harmonic oscillator.

III. COULOMB POTENTIAL

In this section we discuss the virial theorem for the Coulomb problem in three dimensions. We shall see that, with the chosen boundary conditions at the origin, the extra term is required to cancel the divergence that appears in the application of the virial theorem.

Consider the Hamiltonian

$$H = \frac{\mathbf{p}^2}{2m} - \frac{k}{r}.\tag{9}$$

If we separate the angular variables from the radial variable, we obtain the Hamiltonian for the latter

$$H_{l} = \frac{\hbar^{2}}{2m} \left(-\frac{1}{r} \frac{d^{2}}{dr^{2}} r + \frac{l(l+1)}{r^{2}} - \frac{\xi}{r} \right), \tag{10}$$

which depends on the angular momentum number l. Here $\xi = 2mk/\hbar^2$.

We can simplify the Hamiltonian by performing a similarity transformation $\phi = r\psi$. Therefore, the new wave functions ϕ are square integrable in $[0,\infty)$ with respect to the Lebesgue measure. The new Hamiltonian is then

$$\hat{H}_{l} = \frac{\hbar^{2}}{2m} \left(-\frac{d^{2}}{dr^{2}} + \frac{l(l+1)}{r^{2}} - \frac{\xi}{r} \right), \tag{11}$$

and it is symmetric with respect to the Lebesgue measure in $[0,\infty)$ when acting on normalizable wave functions whose support does not contain the origin.

In order to have a self-adjoint Hamiltonian we have to look for extensions of \hat{H}_l with such a property. This issue has been analyzed in detail in [16]. The result is that, if $l \ge 1$, there is a single self-adjoint extension of \hat{H}_l whose domain is the wave functions that vanish at 0, while \hat{H}_0 has an infinite number of self-adjoint extensions characterized by a single parameter α . If we call H^{α} such a self-adjoint extension, its domain is given by

$$\begin{split} D(H^{\alpha}) &= \{ \phi \in D(\hat{H}_0^+) | \quad \text{as} \quad r \to 0, \\ \phi(r) &= \phi(0) [1 - \xi r \ln(|\xi|r) + \alpha r + o(r)] \}. \end{split}$$

where the little-o notation has been used. The case of wave functions that vanish at the origin (the standard boundary conditions for the hydrogen atom) is recovered when $\alpha \to -\infty$ and $\phi(0) \to 0$, while the product $\phi(0)\alpha$ remains finite.

The point spectrum of H^{α} is obtained from the solutions in λ to the equation

$$F_C(\lambda) \equiv \Psi(1-\lambda) - \ln|\lambda| + \frac{1}{2\lambda} + 2\gamma - 1 = \frac{\alpha}{\varepsilon}, \quad (12)$$

such that $\xi/\lambda \geqslant 0$. Here γ is the Euler's constant and Ψ is the digamma function (i.e., the logarithmic derivative of the gamma function). For a solution λ_n of the previous equation the eigenvalue of H^{α} is

$$E_n = -\frac{mk^2}{2\hbar^2} \frac{1}{\lambda_n^2}.$$

In Fig. 1 we represent graphically the solutions of (12) for λ as the intersection of the branches of the curve with the line of constant value α/ξ . On the negative side, $F(\lambda)$ has an asymptotic value $2\gamma - 1 \approx 0.1544$.

Note that, even in the presence of a repulsive Coulomb potential, k < 0, if we take $\alpha/\xi < 2\gamma - 1$ there is a single bounded state associated with the negative solution for λ of the equation $F_C(\lambda) = \alpha/\xi$.

The eigenfunction associated with the eigenvalue E_n is given by the Whittaker function [17] conveniently normalized

$$\phi_n(r) = N_n W_{\lambda_n, 1/2}(\xi r/\lambda_n),$$

with the normalization constant given by

$$N_n^2 = \frac{\xi |\Gamma(-\lambda_n)|^2}{2\lambda_n \Psi'(-\lambda_n) + 2 - \lambda_n^{-1}},$$
 (13)

where the prime denotes derivative.

Now we want to apply the generalized form of the virial theorem to this system. We take the operator $G = -\frac{i}{\hbar} \mathbf{x} \cdot \mathbf{p} - \frac{3}{2}$, then the modified hypervirial theorem reduces to

$$2\langle \psi_n | T\psi_n \rangle + \langle \psi_n | V\psi_n \rangle = \mathcal{A}.$$

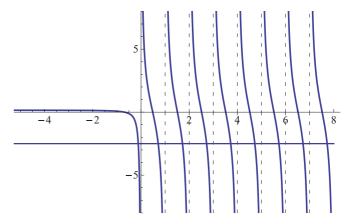


FIG. 1. (Color online) In the figure we represent λ on the horizontal axis and plot $F_C(\lambda)$. The point-spectrum of H^{α} is obtained from the values of λ at the intersection of the graph of F_C with the horizontal line at α/ξ . Only those values of λ with the same sign as ξ are allowed.

Using the energy eigenvalue and integrating out the angular coordinates, we can write for the radial wave function

$$\mathcal{A} - \left\langle \phi_n \left| \frac{k}{r} \phi_n \right\rangle = 2E_n. \tag{14}$$

In this expression we can immediately see the necessity of an additional contribution to the virial theorem: the expectation value of the potential term diverges logarithmically at the origin (recall that we must use the Lebesgue measure in $[0,\infty)$ while the right-hand side is finite). In fact, we shall see that an analogous divergence appears in the extra contribution so that both of them cancel to yield the correct finite result.

We will sketch how to verify it. We first introduce a cutoff ϵ that removes the origin from the integration region. The extra term only depends on the boundary conditions of the wave function. We can take it from (7) and, in its *regularized* version, it reads

$$\mathcal{A}_{\epsilon} = -\frac{\hbar^2}{2m} |\phi_n(0)|^2 \left[\xi \ln(|\xi|\epsilon) + \xi - \alpha + \cdots \right],$$

where the dots stand for terms that vanish when $\epsilon \to 0$. We now compute the potential term

$$\left\langle \phi_n \left| \frac{k}{r} \phi_n \right\rangle_{\epsilon} = N^2 k \int_{\epsilon}^{\infty} \frac{1}{r} W_{\lambda_n, 1/2} (\xi r / \lambda_n)^2 dr. \right.$$

To evaluate the expectation value we use

$$\int_{\epsilon}^{\infty} \frac{1}{x} W_{\lambda, 1/2}(x)^2 dx = \frac{-1}{\Gamma(-\lambda)^2} [\ln \epsilon - \lambda \Psi'(1 - \lambda) + \Psi(1 - \lambda) + 2\gamma + \cdots],$$

where, as before, the dots represent terms that vanish when $\epsilon \to 0$. The integral can be obtained using the identity [18]

$$\int \frac{1}{x} W_{\mu,\sigma}(x) W_{\rho,\sigma}(x) dx$$

$$= \frac{1}{\mu - \rho} [W_{\mu,\sigma}(x) W'_{\rho,\sigma}(x) - W'_{\mu,\sigma}(x) W_{\rho,\sigma}(x)] \quad (15)$$

and the series expansion of $W_{\mu,\sigma}$ [17].

Putting everything together, and given

$$W_{\lambda 1/2}(0) = 1/\Gamma(1-\lambda)$$

and (12) we obtain,

$$A_{\epsilon} - \left\langle \phi_n \left| \frac{k}{r} \phi_n \right\rangle_{\epsilon} = -\frac{mk^2}{\hbar^2} \frac{1}{\lambda_n^2} + \cdots,$$

in agreement with (14).

IV. HARMONIC OSCILLATOR

In the previous section we saw that the verification of the virial theorem is plagued by infinities. This fact is not inherent to the generalized virial theorem itself but is rather related to the singularity in the Coulomb potential. To illustrate this fact we shall discuss now a new instance of the generalized virial theorem in three dimensions that is free of these divergences.

We consider a three-dimensional harmonic oscillator given by the Hamiltonian

$$H = \frac{\mathbf{p}^2}{2m} + \frac{1}{2}m\omega^2 r^2.$$
 (16)

As before, we factorize the angular part and study the radial Hamiltonian for the radial part $\psi(r)$ of the wave function of angular momentum number l:

$$H_{l} = \frac{\hbar^{2}}{2m} \left(-\frac{1}{r} \frac{d^{2}}{dr^{2}} r + \frac{l(l+1)}{r^{2}} \right) + \frac{1}{2} m \omega^{2} r^{2}.$$

Or, acting on $\phi(r) = r\psi(r)$,

$$\hat{H}_{l} = \frac{\hbar^{2}}{2m} \left(-\frac{d^{2}}{dr^{2}} + \frac{l(l+1)}{r^{2}} \right) + \frac{1}{2} m \omega^{2} r^{2},$$

which is a symmetric operator with respect to the Lebesgue measure in $[0,\infty)$ when acting on functions whose support does not include the point 0.

In order to obtain a self-adjoint operator we have to extend \hat{H}_l to a larger domain. As was the case in the previous section if $l \ge 1$, the only self-adjoint extension is to functions that vanish at the origin. In this case the extra term cancels and the standard virial theorem holds. However, if l = 0, we have a whole family of self-adjoint extensions parametrized by β . If we call H^β such an extension its domain corresponds to Robin boundary conditions; that is,

$$D(H^{\beta}) = \{ \phi \in D(\hat{H}_0^+) | \phi'(0) + 2\beta\phi(0) = 0 \}. \tag{17}$$

The normalized eigenfunctions of H^{β} can be written in terms of the Whittaker function as

$$\phi_E(r) = \frac{N}{\sqrt{\sigma r}} W_{\kappa, 1/4}(\sigma^2 r^2), \tag{18}$$

where $\sigma=(m\omega/\hbar)^{1/2}$, $\kappa=\frac{E}{2\hbar\omega}$, and the normalization constant N is obtained from the condition

$$\frac{N^2}{2\sigma} \int_0^\infty \frac{1}{x} W_{\kappa, 1/4}(x)^2 dx = 1.$$

The integral can be performed using (15) and finally we get

$$N^2 = \frac{2\sigma}{\pi} \frac{\Gamma(3/4-\kappa)\Gamma(1/4-\kappa)}{\Psi(3/4-\kappa)-\Psi(1/4-\kappa)}.$$

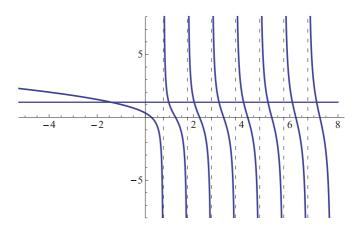


FIG. 2. (Color online) In the figure we represent $F_H(\kappa)$ with κ on the horizontal axis. The point spectrum of H^{β} is obtained from the values of κ at the intersection of the graph of F_H with the horizontal line at β/σ . The possible values of κ are given as the intersection of the curve with the horizontal line at β/σ .

Only for specific values of κ do the eigenstates (18) belong to domain (17), demanding $\phi_E \in D(H^{\beta})$ we obtain the equation for κ :

$$F_H(\kappa) \equiv \frac{\Gamma(3/4 - \kappa)}{\Gamma(1/4 - \kappa)} = \frac{\beta}{\sigma}.$$
 (19)

The solutions are represented in Fig. 2.

In general the eigenvalues $E_n = 2\kappa_n \hbar \omega$ must be computed numerically. One particular case in which we have analytic solutions is $\beta = 0$, which corresponds to Neumann boundary conditions. In this case we get $E_n = n\hbar\omega + 1/2$. We also have exact results for $\beta = -\infty$, Dirichlet boundary conditions and $E_n = n\hbar\omega + 3/2$. The latter corresponds to the harmonic oscillator without point interaction. Note also that for β/σ > $\Gamma(3/4)/\Gamma(1/4) \approx 0.337\,989\ldots$, there is an eigenstate with negative energy. Of course, as discussed in the footnote of Sec. II, there is not a contradiction between the negative eigenvalue and the fact that H^{β} is the sum of squares of operators. The point is that, with Robin boundary conditions, some of those operators are not self-adjoint and there is not any reason to argue that H^{β} is semipositive definite. The same applies to the Coulomb potential with repulsive interaction that we discussed in the previous section.

The generalized virial theorem, in terms of the radial wave function ϕ_n , reads

$$\langle \phi_n | T \phi_n \rangle - \langle \phi_n | V \phi_n \rangle = \mathcal{A},$$

where we have taken G as before. Using the energy eigenvalue we can rewrite the identity above in the following way:

$$2\langle \phi_n | V \phi_n \rangle + \mathcal{A} = E_n, \tag{20}$$

and this is the expression that we will check.

We first compute the additional term

$$\mathcal{A} = \frac{\hbar^2}{4m} \phi_n(0) \phi'_n(0)^* = -\frac{\hbar^2 \beta}{2m} |\phi_n(0)|^2$$
$$= -N^2 \frac{\pi \sigma \hbar^2}{2m} \frac{1}{\Gamma(3/4 - \kappa_n) \Gamma(1/4 - \kappa_n)}.$$
 (21)

In the last equality, condition (19) was used.

The expectation value of the potential can be written as

$$\langle \phi_n | V \phi_n \rangle = N^2 \frac{\sigma \hbar^2}{4m} \int_0^\infty W_{\kappa_n, 1/4}(z)^2 dz, \tag{22}$$

where the change of variable $z = \sigma^2 x^2$ was performed.

Now we can compute the previous integral with the help of the identity

$$zW_{\kappa,1/4}(z) = W_{\kappa+1,1/4}(z) + 2\kappa W_{\kappa,1/4}(z) + (3/4 - \kappa)(1/4 - \kappa)W_{\kappa-1,1/4}(z), \quad (23)$$

which we apply to one of the factors of the square inside the integral. We perform the resulting integrals using (15) and finally we obtain

$$\langle \phi_n | V \phi_n \rangle = \frac{\sigma^2 \hbar^2 \kappa_n}{m} + N^2 \frac{\pi \sigma \hbar^2}{4m} \frac{1}{\Gamma(3/4 - \kappa_n) \Gamma(1/4 - \kappa_n)}.$$
(24)

If we plug this expression and (21) into (20) we can see that the identity holds.

V. CONCLUSION

We proved that, when the dilations symmetry does not keep invariant the domain of definition of the Hamiltonian, the exact virial theorem has an extra term that accounts for this fact. We have shown that this term is the integral of a pure derivative and, therefore, its value depends only of the behavior of the wave function at the boundary. We have verified this generalization of the virial theorem in the cases of Coulomb and harmonic oscillator with point interaction in three dimensions.

As an extension of our work, it could be interesting to study if a similar phenomenon occurs for the local virial theorem for fermions with a harmonic potential [12,13] and the boundary conditions used in this paper.

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