## Quantum Solutions

- 1. (a) Since at t=0 the system is in an eigenstate of the parity operator, we know that the associated wavefunction is either symmetric or anti-symmetric at this time. Moreover, since the potential is symmetric we have that the Hamiltonian operator commutes with  $\hat{\Pi}$ . Hence, the symmetry of the state will remain the same as the system evolves in time. Now the operator A is anti-symmetric (that is,  $\hat{\Pi} A \hat{\Pi} = -A$ ). Therefore if  $|\psi_{\pm}\rangle$  is an eigenstate of  $\hat{\Pi}$  with eigenvalue  $\pm 1$ , then  $A |\psi_{\pm}\rangle$  is an eigenstate of  $\hat{\Pi}$  with eigenvalue  $\mp 1$ . Thus  $\langle A \rangle_{|\psi_{\pm}\rangle} = \langle \psi_{\pm} | A | \psi_{\pm} \rangle = 0$ , from which our result follows.
- (b) Since  $\langle \hat{\Pi} \rangle_{|\psi\rangle} = 0$ , we have that  $|\psi\rangle = \frac{1}{\sqrt{2}}(|\psi_{+}\rangle + |\psi_{-}\rangle)$ , where  $|\psi_{\pm}\rangle$  is an eigenstate of  $\hat{\Pi}$  with eigenvalue  $\pm 1$ . In order to obtain the lowest value of  $\langle H \rangle_{|\psi\rangle}$  in such a state, simple choose  $|\psi_{+}\rangle$  to be the ground state (which is the symmetric state with the lowest energy, namely  $E_{0} = \frac{1}{2}\hbar\omega$ ) and  $|\psi_{-}\rangle$  to be the first excited state (which is the anti-symmetric state with the lowest energy, namely  $E_{1} = \frac{3}{2}\hbar\omega$ ). A simple calculation then yields  $\langle H \rangle_{|\psi\rangle} = (E_{0} + E_{1})/2 = \hbar\omega$ .
- 2. (a) Since A and B are Hermitian, we have that  $(A+B)^2$  is also Hermitian. But 2AB is Hermitian if and only if A commutes with B. Therefore we have that  $(A+B)^2=A^2+B^2+2AB=2AB$  (where the first equality uses  $[A,B]=\hat{0}$ ), and hence  $A^2+B^2=\hat{0}$ . Taking the expectation of both sides of this last equation in an arbitrary state  $|\psi\rangle$  yields  $\langle A\psi|A\psi\rangle+\langle B\psi|B\psi\rangle=0$ , where we have used the Hermiticity of A and B. Since both terms on the LHS of this equation are  $\geq 0$  for any A and  $|\psi\rangle$ , we see that  $\langle A\psi|A\psi\rangle=\langle B\psi|B\psi\rangle=0$ . But  $\langle \phi|\phi\rangle=0$  if and only if the vector  $|\phi\rangle$  is the zero vector  $|0\rangle$ . Hence  $A|\psi\rangle=B|\psi\rangle=|0\rangle$  for all states  $|\psi\rangle$ , which shows that  $A=B=\hat{0}$ .
- (b) Let  $|\phi\rangle$  be a unit vector in  $\mathcal{H}$  which is orthogonal to  $|\psi\rangle$ . (The vector  $|\phi\rangle$  is unique up to an overall phase.) Then  $\{|\psi\rangle, |\phi\rangle\}$  is an orthonormal basis for  $\mathcal{H}$ . We can now (for example) choose  $A = \alpha(|\psi\rangle\langle\phi| + |\phi\rangle\langle\psi|)$  and  $B = \beta(|\psi\rangle\langle\psi| |\phi\rangle\langle\phi|)$ , where  $\alpha$  and  $\beta$  are non-zero real numbers. The eigenvalues of A are  $\pm \alpha$ , and the state of the system after the first measurement is  $|\chi_{\pm}\rangle = \frac{1}{\sqrt{2}}(|\psi\rangle \pm |\phi\rangle)$  if the value  $\pm \alpha$  is obtained (each occurs with probability 1/2). We want to show that if we now measure B (whose eigenvalues are  $\pm \beta$ ), there will be a non-zero probability of obtaining the value  $-\beta$ , in which case the state of the system after the measurement will be  $|\phi\rangle$  (which is orthogonal to  $|\psi\rangle$ ). But it is immediate from the form of the states  $|\chi_{\pm}\rangle$  that the probability of obtaining  $-\beta$  in the measurement of B is 1/2, regardless of the outcome of the first measurement.
- 3. (a) For any Hamiltonian operator H and any state  $|\psi\rangle$ , the variational theorem tells us that  $E_0 \leq \langle \psi | H | \psi \rangle$ , where  $E_0$  is the lowest energy eigenvalue of H. In our case, if we make the choice  $|\psi\rangle = |\psi_0^{(0)}\rangle$  we immediately obtain  $\langle \psi | H | \psi \rangle = E_0^{(0)} + \lambda E_0^{(1)}$  (since  $H_0 | \psi_0^{(0)} \rangle = E_0^{(0)} | \psi_0^{(0)} \rangle$  and  $E_0^{(1)} = \langle \psi_0^{(0)} | H_1 | \psi_0^{(0)} \rangle$ ).
- (b) Plugging the expansions for  $|\psi_n\rangle$  and  $E_n$  into the time-independent Schrödinger equation  $H|\psi_n\rangle=E_n|\psi_n\rangle$  and grouping terms of order  $\lambda^m$   $(m\geq 1)$  yields

$$H_0|\psi_n^{(m)}\rangle + H_1|\psi_n^{(m-1)}\rangle = \sum_{k=0}^m E_n^{(k)}|\psi_n^{(m-k)}\rangle$$
.

Taking the inner product of both sides with  $|\psi_n^{(0)}\rangle$  (and using  $\langle \psi_n^{(0)} | \psi_n^{(m)} \rangle = 0$  for all  $m \geq 1$  and the Hermiticity of  $H_0$ ) then gives the desired result.

4. (a) The energy levels of the particle in the two-dimensional box is just the sum of the energy levels of two one-dimensional square-well problems, one of width L and the other of width L/(N+1). Thus we have

$$E_{n_x,n_y} = \frac{\hbar^2 \pi^2}{2ML^2} (n_x^2 + (N+1)^2 n_y^2)$$

(where  $n_x, n_y = 1, 2, ...$ ). The result then follows by a simple computation.

- (b) For a state with total angular momentum quantum number  $\ell$ , the radial wavefunction satisfies  $R(r) \sim r^{\ell}$  as  $r \to 0$ . Hence we have  $\ell = 3$ . Moreover, if the state has magnetic quantum number m, the  $\phi$ -dependence of its wavefunction if  $e^{im\phi}$ . Hence we have m = 2. Finally, by substituting R(r) in the (time-independent) radial Schrödinger equation and solving for V(r) at large r, we find that V(r) goes to a constant C as  $r \to \infty$ . Hence, there will be a continuum in the spectrum of H for  $E \geq C$ .
- 5. (a) The <sup>11</sup>B atom has 5 protons, 6 neutrons, and 5 electrons, each of which is a fermion. Hence <sup>11</sup>B is made of an even number of fermions, and is therefore a boson.
- (b) In the absence of Coulomb repulsion (and spin-dependent forces), the problem reduces to 5 independent Hydrogenic atoms, each with nuclear charge Z=5, where we also have to use the Pauli Exclusion Principle in order to properly distribute the spin one-half electrons in the Hydrogenic levels. In particular, 2 electrons will be in the 1S state, 2 in the 2S state, and one in the 2P state. Hence, in our approximation, the ground state energy of the Boron atom will be  $E_0 = -Z^2(13.6 \text{ eV})(1+1+1/4+1/4)=-935 \text{ eV}$ .
- (c) The degeneracy of the ground state is just the degeneracy of the 2P state for an electron, which is 6 since the quantum number  $m_{\ell}$  of the 2P electron can be 1,0, or -1 (since this is an  $\ell = 1$  state) and the quantum number  $m_s$  of this electron can be  $\pm 1/2$  (since the electron has s = 1/2).
- (d) Each of the spatial 2P basis states  $|\psi_{2,1,m_\ell}\rangle$  ( $m_\ell=0,\pm 1$ ) is not rotationally symmetric, and the same will be true of any linear combination of these linearly independent degenerate states.
- (e) Each Coulomb repulsion term increases the potential energy of the system, and hence the ground state energy will increase (but decrease in magnitude).