

PH423 Assignment 4

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Question 2.

[Sankalp: Got it]

(a) The given Hamiltonian $H(\epsilon)$ is

$$H(\epsilon) = V_0 \cdot \begin{pmatrix} (1-\epsilon) & 0 & 0 \\ 0 & 1 & \epsilon \\ 0 & \epsilon & 2 \end{pmatrix} \quad (1)$$

and the unperturbed Hamiltonian H_0

$$H_0 = H(0) = V_0 \cdot \begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 2 \end{pmatrix} \quad (2)$$

is diagonal, so the eigenvectors of the unperturbed system are simply $(1, 0, 0)$, $(0, 1, 0)$, and $(0, 0, 1)$ with eigenvalues V_0 , V_0 , and $2V_0$ respectively.

The perturbation H' is thus

$$H' = H(\epsilon) - H_0 = V_0 \cdot \begin{pmatrix} -\epsilon & 0 & 0 \\ 0 & 0 & \epsilon \\ 0 & \epsilon & 0 \end{pmatrix} \quad (3)$$

Eigenvalues of the whole system are given by the equation in ϵ and λ , the eigenvalues themselves,

$$|(H(\epsilon) - \lambda \hat{1})| = 0, \text{ or,} \quad (4)$$

$$\begin{vmatrix} (V_0(1-\epsilon) - \lambda) & 0 & 0 \\ 0 & V_0 - \lambda & V_0\epsilon \\ 0 & V_0\epsilon & 2V_0 - \lambda \end{vmatrix} = 0 \quad (5)$$

$$((1-\epsilon)V_0 - \lambda)((V_0 - \lambda)(2V_0 - \lambda) - V_0^2\epsilon^2) = 0 \quad (6)$$

This gives $\lambda = (1-\epsilon)V_0$, $\frac{V_0}{2}(3 - \sqrt{4\epsilon^2 + 1})$, $\frac{V_0}{2}(3 + \sqrt{4\epsilon^2 + 1})$ as the eigenvalues of H . Call these $\lambda_{1,2,3}$ respectively.

15 Evaluating to second order in ϵ , we get

$$(\lambda_1, \lambda_2, \lambda_3) = V_0(1 - \epsilon, 1 - \epsilon^2, 2 + \epsilon^2) \quad (7)$$

16 **(b)** The last eigenvalue, $2V_0$, is non degenerate, so we attempt to approximate its perturbed analogue
 17 using non-degenerate perturbation theory as required, using both first and second order theories.
 18 The first order correction is given simply by

$$\begin{aligned} \langle v_3 | H' | v_3 \rangle &= (0 \quad 0 \quad 1) \begin{pmatrix} -\epsilon & 0 & 0 \\ 0 & 0 & \epsilon \\ 0 & \epsilon & 0 \end{pmatrix} \begin{pmatrix} 0 \\ 0 \\ 1 \end{pmatrix} \\ &= (0 \quad 0 \quad 1) \begin{pmatrix} 0 \\ \epsilon \\ 0 \end{pmatrix} \\ &= 0. \end{aligned} \quad (8)$$

19 That is, there is no first-order correction to the eigenvalue, which is expected, given the calculation
 20 from part a.

21 Moving on, representing the n^{th} order correction to the i^{th} eigenket as $|v_{i,n}\rangle$, the second-order
 22 correction is

$$\langle v_{3,0} | H' | v_{3,1} \rangle = \langle v_{3,0} | H' \sum_{m \neq 3} \frac{\langle v_{m,0} | H' | v_{3,1} \rangle}{E_3^{(0)} - E_m^{(0)}} | v_{m,0} \rangle \quad (9)$$

23 Since our eigenvectors are basis vectors, this basically filters components. Since we know $H' | v_{3,1} \rangle$,
 24 the calculation trivially reduces to

$$\begin{aligned} \langle v_{3,0} | H' | v_{3,1} \rangle &= (0 \quad 0 \quad 1) \begin{pmatrix} -\epsilon & 0 & 0 \\ 0 & 0 & \epsilon \\ 0 & \epsilon & 0 \end{pmatrix} \begin{pmatrix} 0 \\ \epsilon \\ 0 \end{pmatrix} \\ &= (0 \quad 0 \quad 1) \begin{pmatrix} 0 \\ 0 \\ \epsilon^2 \end{pmatrix} \\ &= \epsilon^2 \end{aligned} \quad (10)$$

25 Thus the second order correction is ϵ^2 , exactly matching the result from part a.

26 **(c)** As for the degenerate levels, corresponding to $|v_{1,0}\rangle$ and $|v_{2,0}\rangle$, we apply the degenerate theory.
 27 That is, compute the matrix W , with elements $\langle v_{i,0} | H' | v_{j,1} \rangle$ for $i, j \in \{1, 2\}$.

28 We get

$$W = \begin{pmatrix} -\epsilon & 0 \\ 0 & 0 \end{pmatrix}. \quad (11)$$

Since this matrix is diagonal, we simply read off the eigenvalues, ϵ and 0. These are the first order corrections to the energy values. Both of them match the exact results we got in part a. The first one matching exactly, and the second one missing a second order correction.

Question 3.

[Parth: Doing question 3, might have to check later]

The 3D isotropic oscillator has a potential profile that looks like the follows -

$$V(x, y, z) = \frac{1}{2}m\omega^2(x^2 + y^2 + z^2)$$

We know that for the 1D harmonic oscillator ($V(x) = 1/2m\omega^2x^2$), the eigenfunctions for the same take the following form

$$\psi_n(x) = \frac{1}{\sqrt{2^n n!}} \left(\frac{m\omega}{\pi \hbar} \right)^{1/4} \cdot e^{-\frac{m\omega x^2}{2\hbar}} \cdot H_n \left(\sqrt{\frac{m\omega}{\hbar}} x \right)$$

where $H_n(x)$ are the Hermite polynomials.

$$H_n(x) = (-1)^n e^{x^2} \frac{d^n}{dx^n} (e^{-x^2})$$

$$\text{with } H_0(x) = 1, H_1(x) = 2x$$

The potential profile in the 3D Harmonic oscillator is a separable function in x, y and z . Which means that the eigenfunctions to the Hamiltonian for the 3D harmonic oscillator can be derived as a direct product of the eigenfunctions of the 1D harmonic oscillators in x, y and z directions.

$$\psi_{n_x n_y n_z} = \frac{1}{\sqrt{2^{n_x+n_y+n_z} n_x! n_y! n_z!}} \left(\frac{m\omega}{\pi \hbar} \right)^{3/4} \cdot e^{-\frac{m\omega(x^2+y^2+z^2)}{2\hbar}} \cdot H_{n_x} \left(\sqrt{\frac{m\omega}{\hbar}} x \right) \cdot H_{n_y} \left(\sqrt{\frac{m\omega}{\hbar}} y \right) \cdot H_{n_z} \left(\sqrt{\frac{m\omega}{\hbar}} z \right)$$

with the energy of the eigenstate being

$$E_{n_x n_y n_z} = \left(n_x + n_y + n_z + \frac{3}{2} \right) \hbar \omega$$

Thus, the ground state for the 3D Harmonic oscillator is -

$$\psi_{000} = \left(\frac{m\omega}{\pi\hbar} \right)^{3/4} e^{-\frac{m\omega(x^2+y^2+z^2)}{2\hbar}}$$

43 with energy $3/2\hbar\omega$

44 The first excited state is triply degenerate, since any of n_x , n_y and n_z can be changed to 1 and thus we
45 have three degenerate eigenstates

$$\begin{aligned}\psi_{100} &= \frac{2}{\pi^{3/4}} \left(\frac{m\omega}{\hbar} \right)^{5/4} x e^{-\frac{m\omega(x^2+y^2+z^2)}{2\hbar}} \\ \psi_{010} &= \frac{2}{\pi^{3/4}} \left(\frac{m\omega}{\hbar} \right)^{5/4} y e^{-\frac{m\omega(x^2+y^2+z^2)}{2\hbar}} \\ \psi_{001} &= \frac{2}{\pi^{3/4}} \left(\frac{m\omega}{\hbar} \right)^{5/4} z e^{-\frac{m\omega(x^2+y^2+z^2)}{2\hbar}}\end{aligned}$$

46 with energies $5/2\hbar\omega$

47 We now introduce a perturbation to this system, given as

$$H' = \lambda x^2 y z$$

48 for a constant λ . Since the ground state is non-degenerate, we can use the non-degenerate perturbation
49 theory to calculate the correction to the energy eigenvalue. So

$$\begin{aligned}E_{000,1} &= \langle \psi_{000} | H' | \psi_{000} \rangle \\ &= \lambda \int \int \int |\psi_{000}(x, y, z)|^2 x^2 y z dx dy dz\end{aligned}$$

50 The integrals over y and z are zero, because ψ_{000} is an even function over x , y and z , and the perturbation
51 H' is odd in y and z . Which makes the integrals over y and z odd, and thus they equate to 0. Thus

$$E_{000,1} = 0$$

52 and thus, there is no correction to the ground state energy. $E'_{000} = E_{000} = \frac{3}{2}\hbar\omega$

53 For the first excited state, we use the degenerate perturbation theory. We thus need to calculate the matrix
54 elements $W_{ab} = \langle n_x n_y n_z | H' | m_x m_y m_z \rangle$. This means

$$W_{ab} = \lambda \langle n_x | x^2 | m_x \rangle \langle n_y | y | m_y \rangle \langle n_z | z | m_z \rangle$$

55 We just consider the matrix elements where one of n_x , n_y , n_z is 1 and the others are 0, and likewise for
56 the m_x , m_y , m_z

57 Now we know that

$$\begin{aligned}\langle n_x | x | n_x \rangle &= \int \psi_{n_x}(x) x \psi_{n_x}(x) dx \\ &= \int |\psi_{n_x}|^2 x dx = 0\end{aligned}$$

58 since the integrand is an odd function.

59 For $n_x \neq m_x$, we know that

$$\begin{aligned}\langle n_x | x | m_x \rangle &= \int \psi_{n_x}(x) x \psi_{m_x}(x) dx \\ &= \sum_{m_x \neq n_x} \frac{|\langle n_x | x | m_x \rangle|^2}{(m_x - n_x) \hbar \omega} \\ &= \sqrt{\frac{\hbar}{2m\omega}} \sum_{m_x \neq n_x} \frac{[\sqrt{m_x + 1} \delta_{n_x, m_x+1} + \sqrt{m_x} \delta_{n_x, m_x-1}]^2}{(m_x - n_x)} \\ &= \sqrt{\frac{\hbar}{2m\omega}} (\sqrt{m_x + 1} \delta_{n_x, m_x+1} + \sqrt{m_x} \delta_{n_x, m_x-1})\end{aligned}$$

60 $\langle n_x | x^2 | m_x \rangle = 0$, if $n_x \neq m_x$, since the integrand becomes an odd function, with an x^3 inside. Thus, for
61 the overall W_{ab} , the only non-zero elements will have $n_x = m_x$, $n_y \neq m_y$ and $n_z \neq m_z$. The elements that
62 satisfy this are

$$\begin{aligned}W_{010,001} &= \lambda \langle 0 | x^2 | 0 \rangle \langle 1 | y | 0 \rangle \langle 0 | z | 1 \rangle \\ W_{001,010} &= \lambda \langle 0 | x^2 | 0 \rangle \langle 0 | y | 1 \rangle \langle 1 | z | 0 \rangle\end{aligned}$$

63 For $\langle 0 | x^2 | 0 \rangle$, we have

$$\begin{aligned}\langle 0 | x^2 | 0 \rangle &= \int \left(\frac{m\omega}{\pi \hbar} \right)^{1/2} \cdot e^{-\frac{m\omega x^2}{\hbar}} x^2 dx \\ &= \left(\frac{m\omega}{\pi \hbar} \right)^{1/2} \int_{-\infty}^{\infty} x^2 e^{-\frac{m\omega x^2}{\hbar}} dx\end{aligned}$$

64 let $\left(\sqrt{\frac{m\omega}{\hbar}} x \right) = y$

$$\begin{aligned}
\left(\frac{m\omega}{\pi\hbar}\right)^{1/2} \int_{-\infty}^{\infty} x^2 e^{-\frac{m\omega x^2}{\hbar}} dx &= \left(\frac{m\omega}{\pi\hbar}\right)^{1/2} * \left(\frac{\hbar}{m\omega}\right) * \left(\sqrt{\frac{\hbar}{m\omega}}\right) \int_{-\infty}^{\infty} y^2 e^{-y^2} dy \\
&= \frac{1}{\sqrt{\pi}} * \left(\frac{\hbar}{m\omega}\right) * \frac{\sqrt{\pi}}{2} \\
&= \frac{\hbar}{2m\omega}
\end{aligned}$$

65 Thus, the W_{ab} matrix becomes -

$$W_{ab} = \lambda \begin{bmatrix} 0 & 0 & 0 \\ 0 & 0 & \left(\frac{\hbar}{2m\omega}\right)^2 \\ 0 & \left(\frac{\hbar}{2m\omega}\right)^2 & 0 \end{bmatrix}$$

66 Thus, the energy corrections, are the eigenvalues of W_{ab} , which are computed as

$$\begin{aligned}
-E \left(E^2 - \left(\frac{\hbar}{2m\omega} \right)^4 \right) &= 0 \\
\Rightarrow E_1 &= 0, \pm \lambda \left(\frac{\hbar}{2m\omega} \right)^2
\end{aligned}$$

67 Thus, the perturbation splits the degeneracy. It leaves $E'_{100} = E_{100} = \frac{5}{2}\hbar\omega$ while the degeneracy in ψ_{010}
68 and ψ_{001} is lifted with their energy eigenvalues changing to $E = \frac{5}{2}\hbar\omega \pm \lambda \left(\frac{\hbar}{2m\omega} \right)^2$.

69 This is the effect of the perturbation on the ground state and the first excited state of the isotropic
70 oscillator.

71 Question 4.

72 We use the following expression for the Kinetic energy T:

$$\begin{aligned}
T &= E - mc^2 \\
&= \sqrt{p^2 c^2 + m^2 c^4} - mc^2 \\
&= (mc^2) \sqrt{1 + \frac{p^2}{m^2 c^2}} - mc^2 \\
&= \frac{p^2}{2m} - \frac{p^4}{8m^3 c^2} + \mathcal{O}(p^6)
\end{aligned} \tag{12}$$

73 Using the first two terms to find the lowest order relativistic correction, we get the following Hamiltonian
74 for the 1-D harmonic oscillator:

$$H = \frac{p^2}{2m} - \frac{p^4}{8m^3c^2} + \frac{1}{2}m\omega^2x^2 \quad (13)$$

75 We treat $H_0 = \frac{p^2}{2m} + \frac{1}{2}m\omega^2x^2$ as our unperturbed Hamiltonian and $H_1 = -\frac{p^4}{8m^3c^2}$ as the perturbation.

76 The resulting energy levels will be of the form $E_n = E_n^0 + E_n^1$ where E_n^1 is the relativistic correction.

77 This requires calculation of terms of the form:

$$\begin{aligned} E_n^1 &= \langle n^0 | H_1 | n^0 \rangle \\ &= \langle n^0 | \left(\frac{-\hbar^2\omega^2}{32mc^2} \right) (a^\dagger - a)^4 | n^0 \rangle \end{aligned} \quad (14)$$

78 Here we have used the ladder operators $a = \sqrt{\frac{m\omega}{2\hbar}} \left(x + \frac{ip}{m\omega} \right)$ and $a^\dagger = \sqrt{\frac{m\omega}{2\hbar}} \left(x - \frac{ip}{m\omega} \right)$

79 Since states of different n are orthogonal to each other, we only select terms that have equal creation and
80 annihilation operators from $(a^\dagger - a)^4$. There are ${}^4C_2 = 6$ terms of this kind:

$$\begin{aligned} &a^\dagger a^\dagger a a \\ &a^\dagger a a^\dagger a \\ &a a^\dagger a^\dagger a \\ &a^\dagger a a a^\dagger \\ &a a^\dagger a a^\dagger \\ &a a a^\dagger a^\dagger \end{aligned} \quad (15)$$

81 Inserting these terms into the equation for E_n^1 and using $a^\dagger |n\rangle = \sqrt{n+1} |n+1\rangle$ and $a |n\rangle = \sqrt{n} |n-1\rangle$
82 we get:

$$\begin{aligned} E_n^1 &= \left(\frac{-\hbar^2\omega^2}{32mc^2} \right) \left(n\sqrt{n^2-1} + n^2 + n(n+1) + (n+1)n + (n+1)^2 + (n+2)(n+1) \right) \\ &= \left(\frac{-\hbar^2\omega^2}{32mc^2} \right) \left(n\sqrt{n^2-1} + 5n^2 + 7n + 3 \right) \end{aligned} \quad (16)$$

83 Thus energies with first order relativistic corrections for the Harmonic oscillator are:

$$E_n = \hbar\omega \left(\frac{1}{2} + n \right) + \left(\frac{-\hbar^2\omega^2}{32mc^2} \right) \left(n\sqrt{n^2-1} + 5n^2 + 7n + 3 \right) \quad (17)$$

Question 6.

$$V(r) = \begin{cases} -\frac{Ze^2}{r}, & r \geq a \\ -\frac{Ze^2}{a} \left(\frac{3}{2} - \frac{r^2}{2a^2} \right), & r \leq a \end{cases}$$

84 where Z is the nuclear charge and a is the nuclear radius. Assuming that $a \ll \frac{\hbar^2}{me^2} = a_0$ (typically
 85 , $a \sim 10^{-13}\text{cm}$, while $a_0 \sim 10^{-8}\text{cm}$ calculate the first order change in the ground state energy from its
 86 value for a point nucleus.

87 Ans. The Hamiltonian for such a system will be given by:

$$\begin{aligned} H &= \frac{p^2}{2m} + V(r) \\ &= \frac{p^2}{2m} - \frac{Ze^2}{r} + \left(V(r) + \frac{Ze^2}{r} \right) \end{aligned} \quad (18)$$

88 We regard $H_0 = \frac{p^2}{2m} - \frac{Ze^2}{r}$ as the unperturbed Hamiltonian and $H_1 = V(r) + \frac{Ze^2}{r}$ as the perturbation. The
 89 perturbation term may be explicitly written as follows to verify that it is small (it is only non zero and
 90 finite over a small region, hence any integrals involving it will be small).

$$\begin{aligned} H_1(r) &= 0 \text{ for } r > a \\ &= -\frac{Ze^2}{a} \left(\frac{3}{2} - \frac{a}{r} - \frac{r^2}{2a^2} \right) \text{ for } r \leq a \end{aligned} \quad (19)$$

91 We can now write the perturbation to the ground state as:

$$E_0^{(1)} = \langle \psi_0 | H_1 | \psi_0 \rangle \quad (20)$$

92 Where ψ_0 is the ground state wavefunction for the hydrogen atom. Substituting the position basis
 93 expressions for relevant quantities, we get the following integral:

$$E_0^{(1)} = \frac{-Ze^2 \times 4\pi}{\pi a_0^3 a} \int_0^a \left(\frac{3}{2} - \frac{a}{r} - \frac{r^2}{2a^2} \right) e^{-2r/a_0} r^2 dr \quad (21)$$

94 Since $a \ll a_0$ we get

$$E_0^{(1)} = \frac{-4Ze^2}{a_0^3 a} \int_0^a \left(\frac{3r^2}{2} - ar - \frac{r^4}{2a^2} \right) (1 + \mathcal{O}(a/a_0)) dr \quad (22)$$

95 Ignoring the $\mathcal{O}(a/a_0)$ term,

$$\begin{aligned}
E_0^{(1)} &= \frac{-4Ze^2}{a_0^3 a} \int_0^a \left(\frac{3r^2}{2} - ar - \frac{r^4}{2a^2} \right) dr \\
&= \frac{2Ze^2 a^2}{5a_0^3}
\end{aligned} \tag{23}$$

96 **Question 8.**

97 [Parth: Doing question 8, might have to check later] [Sankalp: No, me BRUH]

98 The original Hamiltonian is of the form $H_0 = AL^2 + BL_z$. There's a small perturbation added to this
 99 Hamiltonian of the form $H_1 = CL_y$, where C is a constant. Thus, we have

$$\begin{aligned}
H_0 &= AL^2 + BL_z \\
L_y &= \left(\frac{L_+ - L_-}{2i} \right)
\end{aligned}$$

100 The eigenkets of H_0 are the simultaneous eigenkets of L^2 and L_z , $|l, m\rangle$. The energy eigenvalues here
 101 are $E_{l,m} = A\hbar^2 l(l+1) + B\hbar m$.

102 $L_y = \left(\frac{L_+ - L_-}{2i} \right)$, on the other hand, is off-diagonal in the $|l, m\rangle$ basis (owing to the ladder operators seen).

103 Due to this, the first order energy shift $\langle l, m | CL_y | l, m \rangle$ vanishes, and we are left with the second order
 104 energy shifts

$$\begin{aligned}
E'_{l,m} &= E_{l,m}^{(0)} + C^2 \sum_{m' \neq m} \frac{|\langle l, m' | \left(\frac{L_+ - L_-}{2i} \right) | l, m \rangle|^2}{E_{l,m}^{(0)} - E_{l,m'}^{(0)}} + O(C^3) \\
&= E_{l,m}^{(0)} + (C\hbar)^2 \frac{(l-m)(l+m+1)}{-4B\hbar} + (C\hbar)^2 \frac{(l+m)(l-m+1)}{4B\hbar} + O(C^3) \\
&= E_{l,m}^{(0)} + \hbar m \frac{C^2}{2B} + O(C^3)
\end{aligned}$$

Similarly, the shifts in the eigenfunctions are

$$\begin{aligned}
|l, m\rangle &= |l, m\rangle_0 + C \sum_{m' \neq m} \frac{\langle l, m' | \left(\frac{L_+ - L_-}{2i} \right) | l, m \rangle}{E_{l,m}^{(0)} - E_{l,m'}^{(0)}} |l, m'\rangle \\
&= |l, m\rangle_0 + \frac{C}{2i} \sum_{m' \neq m} \frac{\langle l, m' | (L_+ - L_-) | l, m \rangle}{E_{l,m}^{(0)} - E_{l,m'}^{(0)}} |l, m'\rangle \\
&= |l, m\rangle_0 + \frac{C\hbar}{2i} \sum_{m' \neq m} \frac{\sqrt{l(l+1) - m(m+1)} \langle l, m' | l, m+1 \rangle - \sqrt{l(l+1) - m(m-1)} \langle l, m' | l, m-1 \rangle}{E_{l,m}^{(0)} - E_{l,m'}^{(0)}} |l, m'\rangle
\end{aligned}$$

105 We make these calculations for the case $l = 1$. Thus, the corrections in each state corresponding to
 106 $m = -1, 0, +1$ are -

$$\begin{aligned} E'_{1,-1} &= E_{1,-1} - \frac{C^2 \hbar}{2B} = 2A\hbar^2 - \frac{(C^2 + 2B^2)\hbar}{2B} \\ E'_{1,0} &= E_{1,0} = 2A\hbar^2 \\ E'_{1,1} &= E_{1,1} + \frac{C^2 \hbar}{2B} = 2A\hbar^2 + \frac{(C^2 + 2B^2)\hbar}{2B} \end{aligned}$$

$$\begin{aligned} |1, -1\rangle' &= |1, -1\rangle - \frac{C}{\sqrt{2}Bi} |1, 0\rangle \\ |1, 0\rangle' &= |1, 0\rangle - \frac{C}{\sqrt{2}Bi} (|1, 1\rangle + |1, -1\rangle) \\ |1, 1\rangle' &= |1, 1\rangle - \frac{C}{\sqrt{2}Bi} |1, 0\rangle \end{aligned}$$

107 Solving the system exactly, we recognize that the new states must be within the linear span of the $|l, m\rangle$
 108 states. We then write the effect of the Hamiltonian on an arbitrary such state to obtain the eigenvalue
 109 condition.

$$H \sum_{l,m} a_{l,m} |l, m\rangle = (A \hat{L}^2 + B \hat{L}_z + C \hat{L}_y) \sum_{l,m} a_{l,m} |l, m\rangle \quad (24)$$

110 Define $c_{+,l,m}$ and $c_{-,l,m}$ as $\hat{L}_+ |l, m\rangle = \hbar c_{+,l,m} |l, m+1\rangle$ and similarly for the negative. We write $\hat{L}_y = \frac{\hat{L}_+ - \hat{L}_-}{2i}$
 111 and absorb the $2i$ into C as $D = \frac{C}{2i}$ to get

$$\begin{aligned} H \sum_{l,m} |l, m\rangle &= (A \hat{L}^2 + B \hat{L}_z + D \hat{L}_+ - D \hat{L}_-) \sum_{l,m} a_{l,m} |l, m\rangle \\ &= \hbar \sum_{l,m} a_{l,m} ((\hbar A l(l+1) + Bm) |l, m\rangle + D c_{+,l,m} |l, m+1\rangle - D c_{-,l,m} |l, m-1\rangle) \\ &= \hbar \sum_{l,m} (a_{l,m} (\hbar A l(l+1) + Bm) + D a_{l,m-1} c_{+,l,m-1} - D a_{l,m+1} c_{-,l,m+1}) |l, m\rangle \end{aligned} \quad (25)$$

(26)

112 This gives us a form for the matrix representation of H , which has diagonal terms, and terms one above
 113 and one below the diagonal. The eigenvalues of this matrix can be obtained (painfully) via brute force
 114 calculation.

115 We use $l = 1$ for this case to construct the matrix representation of H

$$H = \hbar \begin{pmatrix} 2\hbar A - B & -Dc_{-,1,0} & 0 \\ Dc_{+,1,-1} & 2\hbar A & -Dc_{-,1,1} \\ 0 & Dc_{+,1,0} & 2\hbar A + B \end{pmatrix}. \quad (27)$$

116 All the $c_{\pm,l,m}$ equal $\sqrt{2}$ when appearing in the matrix (zero elements are outside 3×3). And replacing
 117 back $D = \frac{C}{2i}$ we get

$$H = \hbar \begin{pmatrix} 2\hbar A - B & -\frac{C}{\sqrt{2}i} & 0 \\ \frac{C}{\sqrt{2}i} & 2\hbar A & -\frac{C}{\sqrt{2}i} \\ 0 & \frac{C}{\sqrt{2}i} & 2\hbar A + B \end{pmatrix}. \quad (28)$$

118 Eigenvalues of this matrix, using a symbolic solver, were found to be

$$\lambda_1 = 2A\hbar^2, \quad (29)$$

$$\lambda_2 = 2A\hbar^2 - \hbar\sqrt{B^2 + C^2}, \quad (30)$$

$$\lambda_3 = 2A\hbar^2 + \hbar\sqrt{B^2 + C^2}. \quad (31)$$

119 At first sight, these seem very different from the previously calculated values, but a trivial factorisation
 120 of the square root and binomial approximation to first order gives us exactly the values calculated via
 121 perturbation theory!