

Worksheet #12

(Wednesday, February 4, 2026)

Name

Questions (5 pts):

Consider a hydrogen atom in an electric field \mathcal{E} . As we showed, the non-zero first-energy corrections to the unperturbed energies for the $n = 2$ state can be obtained by diagonalizing the following perturbation matrix (in the $|200\rangle, |210\rangle$ basis):

$$\begin{pmatrix} 0 & -3e\mathcal{E}a_0 \\ -3e\mathcal{E}a_0 & 0 \end{pmatrix}$$

- (a) Find the eigenvalues and eigenvectors of the perturbation matrix above. Write the eigenvectors in terms of the basis states.

- (b) Sketch the new energy levels and indicate the new states resulting from the perturbation. Don't forget to include ones for which the energy correction is 0.

(c) **If you have time:** use symmetry arguments to find the first-order energy correction to the ground state (i.e. $n=1$).

(d) **If you have time:** Let's say you are a spectroscopist measuring emission from the $n=2$ to $n=1$ transition. Sketch the transitions and the emission spectra for what you expect to see without and with the applied electric field.