The $n = 2$ level of the hydrogen atom is 4-fold degenerate with energy -0.125 $E_h$. In terms of the $|nlm\rangle$ quantum numbers these states are $|2,0,0\rangle$, $|2,1,0\rangle$, $|2,1,1\rangle$, and $|2,2,-1\rangle$. An electric field in the $z$-direction splits the degeneracy because it mixes the $|2s\rangle$ and the $|2p_z\rangle$ orbitals creating one linear combination polarized in the direction of the field and another polarized against the field.

What happens is that the $|s\rangle$ and $|p\rangle$ wavefunctions "mixed" to produce eigenstates that have shifted centers. This means the atom gets an induced electric dipole moment, whose interaction with the external field either lowers or raises the eigenenergy.

Note in particular that the electronic center of charge has moved from the origin, which means the states have nonzero dipole moments. With the electric field pointing downwards, the state to the left has a lower energy and the one to the right is raised.

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**Degenerate Perturbation Theory**

The Hamiltonian for this perturbation in atomic units is:

\[ H' = \varepsilon_z. \]
which in spherical polar coordinates is:

\[
H' = \varepsilon r \cos(\theta),
\]

where \(\varepsilon\) is the electric field strength.

In this perturbation method treatment the hydrogen atom eigenfunctions are used to evaluate the matrix elements associated with the total Hamiltonian,

\[
[H = H^0 + H'^0]
\]

Since the results for \(H^0\) are known (-0.125 E\(\text{h}\)) only the matrix elements for \(H'\) need to be evaluated and most of these are zero. Below we show that \(\varepsilon = -3\varepsilon\) and that the other matrix elements involving the \(n = 2\) orbitals are equal to zero.

\[
\langle 2s | H' | 2s \rangle = 0
\]

\[
\langle 2p_z | H' | 2p_z \rangle = \langle 2p_y | H' | 2p_y \rangle = \langle 2p_x | H' | 2p_x \rangle = 0
\]

\[
\langle 2s | H' | 2p_z \rangle = -3\varepsilon
\]

\[
\langle 2s | H' | 2p_x \rangle = \langle 2s | H' | 2p_y \rangle = 0
\]
\[ \int_{0}^{\infty} \int_{0}^{\pi} \int_{0}^{2\pi} \psi_{2s}(r) \varepsilon r \cos(\theta) \psi_{2py}(r, \theta, \phi) r^2 \sin(\theta) \, d\phi \, d\theta \, dr \rightarrow 0 \]

\[ \langle 2p_x | H^\prime | 2p_x \rangle = \langle 2p_x | H^\prime | 2p_y \rangle = \langle 2p_y | H^\prime | 2p_z \rangle = 0 \]

\[ \int_{0}^{\infty} \int_{0}^{\pi} \int_{0}^{2\pi} \psi_{2px}(r, \theta, \phi) \varepsilon r \cos(\theta) \psi_{2py}(r, \theta, \phi) r^2 \sin(\theta) \, d\phi \, d\theta \, dr \rightarrow 0 \]

\[ \int_{0}^{\infty} \int_{0}^{\pi} \int_{0}^{2\pi} \psi_{2px}(r, \theta, \phi) \varepsilon r \cos(\theta) \psi_{2pz}(r, \theta) r^2 \sin(\theta) \, d\phi \, d\theta \, dr \rightarrow 0 \]

\[ \int_{0}^{\infty} \int_{0}^{\pi} \int_{0}^{2\pi} \psi_{2py}(r, \theta, \phi) \varepsilon r \cos(\theta) \psi_{2pz}(r, \theta) r^2 \sin(\theta) \, d\phi \, d\theta \, dr \rightarrow 0 \]

The matrix elements of the 4x4 perturbation matrix are

\[ \langle \psi_i | H^o + H^\prime | \psi_j \rangle, \]

where the \( \psi \)'s are the 2s, 2p_z, 2p_x, and 2p_y hydrogen atomic orbitals. Using the values of the integrals evaluated above, the perturbation matrix is formed and its eigenvalues and eigenvectors found.

\[
\begin{pmatrix}
-0.125 - E & -3 \varepsilon & 0 & 0 \\
-3 \varepsilon & -0.125 - E & 0 & 0 \\
0 & 0 & -0.125 - E & 0 \\
0 & 0 & 0 & -0.125 - E
\end{pmatrix}
\begin{pmatrix}
c_1 \\
c_2 \\
c_3 \\
c_4
\end{pmatrix} = 0
\]

This 4x4 energy matrix is clearly one 2x2 and two 1x1 energy matrices. In other words, as we learned from evaluating the matrix elements, the 2p_x and 2p_y are not perturbed by the electric field to first order and have energy -0.125 E_{\text{H}}.

The eigenvectors and eigenvalues of the 2x2 are found as follows.

\[
\begin{pmatrix}
-0.125 - E & -3 \varepsilon & 0 & 0 \\
-3 \varepsilon & -0.125 - E & 0 & 0 \\
0 & 0 & -0.125 - E & 0 \\
0 & 0 & 0 & -0.125 - E
\end{pmatrix}
\begin{pmatrix}
c_1 \\
c_2 \\
c_3 \\
c_4
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\end{pmatrix} = 0
\]

\[
\begin{pmatrix}
-0.125 - E & -3 \varepsilon & 0 & 0 \\
-3 \varepsilon & -0.125 - E & 0 & 0 \\
0 & 0 & -0.125 - E & 0 \\
0 & 0 & 0 & -0.125 - E
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\begin{pmatrix}
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This 4x4 energy matrix is clearly one 2x2 and two 1x1 energy matrices. In other words, as we learned from evaluating the matrix elements, the 2p_x and 2p_y are not perturbed by the electric field to first order and have energy -0.125 E_{\text{H}}.

The eigenvectors and eigenvalues of the 2x2 are found as follows.
The wavefunctions of the perturbed 2s and 2p\(_z\) orbitals are sp\(_z\) hybrid states as shown below.

\[
\begin{pmatrix}
\frac{1}{\sqrt{2}} (2s + 2p_{z}) & E = (-0.125 - 3 \varepsilon) E_{h} \\
\frac{1}{\sqrt{2}} (2s - 2p_{z}) & E = (-0.125 + 3 \varepsilon) E_{h}
\end{pmatrix}
\]

Because the energy of the symmetric 1s state is unaffected by the electric field, the effect of this perturbation on the electronic spectrum of hydrogen is to split the \(n = 1\) to \(n = 2\) transition into three lines of relative intensity 1:2:1.

\[
< 2s | H' | 2s > = 0
\]

\[
\psi_{1s} (r) = \frac{1}{\sqrt{\pi}} \exp(-r)
\]

\[
\int_{0}^{\infty} \int_{0}^{\pi} \int_{0}^{2\pi} \psi_{1s} (r) \varepsilon r \cos (\theta) \psi_{1s} (r) r^2 \sin (\theta) d\phi d\theta dr \rightarrow 0
\]

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**Contributors and Attributions**

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- Emilio Pisanty via StackExchange (Physics)