

Q: For $\ell=0$, how does $\langle 1/r_{nl} \rangle$ scale with n

- A. n^{-1}
- B. n^{-2}
- C. n^{-3}

Q: For $\ell=0$, how does $|\psi(0)|^2$ scale with n

- A. n^{-1}
- B. n^{-2}
- C. n^{-3}
- D. n^{-6}

Q: In helium, for the same electronic configuration, are triplet or singlet states more tightly bound

A: Singlet

B: Triplet

Q: The difference between triplet and singlet binding energy (exchange energy) has its origin in

A: Magnetic energy

B: Spin-spin interactions

C: Electrostatic interactions

Q: Which interaction reflects that the potential between nucleus and electron is not exactly an $1/r$ potential?

- A. The Darwin term
- B. The Lamb shift
- C. The volume isotope effect
- D. A, B, C
- E. B, C

Q: The fine structure adds a term ()LS to the Hamiltonian and affects only states with $L \neq 0$.

- A. True
- B. False

Q: For $L \neq 0$, the orbiting electron creates a magnetic field. Spin-orbit interaction is the energy of the electron's spin in this magnetic field.

- A. True
- B. False

Q: In Dirac theory, which states are degenerate?

- A. $2 P_{1/2}$ and $2 P_{3/2}$
- B. $2 S_{1/2}$ and $2 P_{3/2}$
- C. $2 S_{1/2}$ and $2 P_{1/2}$

Q: The degeneracy is lifted by

- A. The Lamb shift
- B. The size of the proton
- C. The mass of the proton
- D. A and B
- E. A, B, C

Hyperfine structure of hydrogen

Q: At high magnetic fields, what are the magnetic moments of those HF states in units of μ_B ?

A: 1, 1, -1, -1

B: 1, 0, -1, 0

C: $\frac{1}{2}$, 0, 0, $\frac{1}{2}$

D: 0, 0, 0, 0

Q: At low magnetic fields, what are the magnetic moments of those HF states?

A: 1, 1, -1, -1

B: 1, 0, -1, 0

C: $\frac{1}{2}$, 0, 0, $\frac{1}{2}$

D: 0, 0, 0, 0

Hyperfine structure of positronium

Q: How many hyperfine states does the 1s state have?

A: 1

B: 2

C: 3

D: 4

Q: At high magnetic fields, what are the magnetic moments of those HF states in units of μ_B ?

A: 1, 1, -1, -1

B: 1, 0, -1, 0

C: 2, 1, -1, -2

D: 2, 0, 0, -2

Q: At low magnetic fields, what are the magnetic moments of those HF states?

A: 0, 0, 0, 0

B: 1, 0, -1, 0

C: 2, 1, -1, -2

D: 2, 0, 0, -2