

Electron

The particle-in-a-box ground state has quantum number $n = 0$.

The H_2 molecule does not have a pure rotational absorption spectrum.

The particle-in-a-box transition that absorbs the photon of longest wavelength is from the $n = 1$ level to the $n = 2$ level.

If f is an eigenfunction of the linear operator \hat{A} with eigenvalue a , then f is an eigenfunction of \hat{A}^2 with eigenvalue a^2 .

The number of peaks in the graph of the harmonic oscillator $|\psi_v|^2$ equals $v + 1$.

The product of a number and its complex conjugate is always a real number.

If we measure the property B when the system's state function is not an eigenfunction of \hat{B} , then we can get a result that is not an eigenvalue of \hat{B} .

The particle-in-a-box energy levels are equally spaced.

For a one-particle problem with $V = br^3$, where b is a positive constant, the stationary state wavefunctions have the form $\psi = f(r)Y_l^m(\theta, \phi)$.

$\int_{-\infty}^{\infty} \psi dx = 1$ for a one-particle, one-dimensional system.

The wavefunction ψ must be a real function.

For the $n = 1$ particle-in-a-box stationary state, the probability of finding the particle in the left third of the box equals the probability of finding it in the middle third.

$$\langle B \rangle = \int \hat{B}|\psi|^2 d\tau.$$

As the frequency of light increases, its wavelength decreases.

$5x$ is an eigenvalue of the position operator \hat{x} .

The spacing between adjacent vibrational levels of a diatomic molecule remains constant as the vibrational energy increases.

If $z = z^*$, then z must be a real number.

Since $\hat{L}^2 Y_l^m = l(l+1)\hbar^2 Y_l^m$, it follows that $\hat{L}^2 = l(l+1)\hbar^2$.

The ground-state energy of a particle in a box is zero.

If g is an eigenfunction of the linear operator \hat{B} , then cg is an eigenfunction of \hat{B} , where c is an arbitrary constant.

A probability density can never be negative.

The probability density for the particle-in-a-box stationary states is constant along the length of the box.

All harmonic oscillator wavefunctions with v an odd integer must have a node at the origin.

If f_1 and f_2 are eigenfunctions of \hat{B} , then $c_1 f_1 + c_2 f_2$ must be an eigenfunction of \hat{B} , where c_1 and c_2 are constants.

Every function of x , y , and z has the form $f(x)g(y)h(z)$.

The spacings between successive two-particle rigid rotor energy levels remain constant as J increases.

ψ must be an eigenfunction of each operator \hat{B} that represents a physical property of the system.

The $v = 1$ harmonic oscillator wavefunction must be negative for $x < 0$.

The symbol e stands for the charge on an electron.

Increasing particle-in-a-box stationary-state energy corresponds to increasing number of nodes in the wave function.

If \hat{A} and \hat{B} do not commute, it is impossible for an eigenfunction of \hat{A} to be also an eigenfunction of \hat{B} .

For harmonic oscillator wavefunctions, $\int_{-\infty}^{\infty} \psi_v^*(x)\psi_{v+1}(x) dx = 0$.

The \hat{L}^2 eigenvalues are degenerate except for $l = 0$.

If the harmonic oscillator wavefunction ψ_v is an even function, then ψ_{v+1} is also an even function.

The degeneracy of the $J = 4$ two-particle rigid-rotor energy level is 7.