

PARTICLE IN A FINITE BOX

$$V(x) = \begin{cases} V_0 & x \le 0 & (I) \\ V(x) = V_0 & 0 < x < L & (II) \\ V_0 & L \le x & (III) \end{cases}$$

WE DIVIDE THE PHYSICAL SPACE IN 3 REGIONS: I, II.

FROM THE TISE AND EXVO

WE GET
$$\frac{3\Psi}{3\chi^2} = + \frac{2M}{4\chi^2}(V_0 - E)\Psi$$

IN RESCOUL I AND W

$$\frac{dx^2}{dx^2} = \kappa^2 \Phi,$$

WHERE

$$K^{2} = \sqrt{\frac{2m}{\kappa^{2}}}(V_{0} - E) =$$

BUT IN REGION II

$$\frac{\partial X_5}{\partial_s \partial_s} = -k_s \partial_s$$

M14 H

$$R = \sqrt{\frac{2mE}{\kappa^2}}$$

$$\Psi_{\underline{\Pi}}(x) = C \mathcal{Q} + D \mathcal{Q}$$

$$\mathcal{F}(x) = \underline{F} \mathcal{C}^{kx} + F \hat{\mathcal{C}}^{kx}$$

BUT AT X->D AND X-D-D, WE SHOULD NOT FIND THE PARTICE, THUS

B = 0 = E.

NEXT, WE CONSIDER CONTIDUITY OF THE
FUNCTIONS & AND ITS FIRST DERIVATIVE

$$\vec{\mathcal{L}}(r) = \vec{\mathcal{L}}(r)$$

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$$\frac{\partial x}{\partial \hat{\tau}}\Big|_{x=0} = \frac{\partial x}{\partial \hat{T}^{H}}\Big|_{x=0}$$

Table 6-2. A		systems Studied in Ch		
Name of System	Physical Example	Potential and Total Energies	Probability Density	Significant Feature
Zero potential	Proton in beam from cyclotron	EV(x)	Ψ*Ψx	Results used for other systems
Step potential (energy below top)	Conduction electron near surface of metal		√ у *Ψ х	Penetration of excluded region
Step potential (energy above top)	Neutron trying to escape nucleus		Ψ*Ψ	Partial reflec- tion at potential discontinuity
Barrier potential (energy below top)	α particle trying to escape Coloumb barrier		$ \int_{0}^{\psi*\psi} x $	Tunneling
Barrier potential (energy above top)	Electron scat- tering from negatively ionized atom		$\int_{0}^{\sqrt{\psi*\psi}} \frac{1}{a} x$	No reflection at certain energies
Finite square well potential	Neutron bound in nucleus	$ \begin{array}{c c} V(x) \\ E \end{array} $	$\bigcup_{0}^{\psi * \psi} \int_{a}^{w} dx$	Energy quantization
Infinite square well potential	Molecule strictly confined to box	V(x) C		Approximation to finite square well
Simple harmonic oscillator potential	Atom of vibrating diatomic molecule	V(x) x'' E	<u></u>	Zero-point energy

QUESTIONS

- 1. Can there be solutions with E < 0 to the time-independent Schroedinger equation for the zero potential?
- 2. Why is it never possible in classical mechanics to have E < V(x)? Why is it possible in quantum mechanics, providing there is some region in which E > V(x)?
- 3. Explain why the general solution to a one-dimensional time-independent Schroedinger equation contains two different functions, while the general solution to the corresponding Schroedinger equation contains many different functions.
- 4. Consider a particle in a long beam of very accurately known momentum. Does a wave function in the form of a group provide a more or a less realistic description of the particle than a single complex exponential wavefunction like (6-9)?

- 5. Under what c mation to an
- 6. If a potential discontinuities
- 7. By combining ing wave. Wh.
- 8. Just what is a
- 9. How can it be associated par
- ior of an unb can be followed apparatus? W
- 11. Exactly what incident on a by the statement than the step
- 12. Since a real e oscillatory fur
- 13. What do you 6-8 as it reflect
- 14. What is the faturneling three
- 15. A particle is i and it is refle direction of ii changed into
- 16. In the sun, to the Coulomb than the sum is responsible on earth if it
- 17. Are there any which is of ir
- 18. Show from a always has or the eigenfunc
- 19. Why do finit the character
- 20. What would well look like
- 21. Why do the best approximately?
- 22. In the n = 3 positions bet positions?
- 23. Explain in si principle.
- 24. Would you matter at ver
- 25. If the eigenfu has even par