

# Molecular and Solid State Physics

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Calculate the macroscopic properties from the microscopic structure.

# Goals

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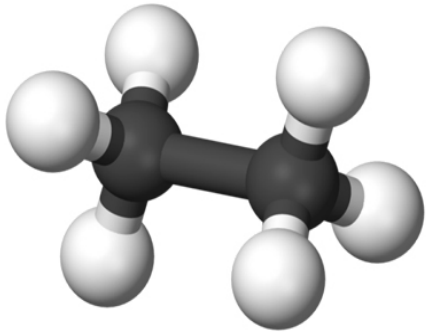
The microscopic structure determines the macroscopic properties.

At the end of this course you should be able to explain how any property of any molecule or solid can be calculated using quantum mechanics and statistical physics.

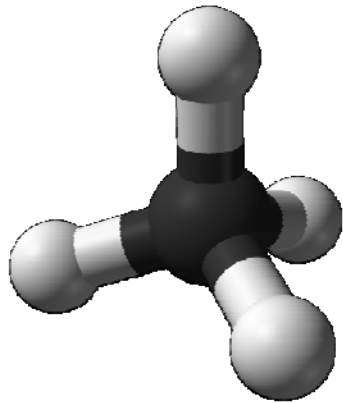
For example: knowing how the atoms are arranged in a crystal, you must be able to say if it is an electrical conductor or not.

# Molecules

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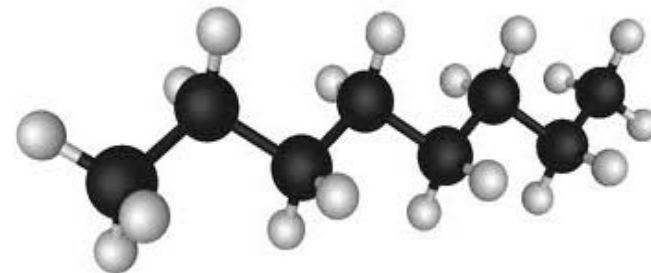
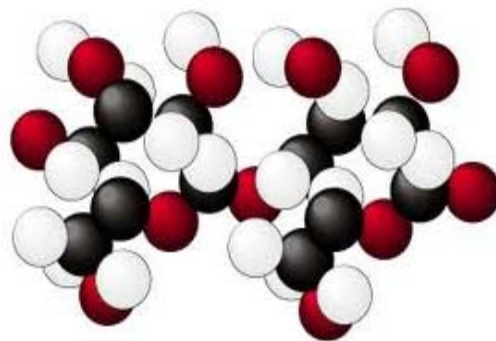


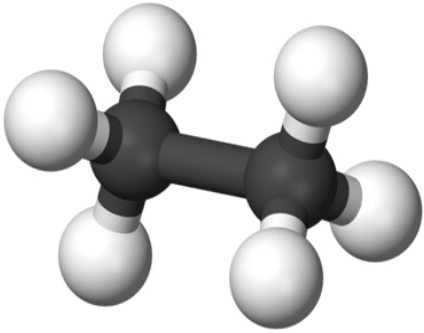
There are billions of useful molecules.



Acids, esthers, alkanes, ...

Biological molecules: DNA, RNA, proteins





# Molecules

Every property of a molecule can be calculated using multi-particle quantum mechanics.

$$H_{\text{mp}} = - \sum_i \frac{\hbar^2}{2m_e} \nabla_i^2 - \sum_a \frac{\hbar^2}{2m_a} \nabla_a^2 - \sum_{a,i} \frac{Z_a e^2}{4\pi\epsilon_0 |\vec{r}_i - \vec{r}_a|} + \sum_{i<j} \frac{e^2}{4\pi\epsilon_0 |\vec{r}_i - \vec{r}_j|} + \sum_{a<b} \frac{Z_a Z_b e^2}{4\pi\epsilon_0 |\vec{r}_a - \vec{r}_b|}$$



We will calculate:

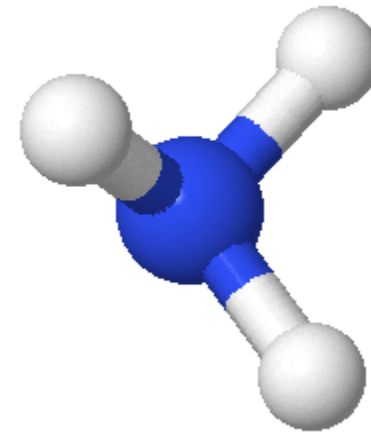
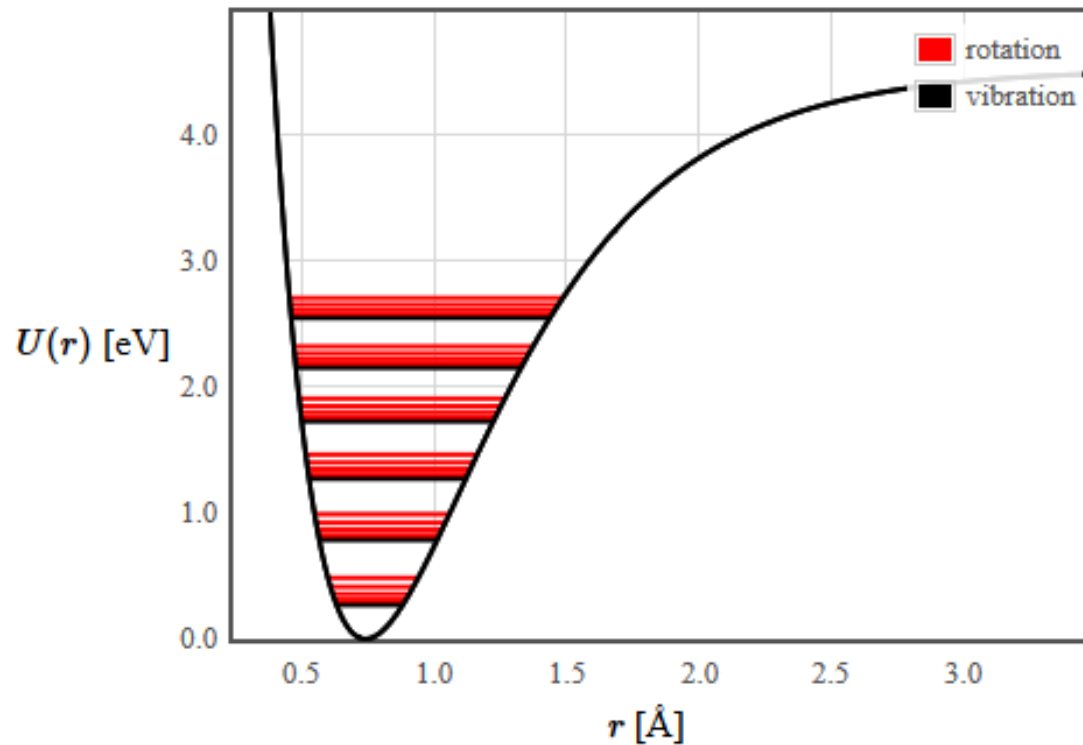
- bond length
- bond strength
- molecular energy levels



# Molecules

$$H_{\text{mp}} = - \sum_i \frac{\hbar^2}{2m_e} \nabla_i^2 - \sum_a \frac{\hbar^2}{2m_a} \nabla_a^2 - \sum_{a,i} \frac{Z_a e^2}{4\pi\epsilon_0 |\vec{r}_i - \vec{r}_a|} + \sum_{i<j} \frac{e^2}{4\pi\epsilon_0 |\vec{r}_i - \vec{r}_j|} + \sum_{a<b} \frac{Z_a Z_b e^2}{4\pi\epsilon_0 |\vec{r}_a - \vec{r}_b|}$$

$$E = \frac{\langle \Psi | H | \Psi \rangle}{\langle \Psi | \Psi \rangle}$$



Bond length: 0.74144 Å.  
 Dissociation energy: 4.52 eV.

# Solids

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Solids are large molecules

Crystal structures

Determining crystal structures with x-ray diffraction

Photons in solids

Phonons in solids (lattice vibrations)

Thermal properties

Free electron model

Band structure (metals, semiconductors, insulators)

Semiconductors

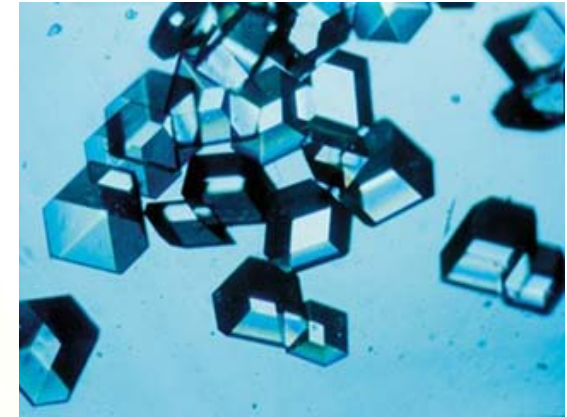
# Crystal = periodic arrangement of atoms



Gallium crystals



quartz



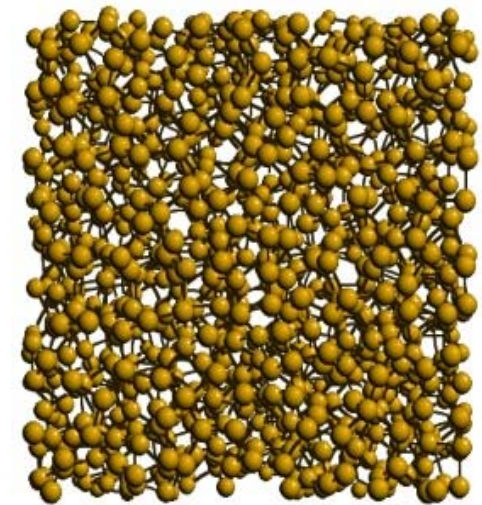
Insulin crystals



amorphous metal



glass

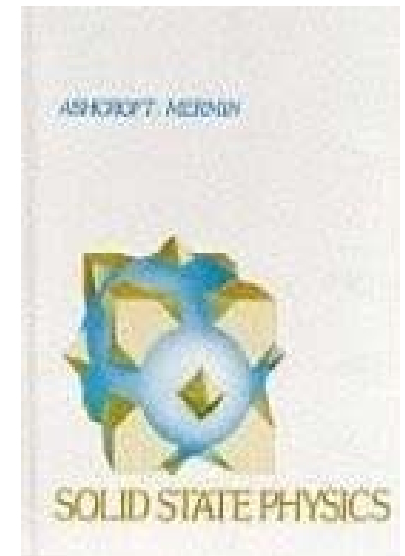
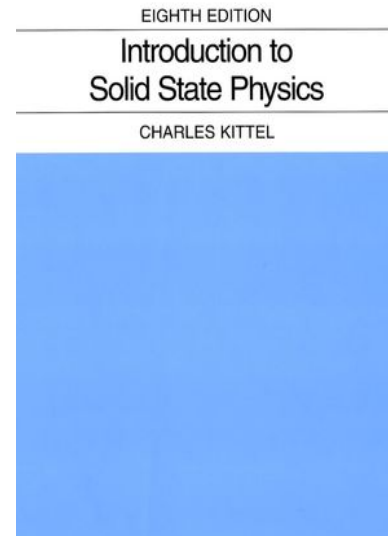
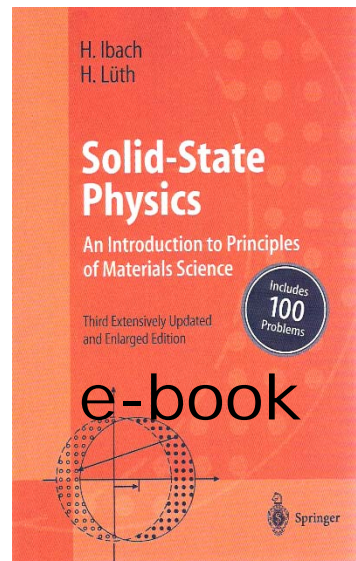
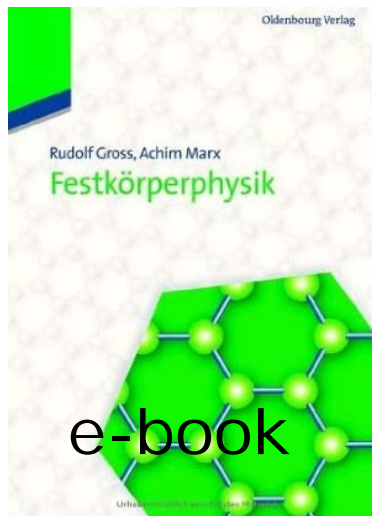
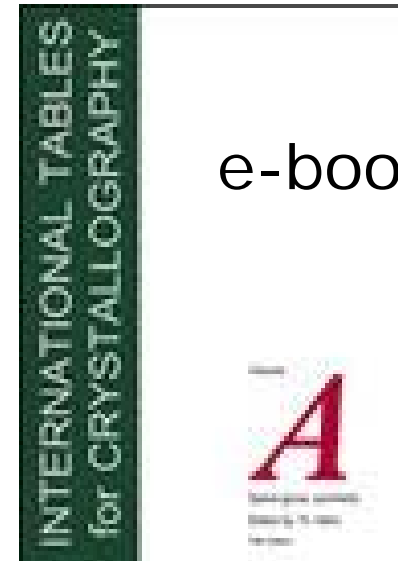
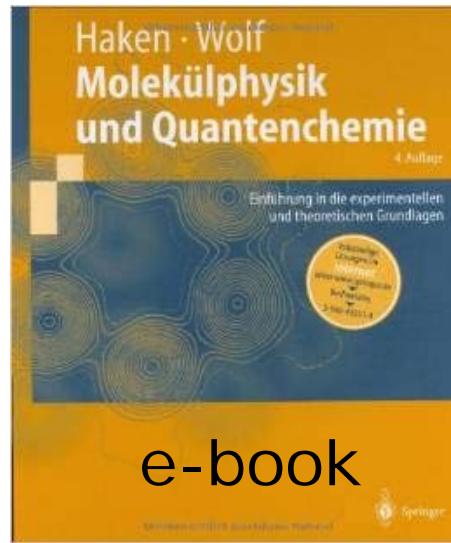


amorphous silicon

<http://www.wikipedia.org>

[http://www.pmc.umontreal.ca/~mousseau/site\\_an/uploads/Main/si1000.jpg](http://www.pmc.umontreal.ca/~mousseau/site_an/uploads/Main/si1000.jpg)

# Books





<http://www.if.tugraz.at/ss1.html>



Menu

Sections

PHY.F20 Molecular and Solid State Physics

## Course outline

- **Introduction** [▶ EN 6:45](#)
- **Review of atomic physics**
  - The solutions to the Schrödinger equation for the hydrogen atom [▶ DE 5:13](#)
    - Plots of the atomic orbitals
  - Helium
  - Many-electron wavefunctions
  - Slater determinants **W**
  - Singlet and triplet states
  - Exchange **W**
  - The intractability of the Schrödinger equation
  - Many-electron atoms
- **Molecules**
  - Molecular orbital theory **W**
    - Solving the total molecular Hamiltonian **W**
      - The Born-Oppenheimer approximation **W**
      - Many-electron wavefunctions
      - Bond potentials
      - Vibrational states
      - Rotational states
    - Solving the molecular orbital Hamiltonian

# Student Projects

Do something that will help other students



Google Custom



Menu

Sections

PHY.F20 Molecular and Solid State Physics

## Student projects

- Over the years there have been many contributions to the course from students and now we need to correct errors and remove unclear contributions before too much more is added. If you find something that is wrong or unclear and can improve it, that would be a suitable project.
- Often in atomic and molecular physics it is necessary to evaluate matrix elements of the form,  $\langle \phi_m | H | \phi_n \rangle$ . Where  $\phi_n$  are the atomic orbitals. To evaluate the integral, we need to know the atomic orbitals and the Laplacians of the atomic orbitals,  $\nabla^2 \phi_n$ . A student made a list of the first few atomic orbitals which can be found at the bottom of the page on [atomic orbitals](#). This list could be expanded.
- There are some examples of using the atomic orbitals to calculate matrix elements such as the [1s orbital](#) and the [2p orbital](#). Code in various languages such as Fortran, c, Java, JavaScript, Matlab, and Mathematica would be useful to have.
- Check the integration and the numerical calculations on the pages for these atomic orbitals [1s](#), [2s](#), [2p<sub>z</sub>](#).
- Plot the bond potential for a hydrogen molecular ion  $H_2^+$ . See the discussion of the [Hydrogen molecular ion  \$H\_2^+\$](#) .
- Calculate the molecular orbitals of ethylene, and butadiene similar to the [calculation for benzene](#). Since these molecules are chain of carbon atoms, refer to the discussion of [molecular chains](#). Separate projects would be the numerical calculation of  $H_{11}$ ,  $H_{12}$ , and  $S_{12}$  for ethylene, butadiene, or benzene.
- Upload a video to YouTube (<10 minutes) that explains some topic in the course outline.
- Upload a video to YouTube that explains how to use a piece of laboratory equipment in the physics building related to this course.
  - x-ray diffractometer

# Review of atomic physics

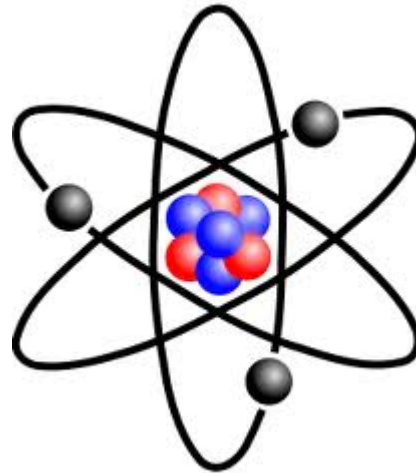
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Estimating the size of an atom

The hydrogen atom

The helium atom

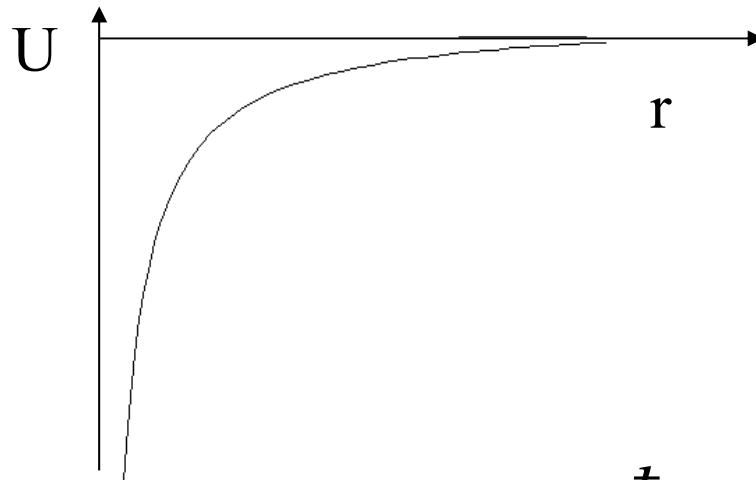
Many electron atoms



# Estimate the size of a hydrogen atom

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Potential energy  $U(r) = -\frac{e^2}{4\pi\epsilon_0 r}$



Uncertainty relation  $\Delta x \Delta p_x \geq \frac{\hbar}{2}$

For an atom:  $\Delta x \sim r_0$

$$\Delta p_x \geq \frac{\hbar}{2r_0}$$

# Estimate the size of a hydrogen atom

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$$\Delta p_x \geq \frac{\hbar}{2r_0}$$

$$\Delta p_x = \sqrt{\langle p_x^2 \rangle - \langle p_x \rangle^2} \quad \langle p_x \rangle = 0$$

$$(\Delta p_x)^2 = \langle p_x^2 \rangle \geq \left( \frac{\hbar}{2r_0} \right)^2$$

$$E_{kin} = \frac{mv^2}{2} = \frac{p^2}{2m}$$

$$\text{Kinetic energy in } x\text{-direction} = \langle E_{kin} \rangle = \frac{\langle p_x^2 \rangle}{2m} \geq \frac{\hbar^2}{8mr_0^2}$$

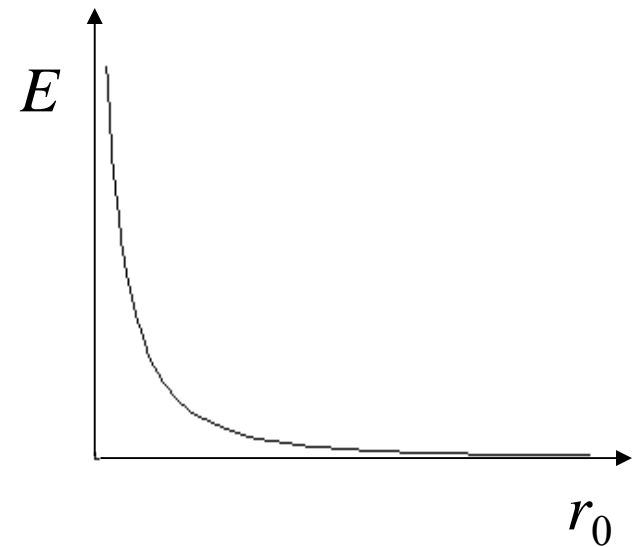
# Confinement energy

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$$\text{Kinetic energy in } x\text{-direction} = \langle E_{kin} \rangle = \frac{\langle p_x^2 \rangle}{2m} \geq \frac{\hbar^2}{8mr_0^2}$$

Confinement energy:

$$\frac{\langle p_x^2 \rangle}{2m} + \frac{\langle p_y^2 \rangle}{2m} + \frac{\langle p_z^2 \rangle}{2m} \geq \frac{3\hbar^2}{8mr_0^2}$$



# Estimate the size of a hydrogen atom

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Total energy = Kinetic + Potential

$$E_{tot} = \frac{3\hbar^2}{8mr^2} - \frac{e^2}{4\pi\epsilon_0 r}$$

$$\frac{dE_{tot}}{dr} = \frac{-3\hbar^2}{4mr^3} + \frac{e^2}{4\pi\epsilon_0 r^2}$$

$$r_0 = \frac{3\hbar^2 \pi\epsilon_0}{me^2} = 4.0 \times 10^{-11} \text{ m}$$

$$a_0 = 5.3 \times 10^{-11} \text{ m}$$

# Confinement energy

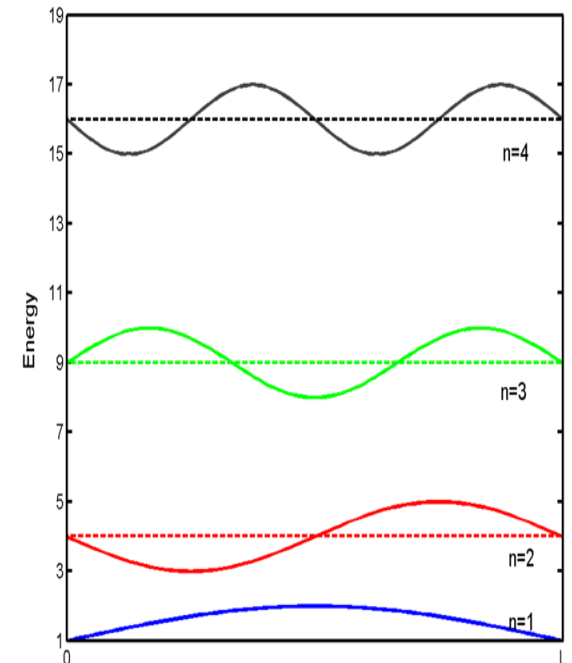
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$$\frac{-\hbar^2}{2m} \nabla^2 \Psi - \frac{e^2}{4\pi\epsilon_0 r} \Psi = E\Psi$$

The kinetic energy term increases as the wavelength gets smaller

$$E_{kin} = \frac{1}{2}mv^2 = \frac{p^2}{2m} = \frac{\hbar^2 k^2}{2m} = \frac{h^2}{2m\lambda^2}$$

$$p = mv \quad p = \hbar k \quad k = \frac{2\pi}{\lambda}$$





# Wave functions of hydrogen

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$$\frac{-\hbar^2}{2m} \nabla^2 \Psi - \frac{e^2}{4\pi\epsilon_0 r} \Psi = E\Psi$$

Solve with the boundary condition  $\Psi \rightarrow 0$  as  $|\vec{r}| \rightarrow \infty$

Assume  $\Psi(r, \theta, \varphi) = R(r)\Theta(\theta)\Phi(\varphi)$

# Hydrogen atom

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$$-\frac{\hbar^2}{2m}\nabla^2\psi - \frac{e^2}{4\pi\epsilon_0 r}\psi = E\psi$$

$$\psi_{nlm}(r, \theta, \varphi) = \sqrt{\left(\frac{2}{na_0}\right)^3 \frac{(n-l-1)!}{2n((n+l)!)} e^{-\rho/2} \rho^l L_{n-l-1}^{2l+1}(\rho) Y_{lm}(\theta, \varphi)}$$

$$\rho = \frac{2r}{na_0} \quad a_0 = \text{Bohr radius}$$

$$L_{n-l-1}^{2l+1}(\rho) = \text{generalized Laguerre polynomials}$$

$$Y_{lm}(\theta, \varphi) = \text{spherical harmonics (appear in centrosymmetric problems)}$$

# Hydrogen wavefunctions

quantum numbers  $n, l, m$

$$l = 0 \dots n-1$$

$$m = -l \dots 0 \dots l$$

$$l = 0 \rightarrow s$$

$$l = 1 \rightarrow p$$

$$l = 2 \rightarrow d$$

$$l = 3 \rightarrow f$$

$$\psi_{1s} = \frac{1}{\sqrt{\pi a_0^3}} e^{-\frac{r}{a_0}},$$

$$\psi_{2s} = \frac{1}{4\sqrt{2\pi a_0^3}} \left(2 - \frac{r}{a_0}\right) e^{-\frac{r}{2a_0}},$$

$$\psi_{2px} = \frac{1}{8\sqrt{\pi a_0^3}} \frac{r}{a_0} e^{-\frac{r}{2a_0}} \sin \theta \cos \varphi,$$

$$\psi_{2py} = \frac{1}{8\sqrt{\pi a_0^3}} \frac{r}{a_0} e^{-\frac{r}{2a_0}} \sin \theta \sin \varphi,$$

$$\psi_{2pz} = \frac{1}{4\sqrt{2\pi a_0^3}} \frac{r}{a_0} e^{-\frac{r}{2a_0}} \cos \theta.$$

Atomic Mass																					
1 amu = 1.66054 × 10 <sup>-27</sup> kg																					
1 H 1.008																	2 He 4.003				
3 Li 6.941	4 Be 9.012															5 B 10.81	6 C 12.01	7 N 14.01	8 O 16	9 F 19	10 Ne 20.18
11 Na 22.99	12 Mg 24.31															13 Al 26.98	14 Si 28.09	15 P 30.97	16 S 32.07	17 Cl 35.45	18 Ar 39.95
19 K 39.1	20 Ca 40.08	21 Sc 44.96	22 Ti 47.88	23 V 50.94	24 Cr 51.94	25 Mn 54.94	26 Fe 55.85	27 Co 58.93	28 Ni 58.69	29 Cu 63.55	30 Zn 65.39	31 Ga 69.72	32 Ge 72.59	33 As 74.92	34 Se 78.96	35 Br 79.9	36 Kr 83.8				
37 Rb 85.47	38 Sr 87.62	39 Y 88.91	40 Zr 91.22	41 Nb 92.91	42 Mo 95.94	43 Tc 98	44 Ru 101.1	45 Rh 102.9	46 Pd 106.4	47 Ag 107.9	48 Cd 112.4	49 In 114.8	50 Sn 118.7	51 Sb 121.8	52 Te 127.6	53 I 126.9	54 Xe 131.3				
55 Cs 132.9	56 Ba 137.3	57 La 138.9	58 Ce 140.9	59 Pr 144.2	60 Nd 147	61 Pm 150.4	62 Sm 152	63 Eu 157.3	64 Gd 158.9	65 Tb 162.5	66 Dy 164.9	67 Ho 167.3	68 Er 168.9	69 Tm 173	70 Yb 173	71 Lu 175					
87 Fr 223	88 Ra 226	89 Ac 227	90 Th 232	91 Pa 231	92 U 238	93 Np 237	94 Pu 242	95 Am 243	96 Cm 247	97 Bk 247	98 Cf 251	99 Es 252	100 Fm 253	101 Md 258	102 No 259	103 Lr 260					