

PHY404- Solid State Physics II

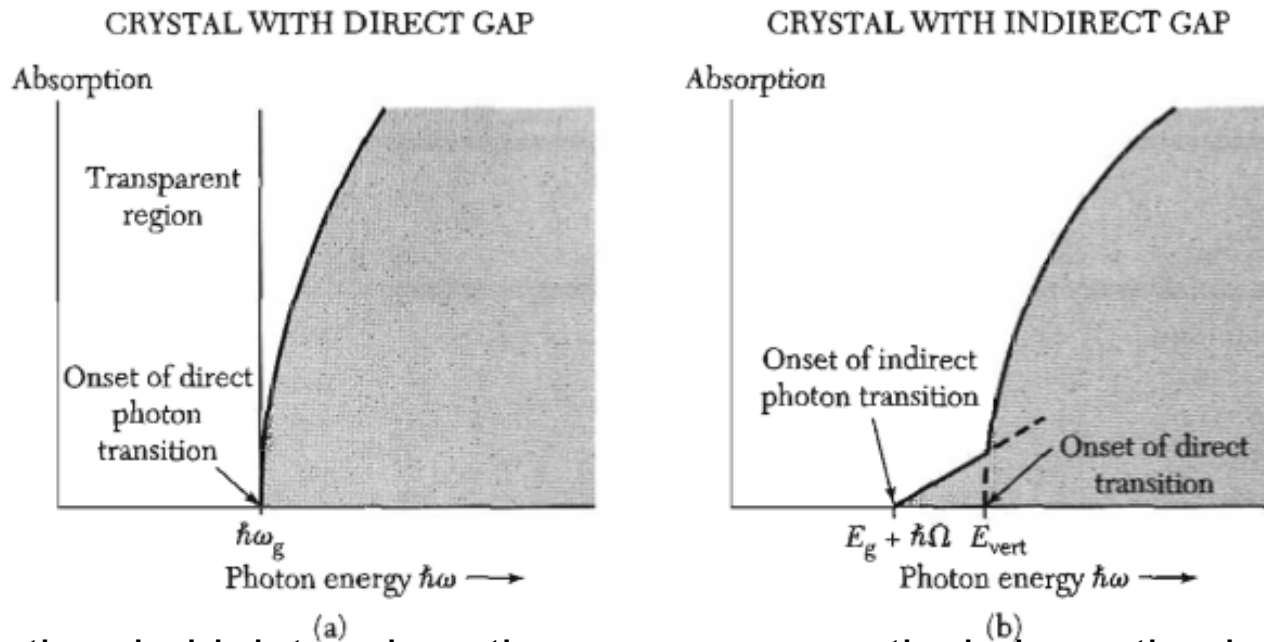
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Energy gap between valence and conduction bands

Crystal	Gap	E_g , eV		Crystal	Gap	E_g , eV	
		0 K	300 K			0 K	300 K
Diamond	<i>i</i>	5.4		SiC(hex)	<i>i</i>	3.0	—
Si	<i>i</i>	1.17	1.11	Te	<i>d</i>	0.33	—
Ge	<i>i</i>	0.744	0.66	HgTe ^a	<i>d</i>	-0.30	
α Sn	<i>d</i>	0.00	0.00	PbS	<i>d</i>	0.286	0.34–0.37
InSb	<i>d</i>	0.23	0.17	PbSe	<i>i</i>	0.165	0.27
InAs	<i>d</i>	0.43	0.36	PbTe	<i>i</i>	0.190	0.29
InP	<i>d</i>	1.42	1.27	CdS	<i>d</i>	2.582	2.42
GaP	<i>i</i>	2.32	2.25	CdSe	<i>d</i>	1.840	1.74
GaAs	<i>d</i>	1.52	1.43	CdTe	<i>d</i>	1.607	1.44
GaSb	<i>d</i>	0.81	0.68	SnTe	<i>d</i>	0.3	0.18
AlSb	<i>i</i>	1.65	1.6	Cu ₂ O	<i>d</i>	2.172	—

i: indirect gap, d: direct gap

Optical Adsorption in Pure Insulators



In (a) The threshold determines the energy gap, optical absorption is weaker near the threshold.

In (b) the energy E , marks the threshold for the creation of a free electron and a free hole, with no phonon involved.

Such a transition is called vertical; it is similar to the direct transition in (a).

In a **direct adsorption process** the threshold of continuous optical adsorption at frequency w_g measures the band gap. A photon is adsorbed by the crystal with the creation of an electron and a hole.

In the **indirect absorption process** the minimum energy gap of the band structure involves electrons and holes seperated by a substantial wavevector \mathbf{k} .

❑ Optical measurements determines whether the gap is direct or indirect.

Ex: The band edges in Ge and in Si are connected by indirect transitions, the band edges in InSb and GaAs are connected by a direct transition.

❑ HgTe and HgSe are semimetals and have negative gaps- the conduction and valence bands overlap.

EQUATIONS OF MOTION

- We derive the equation of motion of an electron in an energy band.
- We look at the notion of a wave packet in an applied electric field.
- Suppose that the wave packet is made up of wavefunctions assembled near a particular wavevector \mathbf{k}

EQUATIONS OF MOTION

The frequency associated with a wavefunction of energy;

$$v_g = \hbar^{-1} d\epsilon/dk \quad \text{or} \quad \mathbf{v} = \hbar^{-1} \nabla_{\mathbf{k}} \epsilon(\mathbf{k})$$

$$v_g = d\omega/dk \quad \text{The group velocity}$$

The work done on the electron by the electric field \mathbf{E} in the time interval is:

$$\delta\epsilon = -eE v_g \delta t$$

External force \mathbf{F} is:

$$\hbar \frac{d\mathbf{k}}{dt} = \mathbf{F} .$$

EQUATIONS OF MOTION

We examine the transfer of momentum between the electron and the lattice when the state \mathbf{k} of the electron is changed to $\mathbf{k} + \Delta\mathbf{k}$ by the application of an external force. We imagine an insulating crystal electrostatically neutral except for a single electron in the state \mathbf{k} of an otherwise empty band.

We suppose that a weak external force is applied for a time interval such that the total impulse given to the entire crystal system is $\mathbf{J} = \int \mathbf{F} dt$. If the conduction electron were free ($m^* = m$), the total momentum imparted to the crystal system by the impulse would appear in the change of momentum of the conduction electron:

$$\mathbf{J} = \Delta\mathbf{p}_{\text{tot}} = \Delta\mathbf{p}_{\text{el}} = \hbar\Delta\mathbf{k}$$