

1:

$$D(E)dE = \frac{1}{(2\pi)^2} \int dS_k \quad \int dS_k = \int dl_E dK_\perp$$

$dl_E$  is the line of constant energy, defined by  $E = \frac{\hbar^2}{2m}(k_x^2 - k_y^2)$

$$dl_E = \sqrt{dk_x^2 + dk_y^2}$$

$$k_y = \sqrt{\frac{2mE}{\hbar^2} - dk_x^2}$$

$$dk_y = \frac{k_x dk_x}{\sqrt{\frac{2mE}{\hbar^2} - k_x^2}}$$

$$dk_y^2 = \frac{k_x^2 dk_x^2}{\frac{2mE}{\hbar^2} - k_x^2}$$

$$\text{So } dl_E = \sqrt{1 + \frac{k_x^2}{\frac{2mE}{\hbar^2} - k_x^2}} = \left( \frac{\frac{2mE}{\hbar^2}}{\frac{2mE}{\hbar^2} - k_x^2} \right)^{\frac{1}{2}}$$

$$\text{and } dE = \frac{\hbar^2}{m} \sqrt{\frac{2mE}{\hbar^2}} dK_\perp$$

$$\text{So } D(E) = \frac{1}{(2\pi)^2} \int \frac{\frac{2mE}{\hbar^2}}{\sqrt{\frac{2mE}{\hbar^2} - k_x^2}} \frac{dk_x dK_\perp}{\frac{\hbar^2}{m} \sqrt{\frac{2mE}{\hbar^2}} dK_\perp} = \frac{m}{(2\pi\hbar)^2} \int \frac{dk_x}{\sqrt{\frac{2mE}{\hbar^2} - k_x^2}} = \frac{m}{(2\pi\hbar)^2} \ln(k_x + \sqrt{\frac{2mE}{\hbar^2} - k_x^2})$$

Look at the edge of the Brillion Zone:  $E = E - E_0$ ,  $k_x \gg E - E_0$

$$\text{So } D(E) = \frac{m}{(2\pi\hbar)^2} \ln(k_x + \frac{2m(E-E_0)}{\hbar^2 k_x} - k_x)$$

$$\text{So } D(E) \approx \ln(E - E_0)$$



2:

$\epsilon_2 \propto (E - E_0)^{1/2}$  at the van hove singularity means that

$$\epsilon_2 \propto \frac{1}{w^2} (E - E_0)^{1/2}$$

$$\text{Then } \epsilon_1 = 1 + \frac{A}{\pi} P \int_{-\infty}^{\infty} \frac{\sqrt{w' - w_0}}{w'^2 (w' - w)} dw'$$

2 residues, at  $w' = 0$  and  $w' = w$

$w' = 0$  :

$$\begin{aligned} \pi \mathbf{i} \lim_{w' \rightarrow 0} \left( \frac{\partial}{\partial w'} \frac{\sqrt{w' - w_0}}{(w' - w)} \right) &= \pi \mathbf{i} \lim_{w' \rightarrow 0} \left( \frac{1}{2\sqrt{w' - w_0}(w' - w)} - \frac{\sqrt{w' - w_0}}{(w' - w)^2} \right) \\ &= \pi \frac{(-w + 2w_0)}{2w^2 \sqrt{w_0}} \end{aligned}$$

$w' = w$  :

$$\pi \mathbf{i} \frac{\sqrt{w - w_0}}{w^2} = \pi \frac{\sqrt{w_0 - w}}{w^2}$$

$$\text{So } \epsilon_1 = 1 + A (w - 2w_0 + 2\sqrt{w_0}\sqrt{w_0 - w}) / (2w^2 \sqrt{w_0})$$

3. Near  $k=0$ ,  $k$  is approximately constant, so we can take it out:  
$$\sum_n |\langle u_v | \vec{k} \cdot \hat{p} | u_n \rangle|^2 \approx \sum_n k^2 |\langle u_v | \vec{k} \cdot \hat{p} | u_n \rangle|^2$$

Let us also assume all but one band is sufficiently far away to be ignored, so that we ignore the series. Then

$$\frac{1}{m^*} = \frac{1}{m} + \frac{2}{m^2} \frac{|\langle u_v | \hat{p} | u_c \rangle|^2}{E_v - E_n}$$

$$|\langle u_v | \hat{p} | u_c \rangle|^2 = P, \text{ so}$$

$$\frac{1}{m^*} = \frac{1}{m} + \frac{2}{m^2} \frac{P^2}{E_g}$$

$$\frac{E_g}{4} \left( \frac{1}{m^*} - \frac{1}{m} \right) = \frac{P^2}{2m^2}$$

$$\frac{P^2}{2m} = \frac{E_g}{4} \left( \frac{m}{m^*} - 1 \right)$$

b) These do not work in this model, as Si and Ge are indirect band gap materials. Therefore we cannot remove  $k$  from the sum.

3.

	Energy Gap (eV) m*	Lattice Constant	P^2	P^2/2m
Si	1.107	1.100	5.43072	-2.571E+04 -0.025159091
Ge	0.670	0.550	5.65754	1.401E+05 0.137045455
Alpha - Sn	0.080	0.020	6.49120	1.002E+06 0.98
Te	0.330	0.190	4.45000	3.594E+05 0.351710526
III-V compounds				
AlSb	1.600	0.090	6.13550	4.133E+06 4.044444444
GaP	2.240	0.350	5.49050	1.063E+06 1.04
GaAs	1.350	0.068	5.65315	4.728E+06 4.625735294
GaSb	0.670	0.050	6.09540	3.253E+06 3.1825
InP	1.270	0.067	5.86875	4.519E+06 4.42130597
InAs	0.360	0.022	6.05838	4.089E+06 4.000909091
InSb	0.165	0.014	6.47877	2.969E+06 2.905178571
II-VI compounds				
ZnO	3.200	0.380	4.63000	1.334E+06 1.305263158
CdO	2.500	0.100	4.69530	5.749E+06 5.625
CdS	2.400	0.165	5.83200	3.103E+06 3.036363636
CdSe	1.740	0.130	6.05000	2.975E+06 2.911153846
HgSe	0.300	0.030	6.08400	2.478E+06 2.425
HgTe	0.150	0.017	6.46230	2.216E+06 2.168382353
Halite structure compounds				
PbS	0.370	0.160	5.93620	4.963E+05 0.485625
PbSe	0.260	0.300	6.12430	1.550E+05 0.151666667
PbTe	0.250	0.210	6.45400	2.403E+05 0.235119048
Others				
Bi2Te3	0.130	0.580	4.38000	2.405E+04 0.023534483
Mg2Si	0.770	0.460	6.33800	2.309E+05 0.225978261
Mg2Sn	0.210	0.370	6.76500	9.136E+04 0.089391892

Average: 2.016877623 eV