


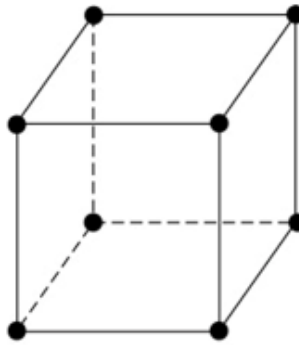
HW1
Pete Mills
Sept 7, 2023

Problem 1

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Problem 1

- (a) A simple cubic structure consists of a single atom in the center of the cube. The lattice constant is a_0 . Determine the volume density of atoms and the surface density of atoms in the $\{110\}$ plane.
- (b) Compare the results of part (a) to the results for the case of the simple cubic structure shown below with the same lattice constant.



a) $\rho_{\text{bcc}} = 2 \text{ atoms} / a_0^3$, $\sigma_{\text{bcc}} = 5 \text{ atoms} / a_0^2$

b) $\rho_{\text{sc}} = 1 \text{ atoms} / a_0^3$, $\sigma_{\text{sc}} = 0.5 \text{ atoms} / a_0^2$,

Therefore,
BCC structure has 2x the volume density as SC structure.
BCC structure has 10x the surface density of SC structure.

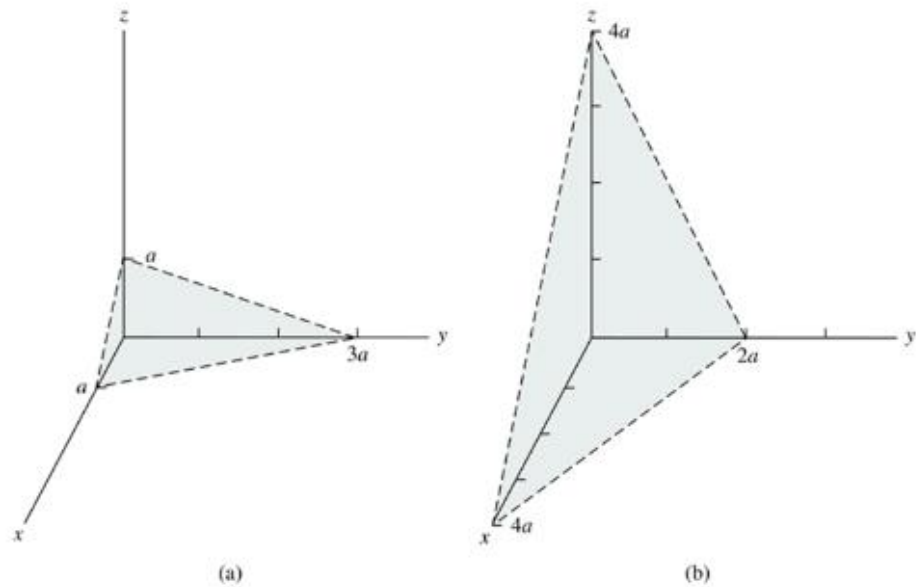
Problem 2

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Problem 2

For a simple cubic lattice, determine the Miller indices for the planes shown in the figure below.



a) $p=1, q=3, s=1 \rightarrow (1/1, 1/3, 1/1) \rightarrow (3/3, 1/3, 3/3) \rightarrow (3, 1, 3)$

b) $p=4, q=2, s=4 \rightarrow (1/4, 1/2, 1/4) \rightarrow (1/4, 2/4, 1/4) \rightarrow (1, 2, 1)$

Problem 3

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Problem 3

Calculate the de Broglie wavelength for

- a) An electron with kinetic energy of
 - i. 1.0 eV
 - ii. 100 eV
- b) A proton with kinetic energy of 1 eV.

```
eoo-331 > hw > hw1 > broglie.py > ...
1  import numpy as np
2
3  # Planck's constant J/s
4  h = 6.62607015e-34
5
6  # Define 1 eV in Joules
7  eV_to_Joules = 1.60219e-19
8
9  def calculate_de_broglie_wavelength(mass_kg, kinetic_energy_J):
10     # Calculate momentum from kinetic energy and mass
11     momentum = np.sqrt(2 * mass_kg * kinetic_energy_J)
12     # Calculate Broglie wavelength using momentum
13     wavelength = h / momentum
14     return wavelength
15
16 # Input parameters
17 mass_proton_kg = 1.6726219e-27 # Mass of a proton in kilograms
18 mass_electron_kg = 9.10938356e-31 # Mass of an electron in kilograms
19 energies = [1.0, 100.0] # Kinetic energy in electronvolts
20
21 for kinetic_energy_eV in energies:
22     # Convert kinetic energy from eV to Joules
23     kinetic_energy_J = kinetic_energy_eV * eV_to_Joules
24
25     # Calculate Broglie wavelengths
26     wavelength_electron = calculate_de_broglie_wavelength(mass_electron_kg, kinetic_energy_J)
27     wavelength_proton = calculate_de_broglie_wavelength(mass_proton_kg, kinetic_energy_J)
28
29     # Print results
30     print(f"Broglie wavelength for an electron with {kinetic_energy_eV} eV kinetic energy: {wavelength_electron:.6e} meters")
31     print(f"Broglie wavelength for a proton with {kinetic_energy_eV} eV kinetic energy: {wavelength_proton:.6e} meters")
32
```

PROBLEMS OUTPUT DEBUG CONSOLE TERMINAL GITLENS

```
pete@pete-Precision-5550:~/code/eoo$ ls
books eeo-224 eeo-315 eeo-331 eeo-352 README.md
pete@pete-Precision-5550:~/code/eoo$ cd eeo-331/hw/hw1/
pete@pete-Precision-5550:~/code/eoo/eoo-331/hw/hw1$ ls
bandgap.py broglie.py
pete@pete-Precision-5550:~/code/eoo/eoo-331/hw/hw1$ python ./broglie.py
Command 'python' not found, did you mean:
  command 'python3' from deb python3
  command 'python' from deb python-is-python3
pete@pete-Precision-5550:~/code/eoo/eoo-331/hw/hw1$ python3 ./broglie.py
Broglie wavelength for an electron with 1.0 eV kinetic energy: 1.226421e-09 meters
Broglie wavelength for a proton with 1.0 eV kinetic energy: 2.862102e-11 meters
Broglie wavelength for an electron with 100.0 eV kinetic energy: 1.226421e-10 meters
Broglie wavelength for a proton with 100.0 eV kinetic energy: 2.862102e-12 meters
pete@pete-Precision-5550:~/code/eoo/eoo-331/hw/hw1$
```

Problem 4

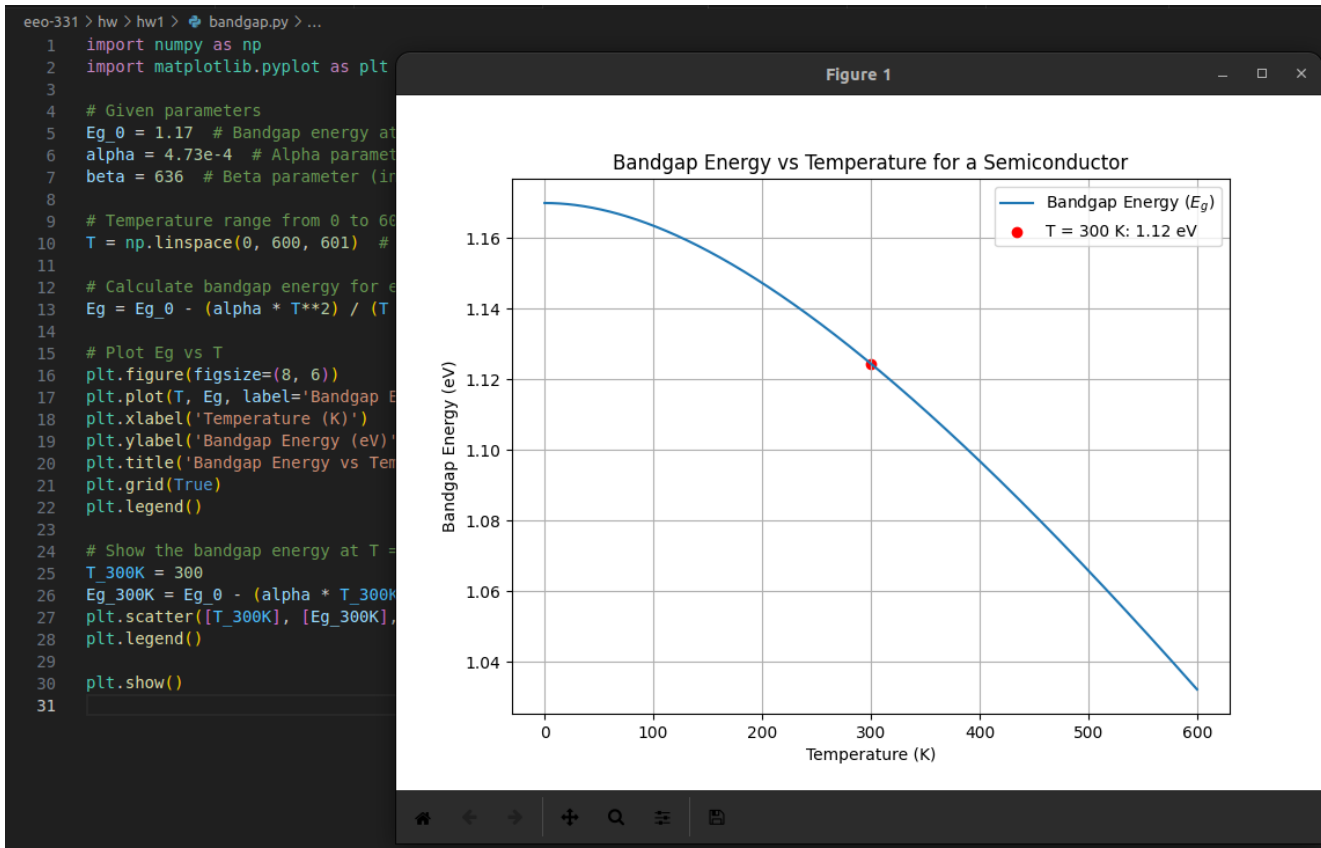
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Problem 4

The bandgap energy in a semiconductor is usually a slight function of temperature. In some cases, the bandgap energy versus temperature can be modeled by

$$E_g = E_g(0) - \frac{\alpha T^2}{(\beta + T)}$$

Where $E_g(0)$ is the value of the bandgap energy at $T = 0$ K. For silicon, the parameter values are $E_g(0) = 1.17$ eV, $\alpha = 4.73 \times 10^{-4}$ eV/K and $\beta = 636$ K. Plot E_g versus T over the range $0 \leq T \leq 600$ K. In particular, note the value at $T = 300$ K.



Problem 5

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Problem 5

The forbidden energy band of GaAs is 1.42 eV.

- a) Determine the minimum frequency of an incident photon that can interact with a valence electron and elevate the electron to the conduction band.
- b) What is the corresponding wavelength?

$$E = hf$$

$$E = 1.42 \text{ eV} \times 1.60217663 \times 10^{-19} \text{ J/eV}$$

$$f_{\min} = E/h$$

$$f_{\min} = E / 6.62607015 \times 10^{-34}$$

$$f_{\min} = 2.854 \times 10^{14} \text{ Hz is the minimum frequency of an incident photon.}$$

$$\lambda = c/f_{\min} \text{ where } c = 3 \times 10^8 \text{ m/s}$$

$$\lambda = 1.05 \times 10^{-6} \text{ m is the corresponding wavelength.}$$