So scattering REALLY attorvates net charge movement rate in a metal.

ii
$$\lambda = \frac{(9.11 \times 10^{-21} \log (1.56 \times 10^{6} \text{m/s}) (5.9 \times 10^{7} \text{sz}^{-1} \text{m}^{-1})}{(8.45 \times 10^{28} \text{m}^{-3})(1.602 \times 10^{-19})^{2}}$$

Internet tells me lattre constant in copper is ~3.61 å (FCC lattre so mean spacing is probably even smaller but same order of mag).

Moughly speaking a conduction electron moves ~10° atomiz distances before scattering?

4.3) a & bac) Read the book

d) The linear ferm comes from the electrons and the T3 term comes from phanons (remember Cph Td from hah #1)

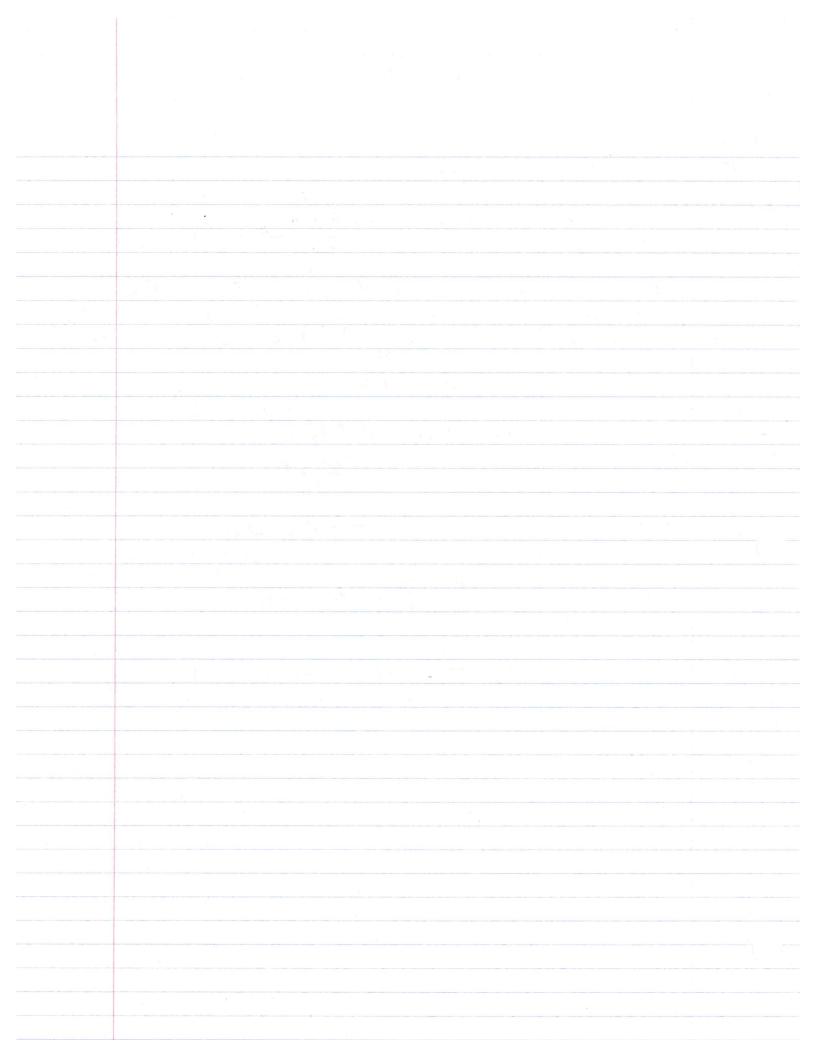
A variety of ways to do this, one is to use the coefficient y along of book derivations of the heat capacity due to elec, which depends on g(E=) to solve for E=

(5) $C_{chc} = \frac{\pi^2}{3} k_B^2 g(E_F) T V$ Now $\frac{\pi^2}{3} k_B^2 g(E_F) V = \frac{\pi^2}{3} k_B^2 \frac{3\pi V}{2E_F} \left(\frac{E_F}{E_F}\right)^{1/2}$ $= \frac{\pi^2}{2} k_B^2 E_F$

= R2 k3 = rmNA
2 # Moles

So $\gamma = \frac{\Re^2}{2 \, \text{kg}} \frac{\text{Na}}{\text{Ep}} = 2.08 \times 10^{-3} \frac{\text{T}}{\text{mol·K}^2}$

 $E_{F} = \frac{\pi^{2}}{2} \frac{k_{B}^{2} N_{A}}{2.08 \times 10^{-3} \text{ J/}} = 2.72 \times 10^{-19} \text{ J} = 1.70 \text{ eV}$



47) Can use con 48 in kext

6)
$$E = V \int de e g(e) n_{\mu} (\beta(e-e_{\mu})) = V \int de e \frac{3n_{\mu} - e_{\mu}}{E_{\mu}} \frac{1}{2} e^{N(e-e_{\mu})} + 1$$

$$= V \int de e \frac{3n_{\mu}}{2E_{\mu}} \left(\frac{e^{-N(e-e_{\mu})}}{E_{\mu}} + \frac{$$

Different periods in table, perhaps core electrons are contributing to difference?

Eurobrestmate

B_K = 2.88×109 Pa = [2.88 GPa]

Estimate Pressure needed to squeeze electrons. B=Vav ~ Y AV -> B = DOIN AP DP=0.013 DPNa = 0.01 (8.42GPa) = 84.2MPa

6.6

H, =
$$\chi$$
 ($|R| + |R-\bar{r}| +$

Recall freld of dipole is

$$E = \frac{1}{4\pi\epsilon_0} \left(\frac{3(\vec{p}_0 \cdot \hat{p}_1) \cdot \vec{r}_0 - \vec{p}_1}{3} \right)$$

So
$$H \sim -(\vec{p}_2 \cdot \vec{E}) = \frac{1}{4\pi\epsilon_0} \left(\frac{-3(\vec{p}_1 \cdot \vec{r})(\vec{p}_2 \cdot \vec{r}) + \vec{p}_2 \cdot \vec{p}_1}{r^3} \right)$$

$$= \frac{e^2}{4\pi\epsilon_0} \left(\frac{-3x_1x_2 + x_1x_2 + y_1y_2 + z_1z_2}{r^3} \right)$$
Since $\vec{r} = -\hat{r}$

b) The first order correction would be AE(1)=(1,0,0;1,0,0) H, (1,0,0;1,0,0) However H, is add in each in each atom's position so it's add under inversion operator if. Note that G.S. is spherically symmetric and even under inversion. Generally, for each atom, 11,0,0> = 7/1,0,0> Dinverted State So it I take the expectator value of say x in the inverted state, I should not solve thing as normal state $\langle 1,0,0| \times, |1,0,0 \rangle = \langle 1,0,0| \times, |1,0,0 \rangle$ However X, as an operator is odd under inversion, wearing (1,0,01x,11,0,0) = (1,0,01(mx,m)1,0,0) = $\langle 1,0,0|X,11,0,0\rangle = \langle 1,0,0|X,11,0,0\rangle$ The only way this can be true is if the expec. value is zero. The first non zero term will come from expec values of form K1,0,0;11,0,01H, In, L, m; n2, l2, m2>1 ~ | R3 ~ R6 The denominator of SE contains E-E, which is always regative. Since the models squared in the numerator is always positive, each term in the sum for SE 13 negative. The perturbation drops the energy implying an attractive force.

The numerator is zero for the same reason as previous problem. To get finite expectation value from $\langle 1,0,0|X|n,l,m\rangle$ I must be I. The is impossible unless n > 2. Uper bound. SE = 57 <1,0,0;1,0,0 | H, In, I, m; n, I, m, in, I, m, i = <1,0,0;1,0,01H2 11,0,0;1,0,0> E, -E,... only the even terms in H, (i.e. x2x2 and not Z,Zzy, y2) contribute x(40x p3)(62)

So $|SE_3| = 2 \frac{e^2 a_0}{HE_0 R^6}$ My Not sove they book doesn't simplify $\frac{2}{4} \Rightarrow 2?$

ISE, I is almost exactly same except sur in denominator changes slightly.