ECE 659 EXAM V Monday May 1 2017, 8-10AM

NAME : SOLUTION

CLOSED BOOK

All four questions carry equal weight

Please show all work. No credit for just writing down the answer, even if correct. **5.1.** Consider a system having a very large number of *degenerate* levels with energy ε with an electron-electron interaction energy related to the number of electrons N by the relation

$$U(N) = \frac{U_0}{2} N(N-1)$$

At what value of the electrochemical potential μ , will the equilibrium number of electrons change from 100 to 101?

SOLUTION:

States	E - μN			
N = 100	$100(\epsilon - \mu) + U_0 \times 50 \times 99$			
N = 101	$101(\epsilon - \mu) + U_0 \times 50 \times 101$			

Transition from N=100 to N=101 occurs when

$$(E - \mu N)_{100} = (E - \mu N)_{101}$$

that is, when

$$\epsilon - \mu + 100U_0 = 0$$

 $\rightarrow \mu = \epsilon + 100U_0$

5.2. Consider a quantum dot with 4 degenerate levels, all with the same energy ε , in equilibrium with a reservoir with electrochemical potential μ and temperature T. If there were no interactions, the average number of electrons at equilibrium would be given by

$$< n > = \frac{4}{1 + e^{(\varepsilon - \mu)/kT}}$$

Obtain an expression for the average number of electrons if the electron-electron interaction is so large that no more than one of the 4 levels can be occupied at the same time.

SOLUTION:

There is one zero-electron state with probability

$$P_0 = \frac{1}{Z}$$

and 4 one-electron states each with a probability

$$P_{1} = \frac{e^{-x}}{Z}, \quad x \equiv \frac{\varepsilon - \mu}{kT}$$

$$< n > = \frac{\sum_{n} nP_{n}}{\sum_{n} P_{n}} = \frac{4P_{1}}{P_{0} + 4P_{1}}$$

$$< n > = \frac{4e^{-x}}{1 + 4e^{-x}} = -\frac{4}{4 + e^{x}} = -\frac{4}{4 + e^{(\varepsilon - \mu)/kT}}$$

5.3. Consider a single quantum dot with *just one level* with energy ε . We wish to find the current as a function of the electrochemical potential $\mu_2 = \mu_1 + qV$ (keeping μ_1 fixed) using the Fock space picture by writing

$$\frac{d}{dt} \begin{cases} P_0 \\ P_1 \end{cases} = \begin{cases} 0 \\ 0 \end{cases} = \begin{bmatrix} W_1 + W_2 \end{bmatrix} \begin{cases} P_0 \\ P_1 \end{cases}$$

where the W_1 and W_2 matrices describe the transition rates between the different states due to electron exchange with contacts 1 and 2 respectively which are held in equilibrium with electrochemical potentials μ_1 and μ_2 .

(a) Write down the matrices W_1 and W_2 is terms of the Fermi functions in contacts 1, 2.

- (b) Obtain the steady-state probabilities P_0 and P_1 .
- (c) Obtain the steady-state current.

SOLUTION:

(a)

$$W_1 = \gamma_1 \begin{bmatrix} -f_1 & 1-f_1 \\ +f_1 & -(1-f_1) \end{bmatrix} \qquad \qquad W_2 = \gamma_2 \begin{bmatrix} -f_2 & 1-f_2 \\ +f_2 & -(1-f_2) \end{bmatrix}$$

(b)

$$\begin{cases} 0 \\ 0 \\ 0 \\ \end{cases} = \begin{bmatrix} -\gamma_1 f_1 - \gamma_2 f_2 & \gamma_1 (1 - f_1) + \gamma_2 (1 - f_2) \\ \gamma_1 f_1 + \gamma_2 f_2 & -\gamma_1 (1 - f_1) - \gamma_2 (1 - f_2) \end{bmatrix} \begin{cases} P_0 \\ P_1 \\ \end{cases}$$

$$\frac{P_1}{P_0} = \frac{\gamma_1 f_1 + \gamma_2 f_2}{\gamma_1 (1 - f_1) + \gamma_2 (1 - f_2)} \rightarrow P_1 = \frac{\gamma_1 f_1 + \gamma_2 f_2}{\gamma_1 + \gamma_2}, \quad P_0 = 1 - P_1$$

(c)

$$I = q\gamma_1 \{ 0 \ 1 \} \begin{bmatrix} -f_1 & 1-f_1 \\ +f_1 & -(1-f_1) \end{bmatrix} \begin{cases} P_0 \\ P_1 \end{bmatrix} = q\gamma_1 (f_1 P_0 - (1-f_1)P_1) = \gamma_1 (f_1 - P_1)$$
$$= q\gamma_1 \left(f_1 - \frac{\gamma_1 f_1 + \gamma_2 f_2}{\gamma_1 + \gamma_2} \right) = q \frac{\gamma_1 \gamma_2}{\gamma_1 + \gamma_2} (f_1 - f_2)$$

5.4. Consider two coupled quantum dots each with two spin-degenerate levels, described by the one-electron Hamiltonian,

$$h = \begin{bmatrix} \varepsilon & t & 0 & 0 \\ t & \varepsilon & 0 & 0 \\ 0 & 0 & \varepsilon & t \\ 0 & 0 & t & \varepsilon \end{bmatrix} \qquad \qquad \mu \qquad \underbrace{\varepsilon}_{t} \underbrace{\varepsilon}_{t} \qquad \underbrace{\varepsilon}_{t} \qquad \mu \qquad \underbrace{\varepsilon}_{t} \underbrace{\varepsilon}_{t} \qquad \underbrace{\varepsilon}_{t} \qquad \mu \qquad \underbrace{\varepsilon}_{t} \underbrace{\varepsilon}_{t} \qquad \underbrace{\varepsilon}_{$$

having an intra-dot interaction energy U and zero inter-dot interaction energy.

We have seen that the Hamiltonian for two-electron states can be written as

	u_1d_1	u_2d_2	u_1d_2	$u_2 d_1$	u_1u_2	$d_1 d_2$
<i>H</i> ₂ =	$\left\lceil 2\varepsilon + b \right\rceil$	U = 0	t	t	0	0]
	0	$U 0 \\ 2\varepsilon + U$	t	t	0	0
	t	t	2ε	0	0	0
	t	t	0	2ε	0	0
	0	0	0	0	2ε	0
	0	0	0	0	0	2ε

which can be transformed into

Explain how you would obtain the lowest energy eigenvalue and the corresponding eigenvector making use of the result that the eigenvectors of

 $\vec{\sigma}.\hat{n} \equiv \sigma_x \sin\theta \cos\phi + \sigma_y \sin\theta \sin\phi + \sigma_z \cos\theta$

corresponding to eigenvalues +1 and -1 can be written as

$$\begin{cases} c \\ s \\ \end{cases}, \begin{cases} -s^* \\ c^* \end{cases}$$

respectively, where

$$c \equiv \cos(\theta/2) e^{-i\varphi/2}$$
, $s \equiv \sin(\theta/2) e^{+i\varphi/2}$

SOLUTION:

The lowest energy eigenvalue comes from the 2x2 block

$$\begin{array}{c|c} \frac{u_1d_1+u_2d_2}{\sqrt{2}} & \frac{u_1d_2+u_2d_1}{\sqrt{2}} \\ \\ \left[2\epsilon+U & 2t \\ 2t & 2\epsilon \end{array} \right] \end{array}$$

which can be written as

$$\begin{pmatrix} 2\epsilon + \frac{U}{2} \end{pmatrix} I + \frac{U}{2}\sigma_z + 2t\sigma_x \\ = \left(2\epsilon + \frac{U}{2}\right)I + B\cos\theta \ \sigma_z + B\sin\theta \ \sigma_x \\ = \left(2\epsilon + \frac{U}{2}\right)I + B \ \vec{\sigma}.\hat{n} \\ where \ B = \sqrt{\left(\frac{U}{2}\right)^2 + (2t)^2} \\ and \ tan\theta = \frac{4t}{U} \end{cases}$$

The lowest eigenvalue is

$$\left(2\epsilon + \frac{U}{2}\right) - B = \left(2\epsilon + \frac{U}{2}\right) - \sqrt{\left(\frac{U}{2}\right)^2 + (2t)^2}$$

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And the corresponding eigenvector is

$$-sinrac{ heta}{2}\;rac{u_1d_1+u_2d_2}{\sqrt{2}}+cosrac{ heta}{2}\;rac{u_1d_2+u_2d_1}{\sqrt{2}}$$

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Have a great summer !!