

EXERCISE 1

CHARGE STABILITY DIAGRAM OF A SINGLE QD

$$1) Q_G = C_G (V_G - V_A) \quad (a)$$

$$Q_A = C_A V_A \quad (b)$$

$$Q_2 = C_2 V_2 \quad (c)$$

$$Q_A - Q_2 - Q_G = -N|e| \quad (d)$$

Sum eq. from the circuit: $V_A + V_2 = V_D \quad (e)$

$$E_{ex} = \frac{1}{2} (Q_G V_G - N|e| V_A + Q_2 V_D) \quad (f)$$

(a), (b), (c) into (d):

$$C_A V_A - C_2 V_2 - C_G (V_G - V_A) = -N|e|$$

$\hookrightarrow V_D - V_A \quad (e)$

$$(C_A + C_2 + C_G) V_A = C_2 V_D + C_G V_G - N|e|$$

$$\rightarrow V_A = \frac{C_G V_G + C_2 V_D - N|e|}{C_A + C_2 + C_G} \quad (g)$$

$$\rightarrow V_2 = V_D - V_A = \frac{C_A V_A + C_2 V_D + C_G V_G - C_G V_G - C_2 V_A + N|e|}{C_A + C_2 + C_G} =$$
$$= \frac{(C_A + C_G) V_D - C_G V_G + N|e|}{C_A + C_2 + C_G} \quad (h)$$

(g), (h), (e), (c) into (f)

$$\begin{aligned}
 E_{el} &= \frac{1}{2} \left(C_G \left(V_G - \frac{C_G V_G + C_2 V_D - N|e|}{C_1 + C_2 + C_G} \right) V_G - N|e| \frac{C_G V_G + C_2 V_D - N|e|}{C_1 + C_2 + C_G} + \right. \\
 &+ C_2 \frac{(C_1 + C_G) V_D - C_G V_G + N|e|}{C_1 + C_2 + C_G} V_D = \\
 &= \frac{1}{2} \left(\frac{C_1 C_G V_G^2 + C_2 C_G V_G^2 + C_G^2 V_G^2 - C_G^2 V_D^2 - C_G C_2 V_D V_G + N|e| C_G V_G}{C_1 + C_2 + C_G} + \right. \\
 &+ \left. \frac{-N|e| C_G V_G - N|e| C_2 V_D + N^2 e^2 + C_1 C_2 V_D^2 + C_2 C_G V_D^2 - C_G C_2 V_G V_D + N|e| C_G V_D}{C_1 + C_2 + C_G} \right. \\
 &= \left. \frac{C_1 C_G V_G^2 + C_1 C_2 V_D^2 + C_2 C_G (V_G - V_D)^2 + N^2 e^2}{2(C_1 + C_2 + C_G)} \right)
 \end{aligned}$$

$$E_{el}(N) = \frac{N^2 e^2}{2(C_1 + C_2 + C_G)}$$

If I want to add an extra electron:

$$\Delta E_{el}(N) = E_{el}(N+1) - E_{el}(N) = \frac{(N+1)^2 e^2 - N^2 e^2}{2(C_1 + C_2 + C_G)} =$$

$$= \frac{N^2 e^2 + 2N e^2 + e^2 - N^2 e^2}{2(C_1 + C_2 + C_G)} = \frac{e^2}{C_1 + C_2 + C_G} \left(N + \frac{1}{2} \right) = \Delta E_e(N)$$

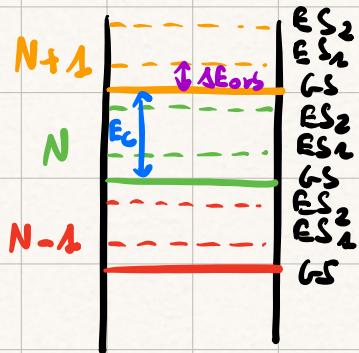
This quantity is called "charging energy", because it tells how much energy we have to spend to add an extra electron to the dot if we have already N electrons inside.

This value corresponds to the "addition energy", in case we are neglecting excited states which arise from quantum confinement:

$$E_{\text{add}} = E_c + \Delta E_{\text{orb}}$$

$$\sim \frac{\hbar^2}{m^* L^2}$$

(particle in a box like)



Notice that in principle there are infinite excited states

Since $E_c > \Delta E_{\text{orb}}$ usually, in all these computations coming from this simple capacitance model we are neglecting effects of quantum confinement ($\Delta E_{\text{orb}} = 0$, meaning that discretization rules are dictated just by coulomb repulsion, since it usually dominates).

$$2) \Delta E_{\text{gen}} = - \delta Q_0 V_0 - \delta Q_2 V_0 = - C_0 \delta (V_0 - V_0) V_0 - C_2 \delta V_2 V_0 \\ = + C_0 \delta V_1 V_0 - C_2 \delta V_2 V_0 = C_0 V_0 - \frac{-(N+1)|e| + N|e|}{C_0 + C_2 + C_0} +$$

$$- C_2 V_0 \frac{(N+1)|e| - N|e|}{C_0 + C_2 + C_0} = - \frac{|e|}{C_0 + C_2 + C_0} (C_0 V_0 + C_2 V_0)$$

$$\rightarrow \Delta E_{\text{tot}} = \frac{1}{C_0 + C_2 + C_0} \left(N|e|^2 + \frac{1}{2}|e|^2 - |e| (C_0 V_0 + C_2 V_0) \right) = \\ = \frac{|e|}{C_0 + C_2 + C_0} \left(\left(N + \frac{1}{2} \right) |e| - C_0 V_0 - C_2 V_0 \right)$$

3) To inject one electron from the source to the dot:

$$\Delta E_{\text{tot}} < 0 \rightarrow (N + \frac{1}{2})|e| - C_a V_D - C_s V_s < 0$$

$$-C_s V_D < -\left(N + \frac{1}{2}\right)|e| + C_a V_s$$

$$\rightarrow V_D > -\frac{C_a}{C_s} V_s + \left(N + \frac{1}{2}\right) \frac{|e|}{C_s} \quad (i)$$

4) Inject an electron from dot to the drain:

$$\Delta E_{\text{tot}} (N \rightarrow N-1) < 0$$

$$\Delta E_c (N \rightarrow N-1) = - \left[\frac{N^2 e^2}{2(C_a + C_s + C_r)} - \frac{(N-1)^2 e^2}{2(C_a + C_s + C_r)} \right] = - \frac{e^2}{C_a + C_s + C_r} \left(N - \frac{1}{2}\right)$$

$$\Delta E_{\text{gen}} (N \rightarrow N-1) = - \delta Q_a V_s + \delta Q_D V_D = - \delta Q_a V_s + \delta(-Q_s + Q_a) V_D =$$

$$= - \cancel{\delta Q_a V_s} - \cancel{\delta Q_s V_D} + \cancel{\delta Q_a V_D} \stackrel{(\delta V_D = 0)}{=} + \cancel{C_a \delta V_s V_s} - \cancel{C_s \delta V_s V_D} +$$

$$- \cancel{C_a \delta V_s V_D} = \delta V_s (C_a V_s - C_s V_D - C_a V_D) =$$

$$= \frac{N|e| - (N-1)|e|}{C_a + C_s + C_r} (C_a V_s - (C_s + C_a) V_D) = \frac{|e|}{C_a + C_s + C_r} (C_a V_s - (C_s + C_a) V_D)$$

$$\Delta E_{\text{tot}} = \frac{|e|}{C_a + C_s + C_r} \left(-\left(N - \frac{1}{2}\right)|e| + C_a V_s - (C_s + C_a) V_D \right)$$

5) $\Delta E_{\text{tot}} < 0 \rightarrow - (C_s + C_a) V_D < \left(N - \frac{1}{2}\right)|e| - C_a V_s$

$$\rightarrow V_D > -\left(N - \frac{1}{2}\right) \frac{|e|}{C_s + C_a} + \frac{C_a}{C_s + C_a} V_s \quad (s)$$

6-7) (i), (s) are eq. of a line in the $V_D - V_G$ plane
(one line for each N value)

However, to get the full picture, we are missing one point: what if $V_D < 0$?

In fact, if $V_D < 0$, it is more energetically favorable to inject one electron from the drain to the dot and then from the dot to the source (the opposite with respect to what we got so far).

If we do the computation, we get two new inequalities:

$$\cdot D \text{ to dot} \Leftrightarrow V_D < \frac{C_0}{C_0 + C_2} V_G - \left(N + \frac{1}{2}\right) \frac{|e|}{C_0 + C_2}$$

$$\cdot \text{dot to } S \Leftrightarrow V_D < -\frac{C_0}{C_2} V_G + \left(N - \frac{1}{2}\right) \frac{e}{C_2}$$

So, in the end, we get 4 inequalities which have to be satisfied to get a current from S to D (for $V_D > 0$) or from D to S (for $V_D < 0$)

$$\cdot V_D > -\frac{C_0}{C_2} V_G + \left(N + \frac{1}{2}\right) \frac{|e|}{C_2} \quad (V_D > 0)$$

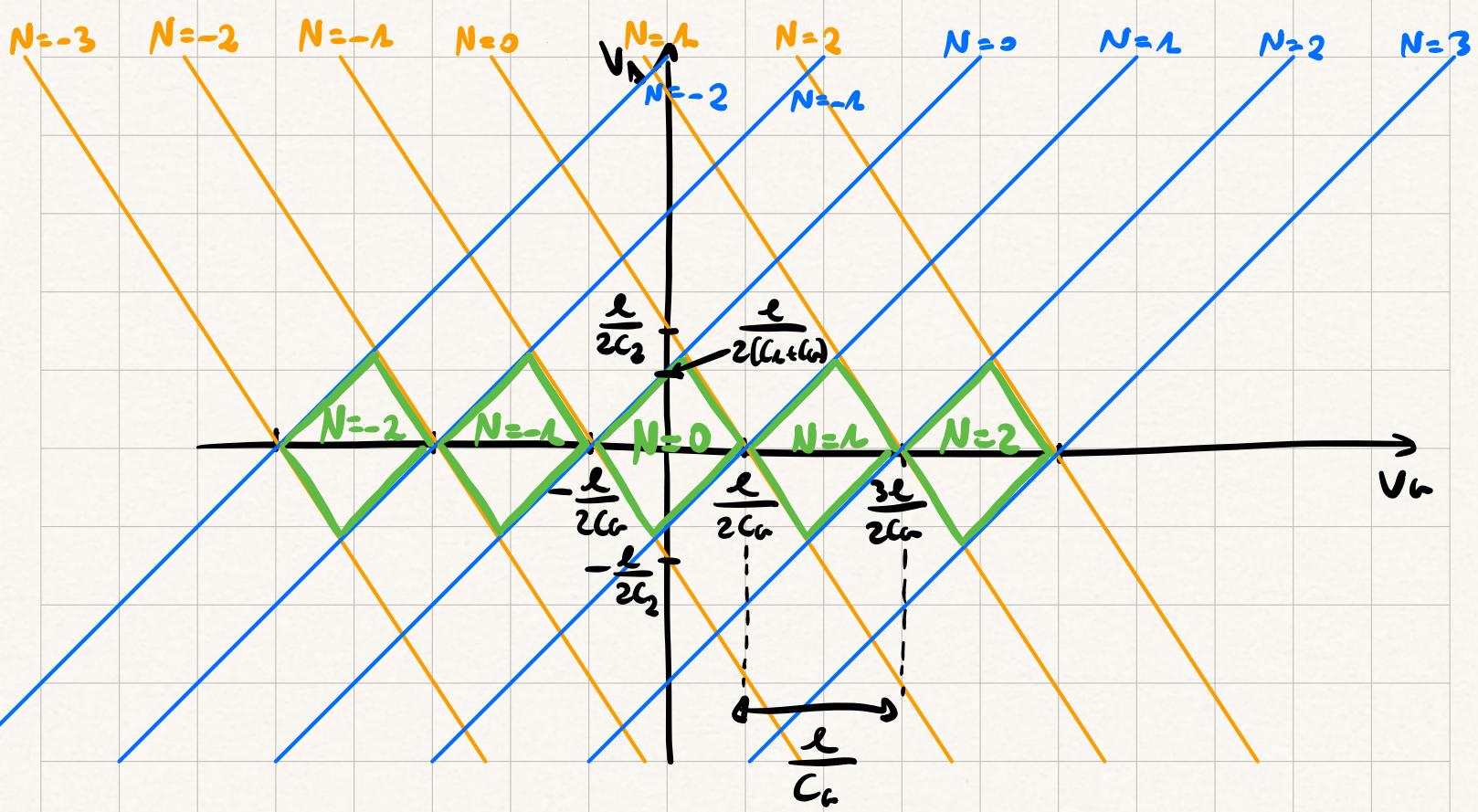
$$\cdot V_D < -\frac{C_0}{C_2} V_G + \left(N - \frac{1}{2}\right) \frac{e}{C_2} \quad (V_D < 0)$$

$$\cdot V_D > \frac{C_0}{C_0 + C_2} V_G - \left(N - \frac{1}{2}\right) \frac{|e|}{C_0 + C_2} \quad (V_D > 0)$$

$$\cdot V_D < \frac{C_0}{C_0 + C_2} V_G - \left(N + \frac{1}{2}\right) \frac{|e|}{C_0 + C_2} \quad (V_D < 0)$$

Source
↓
dot

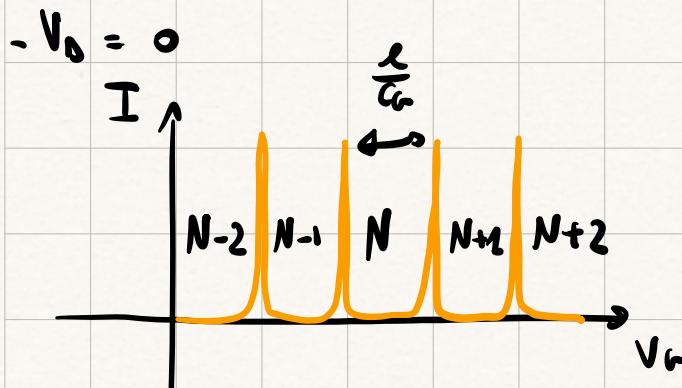
drain
↑
dot



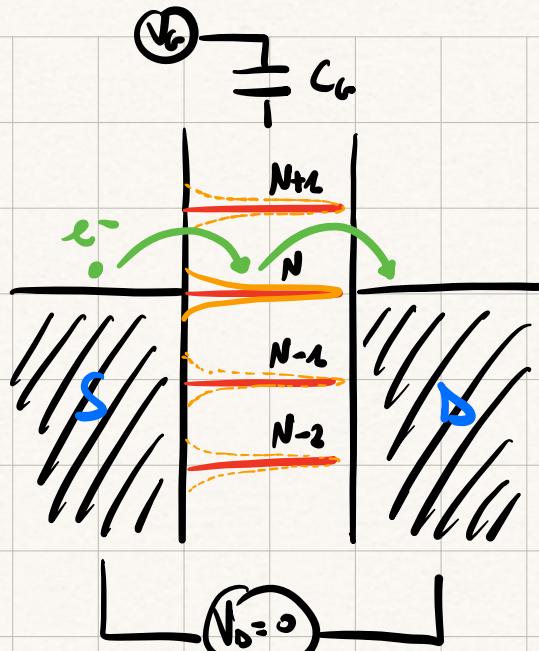
So, if we plot these impurities for each value of N , we get finally the "charge stability diagram" of a single quantum dot. The green regions, called "Coulomb diamonds" because of their shape, are regions where electron transport is completely forbidden by Coulomb repulsion. Outside the diamonds, a current can flow from S to D and vice versa (according to the sign of V_G).

In conclusion, we have to provide energy to electrons in order to overcome Coulomb repulsion. We can do it through V_G or/and V_D .

- Effect of V_G :



Coulomb oscillations

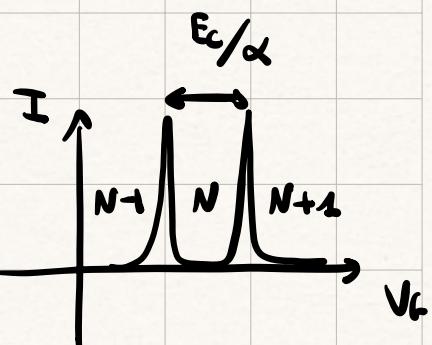


V_G is able to move up and down the chemical potential of the dot. When this potential is aligned to the chemical potentials of the reservoirs, electron tunneling is possible.

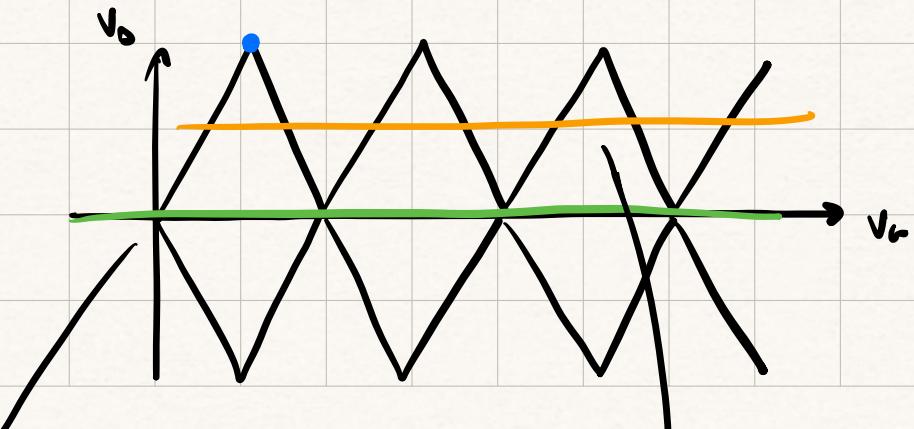
How to convert gate voltage V_G directly into energy provided to the electrons? Through the lever arm α :

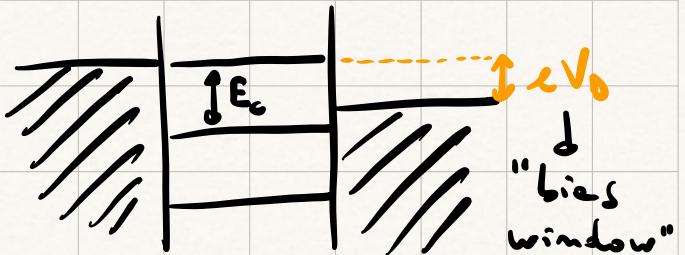
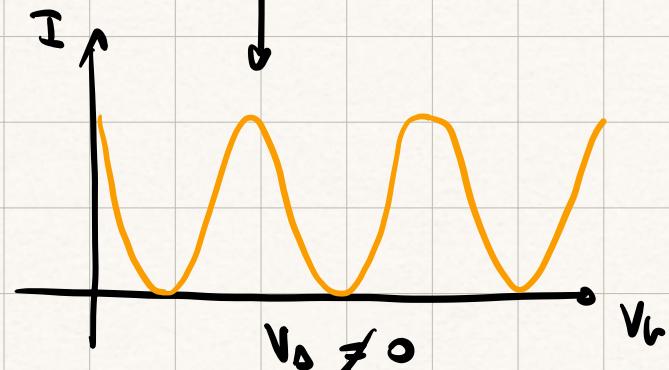
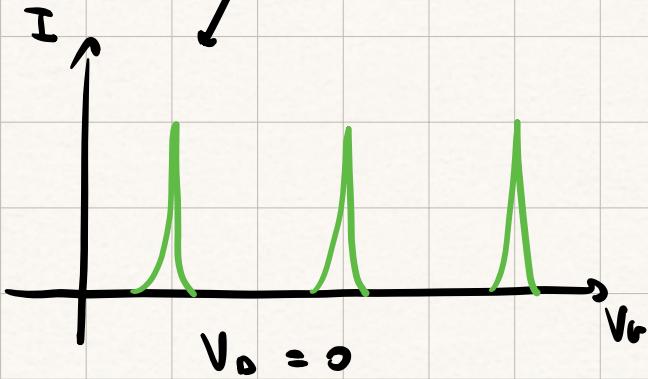
$$\alpha = \frac{C_G}{C_1 + C_2 + C_G} |e| \quad (\text{in } \frac{\text{meV}}{\text{mV}})$$

Through α : $E_C = \alpha \frac{e}{C_G}$ $\xrightarrow{V_G \text{ periodicity}}$



• Effect of V_0

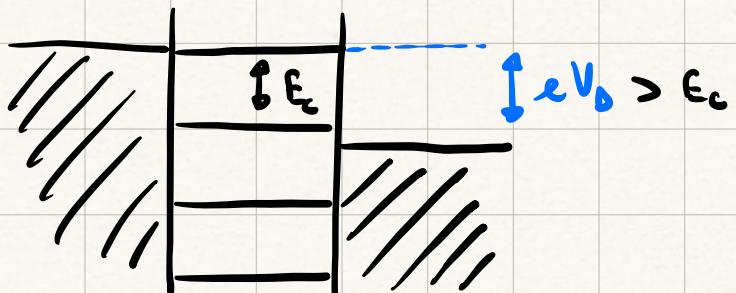




Blue point: no confinement any more

$$\rightarrow E_c = eV_D$$

The lies window is so large that there is always an energy level in the dot within the window, independently on V_G



How much is this value?

From eq. (i) and (j):

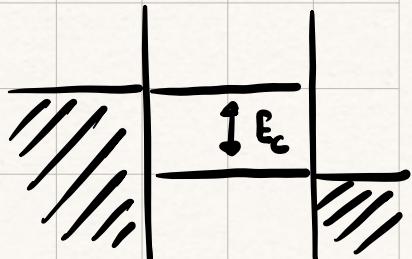
$$\left\{ V_D = -\frac{C_R}{C_L} V_G + \left(N + \frac{1}{2}\right) \frac{|e|}{C_L} \rightarrow V_G = -\frac{C_L}{C_R} V_D + \left(N + \frac{1}{2}\right) \frac{|e|}{C_R} \right.$$

$$\left. V_D = -\left(N - \frac{1}{2}\right) \frac{|e|}{C_L + C_R} + \frac{C_R}{C_L + C_R} V_G \rightarrow V_G = \frac{C_L + C_R}{C_R} V_D + \left(N - \frac{1}{2}\right) \frac{|e|}{C_R} \right.$$

$$- - \frac{C_2}{C_0} V_0 + \left(N + \frac{1}{2}\right) \frac{1}{C_0} = \frac{C_2 + C_0}{C_0} V_0 + \left(N - \frac{1}{2}\right) \frac{1}{C_0}$$

$$(C_2 + C_0) V_0 = 1 \rightarrow V_0 = \frac{1}{C_2 + C_0}$$

$$\rightarrow E_C = \epsilon V_0 = \frac{\epsilon^2}{C_2 + C_0}$$

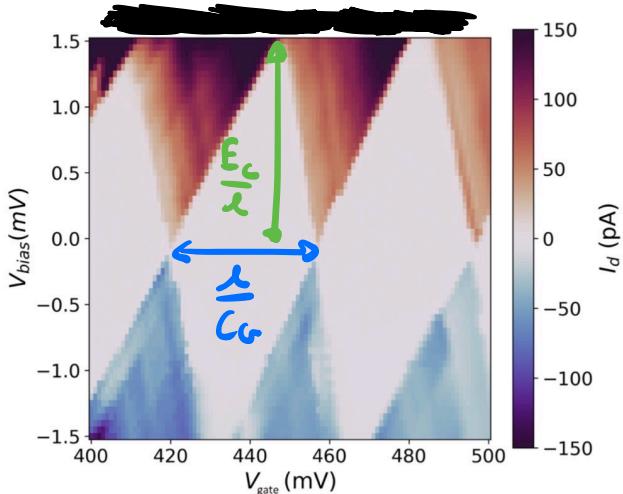


Notice that this value of E_C is different from the expression we found previously in the exercise, which was N -dependent ($\Delta E_C(N) = \frac{\epsilon^2}{C_2 + C_0} \left(N + \frac{1}{2}\right)$)

This is because it is more convenient to re-define an expression which is N -independent, in order to think in terms of energy differences and not absolute values defined starting from an arbitrary reference (ex. $N=0$)

$$E_C = \Delta E_C(N + \epsilon) - \Delta E_C(N) = \frac{\epsilon^2}{C_2 + C_0 + C_{dot}} - \frac{\epsilon^2}{C_{dot}}$$

8)



$$E_C \approx 1.5 \text{ meV}$$

$$\Delta V_G = \frac{\epsilon}{C_0} \approx 36 \text{ mV}$$

$$\alpha = \epsilon \frac{C_0}{C_{dot}} = \frac{\epsilon^2 / \Delta V_G}{\epsilon^2 / E_C} =$$

$$= \frac{E_C}{\Delta V_G} \quad (\rightarrow \Delta V_G = E_C / \alpha)$$

$\rightarrow \alpha \approx 0.04 \frac{\text{meV}}{\text{mV}}$ \rightarrow if we apply $V_R = 1 \text{ mV}$, we shift the chemical potential in the dot of $40 \mu\text{V}$

Notice that now that we know E_c , we can also compute the total capacitance of the dot:

$$E_c = \frac{e^2}{C_{\text{dot}}} \rightarrow C_{\text{dot}} = \frac{e^2}{E_c}$$

And by approx. the dot as a conductive disk, we can estimate the dot radius:

$$C_{\text{dot}} = 8\pi\epsilon_0\epsilon_r R_{\text{dot}}$$

Notice: in the picture you can also see other transition lines outside the diamond, which lead to an increase in current when matching this transition by V_R and V_S \rightarrow excited states due to quantum confinement, as anticipated at the beginning $\rightarrow \Delta E_{\text{orb}}$ can be extracted from the graph as well

Max T?

$$E_c > k_B T \rightarrow T < \frac{E_c}{k_B} = \frac{1.5 \text{ meV}}{8.6 \cdot 10^{-5} \text{ eV/K}} = 17 \text{ K}$$