

EPFL



MICRO-435 Quantum and Nanocomputing

Edoardo Charbon
Mariagrazia Graziano

MIT BEHAVIOR & CHARACTERISTICS
PART 3

OBJECTIVE

- a) CHARGING EFFECT & RESONANCE
- b) EFFECT OF GATE

b) CHARGING EFFECT & RESONANCE

ELECTROSTATIC EFFECT

APPLY V_{DS}

→ CREATE BC

→ VARIATION CHANNEL
CHARGE

↳ E_L, H, L

⇓

I_{DS}



CHARGING EFFECT

WHEN A CHARGE ENTERS IN THE CHANNEL

THE ENERGY LEVELS IN THE CHANNEL ARE

ALTERED

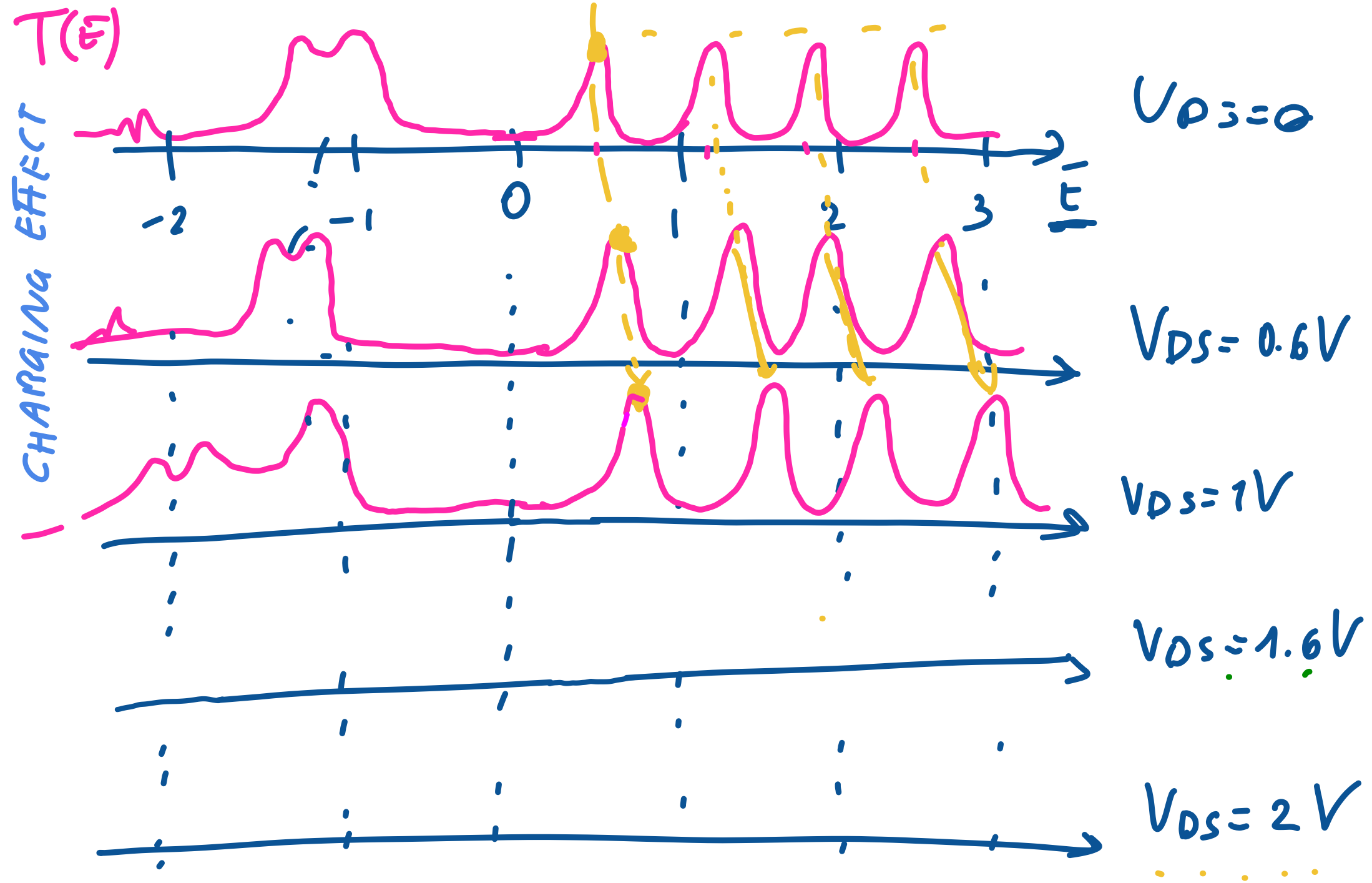
CHARGING EFFECT

WHEN A CHARGE ENTERS IN THE CHANNEL
THE ENERGY LEVELS IN THE CHANNEL ARE
ALTERED

$V_{DS} = 0 \rightarrow T(E) @ \text{EQUILIBRIUM}$
DUE TO N_0 \nearrow N_0 N° of el. IN CHANNEL
@ EQUILIBRIUM

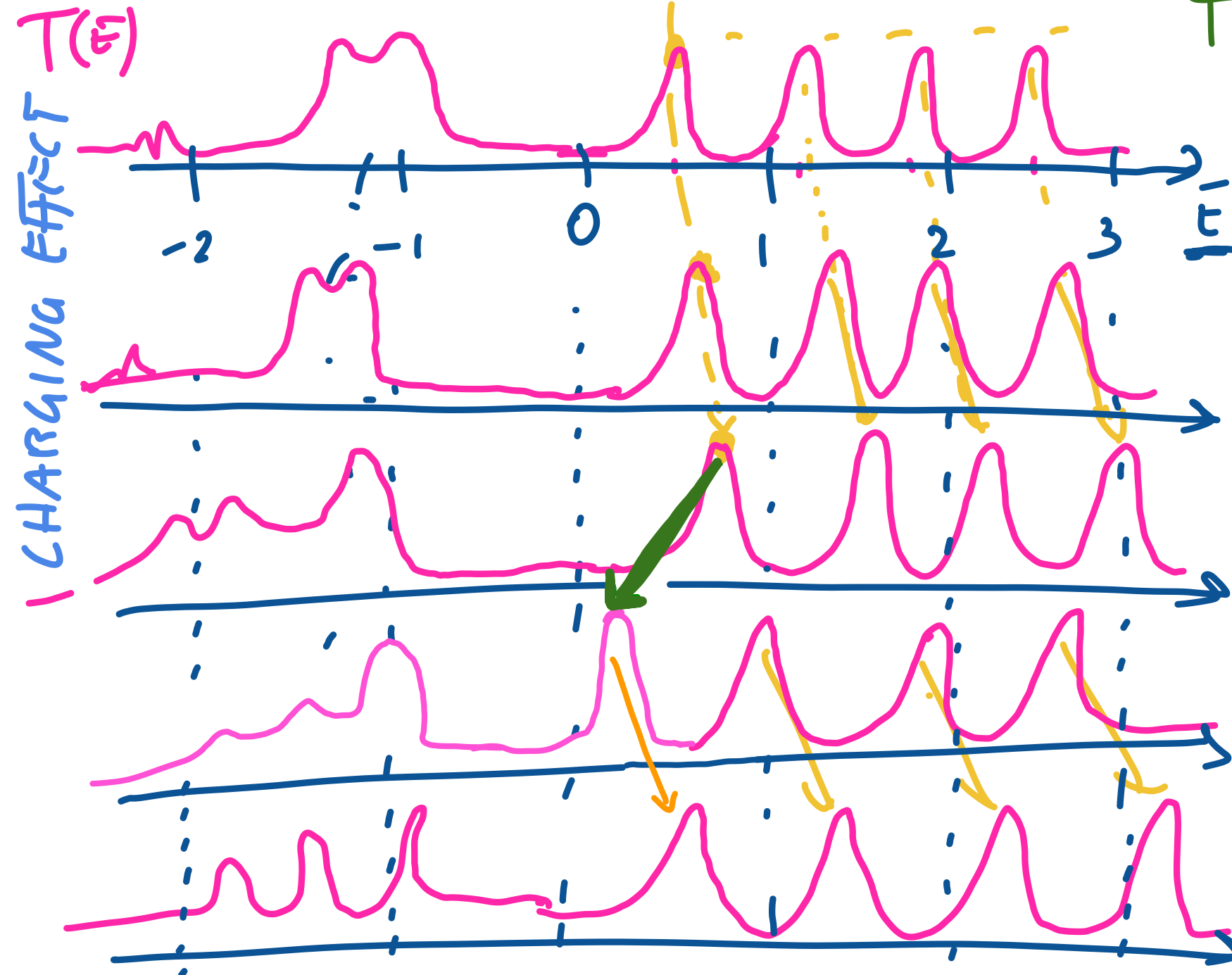
if $V_{DS} \uparrow$ or \downarrow
 $N_0 \rightarrow N \neq N_0$ N° of el. CHANGES
 $T(E) \rightarrow T(E, \mu)$ T. S. CHANGES
 $I_{DS} \rightarrow I_{DS}(T(E, \mu))$ CURRENT CHANGES

EXAMPLE : TIOPHENE BRINGS



EXAMPLE : TIOPHENE BRINGS

WHAT IS THIS CHANGE?



$V_{DS}=0$

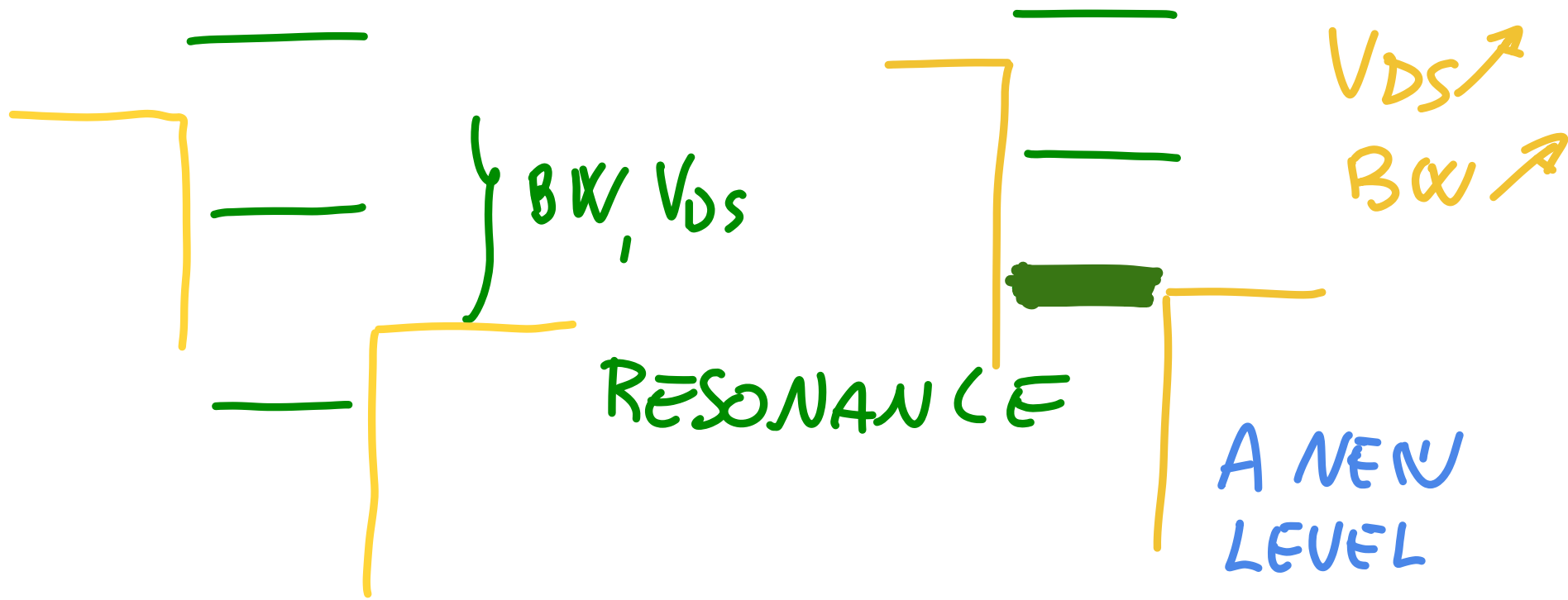
$V_{DS}=0.6V$

$V_{DS}=1V$

$V_{DS}=1.6V$

$V_{DS}=2V$

...



A NEW LEVEL ENTERS IN THE B.W.

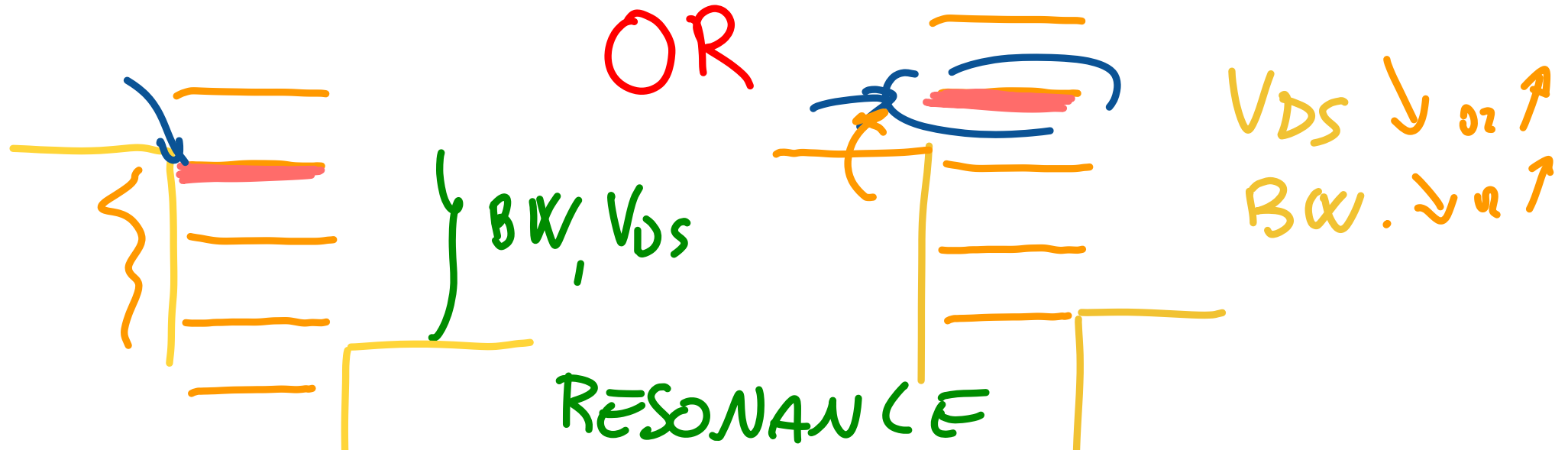


A NEW ELECTRON JUMPS IN THE CHANNEL

NEGATIVE CHARGES ENTERS



THE T_s SHIFTS TOWARD LOWER ENERGIES



if a NEW LEVEL OF ENERGY
EXITS FROM THE CHANNEL

↓
 1eL jumps out

↓
 NEW POSITIVE CHARGE

TS SHIFT TOWARD HIGHER ENERGIES

RECAP

I_{OS}



ELECTROSTATIC
CONTRIBUTION

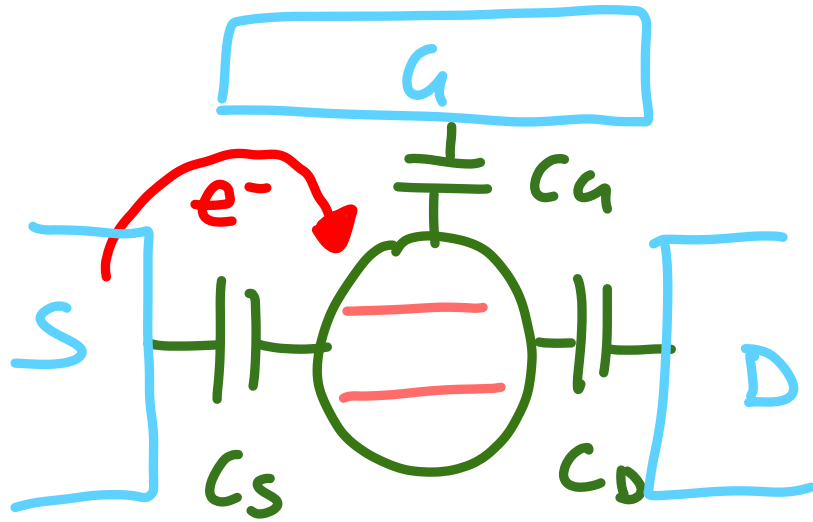


CHARGING
EFFECT CONTRIBUTION

(w/o or w RESONANCE)

HOW IS THIS TAKEN INTO ACCOUNT
IN THE MODEL ?

CAPACITIVE MODEL OF QDOT



$$\underline{C_{ES}} = C_S + C_D + C_G$$

$$U_{\text{dot}} = -q \underset{\uparrow}{V_{DS}} \frac{C_D}{C_{ES}} - q \underset{\uparrow}{V_{GS}} \cdot \frac{C_G}{C_{ES}} + \frac{q^2}{E_S} \cdot (N - N_0)$$

ACTUAL N : NO OF e^- IN THE DOT
 NO OF e^- @ Eq.

$$\underline{N(V_{GS}, V_{DS})} = \sum_{i=1}^M \int \frac{L}{\gamma_{i,1} + \gamma_{i,2}} \cdot D_{E_L}(E - U_{DS}) \cdot [\gamma_{i,1} f(\bar{E}_L, \bar{E}_{TS}) + \gamma_{i,2}(\bar{E}_L, \bar{E}_{FO})] d\bar{E}$$

TO CORRECTLY EVALUATE N

SCF → SELF CONSISTENT FIELD

→ SEE NEXT SECTION OF THE LESSON

NOW LET'S FOCUS ON KEY POINTS

KEY POINTS

V_{DS}

$$\rightarrow \Delta Q_m = -q \Delta N$$

VARIATION
OF CHARGE
IN CHANNEL

$$\Delta V_{\text{CHARGE}} = \frac{-q \Delta N}{C_{ES}}$$

→ MODEL
WITH
CAP.

$$U_{\text{CHARGE}} = -q \Delta V_{\text{CHARGE}} = \frac{q^2}{C_{ES}} \Delta N$$

→ V_{DS} → ΔN → $U \propto (\Delta N)$ → $D \propto (\epsilon, U)$

→ $T(E, U) \propto D$ $I_{DS} \propto T(E, U)$

c) EFFECT OF GATE

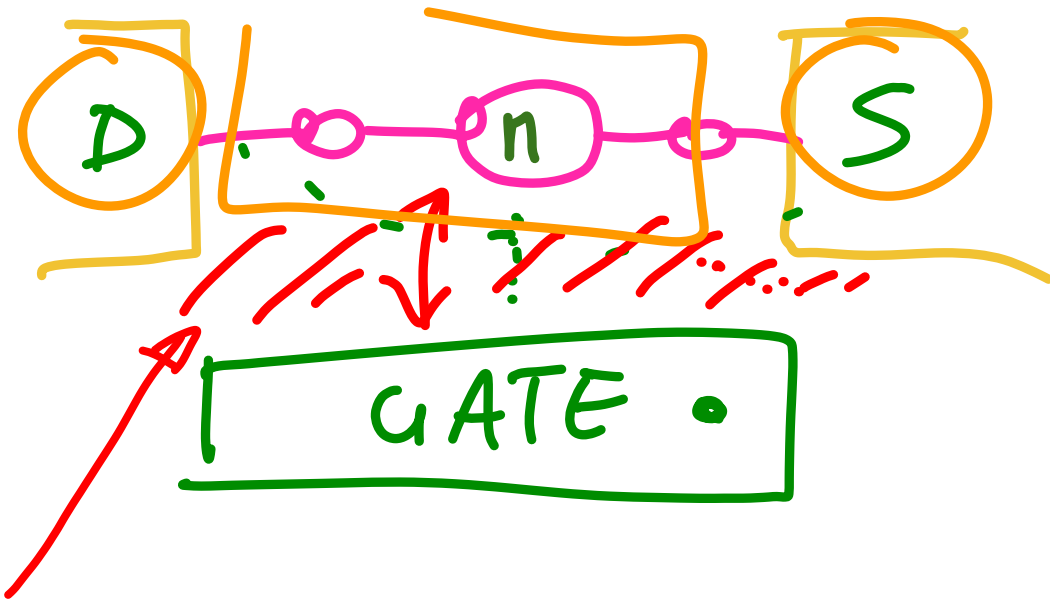
- GENERAL

- BENZENE

- OPV7

- PCP

EFFECT OF GATE



ACERTAIN MATERIAL
AND DISTANCE DETERMINES
THE COUPLING FACTOR

V_{GS} APPLIED

↓ COUPLING FACTOR

D - M - S

SYSTEM CHANGES

POTENTIAL ENERGY

U

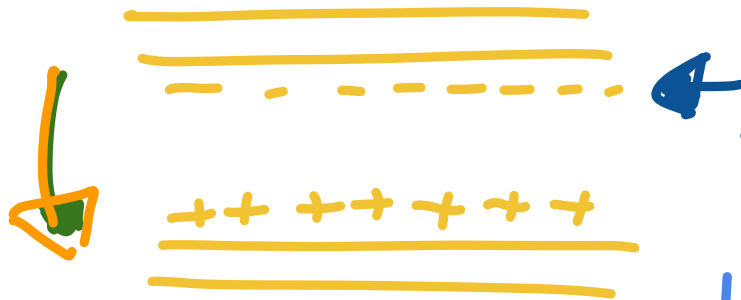
↓

$T(E)$ CHANGES

THE STRUCTURE CAN BE SEEN AS A CAPACITOR

S-M-D

q

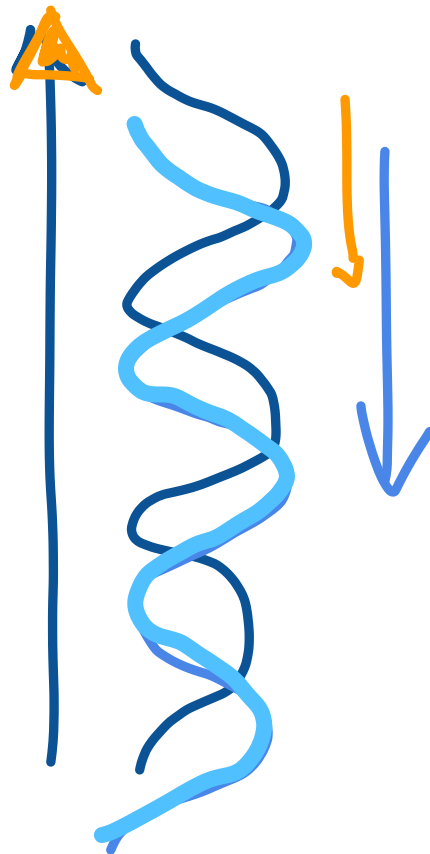


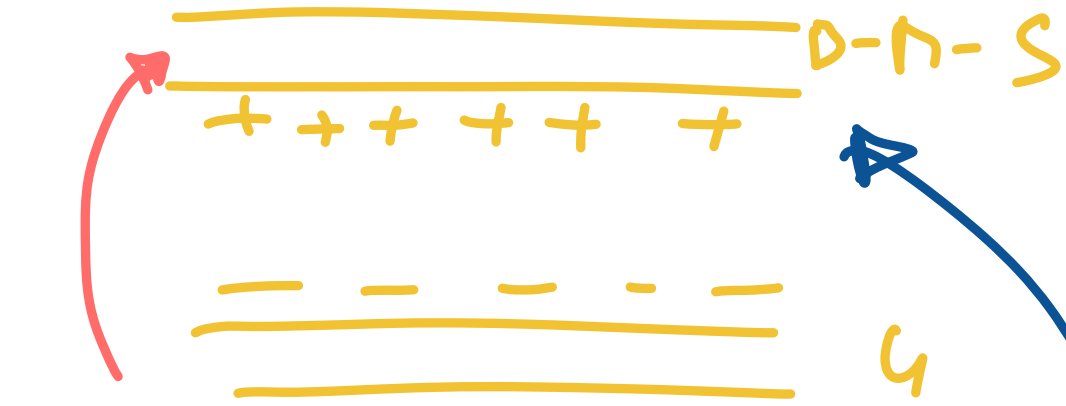
$T_s(E)$

$V_{GS} > 0$

SHIFTING
LOWER
ENERGY

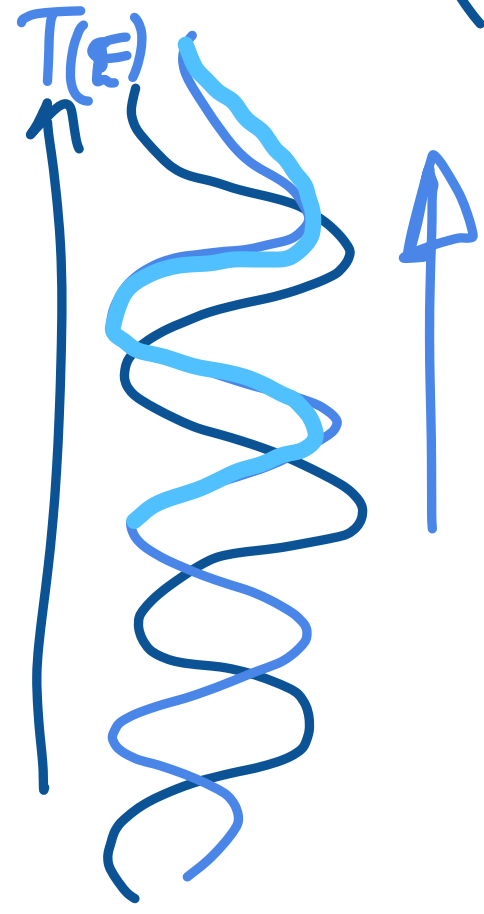
T_s

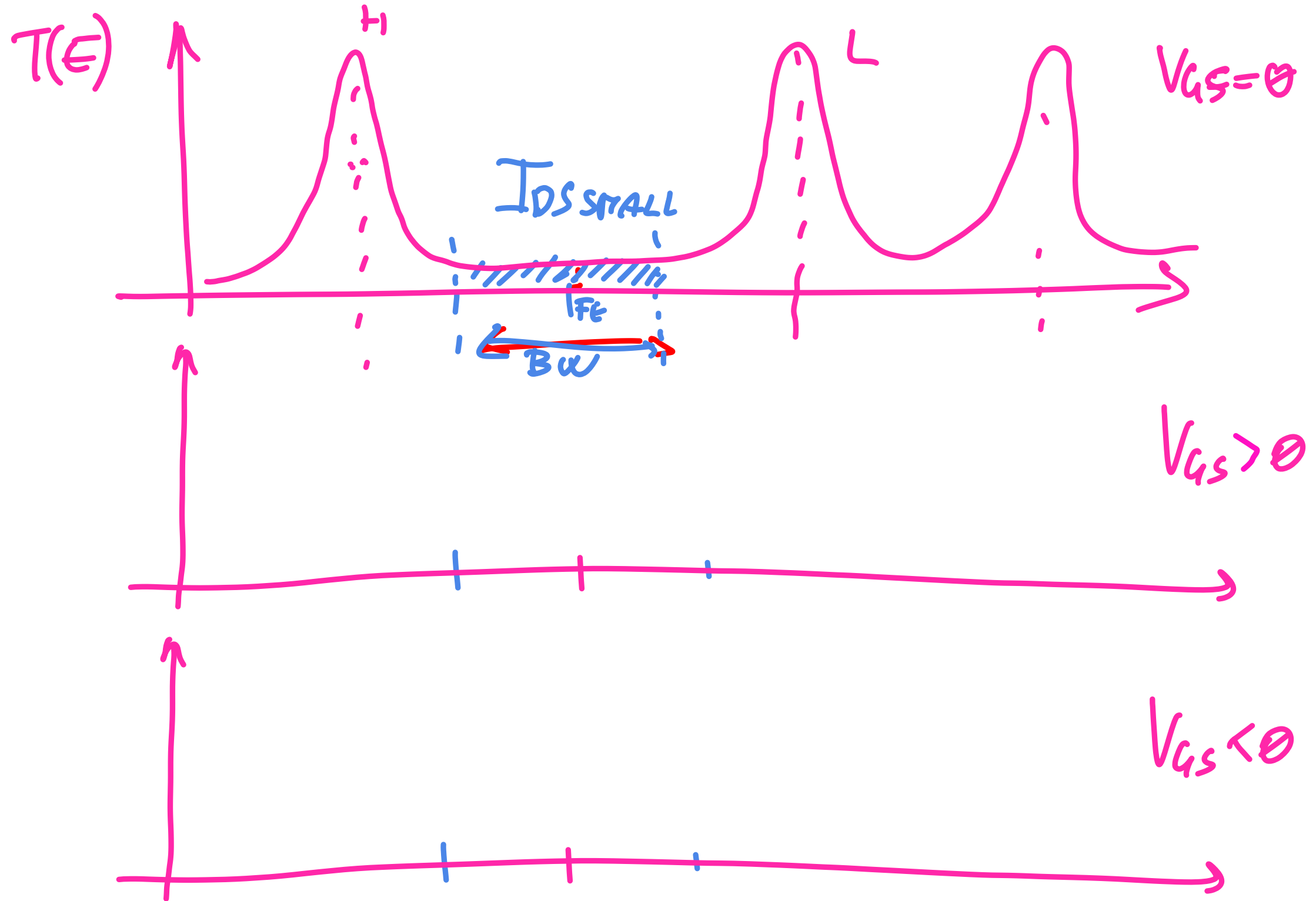


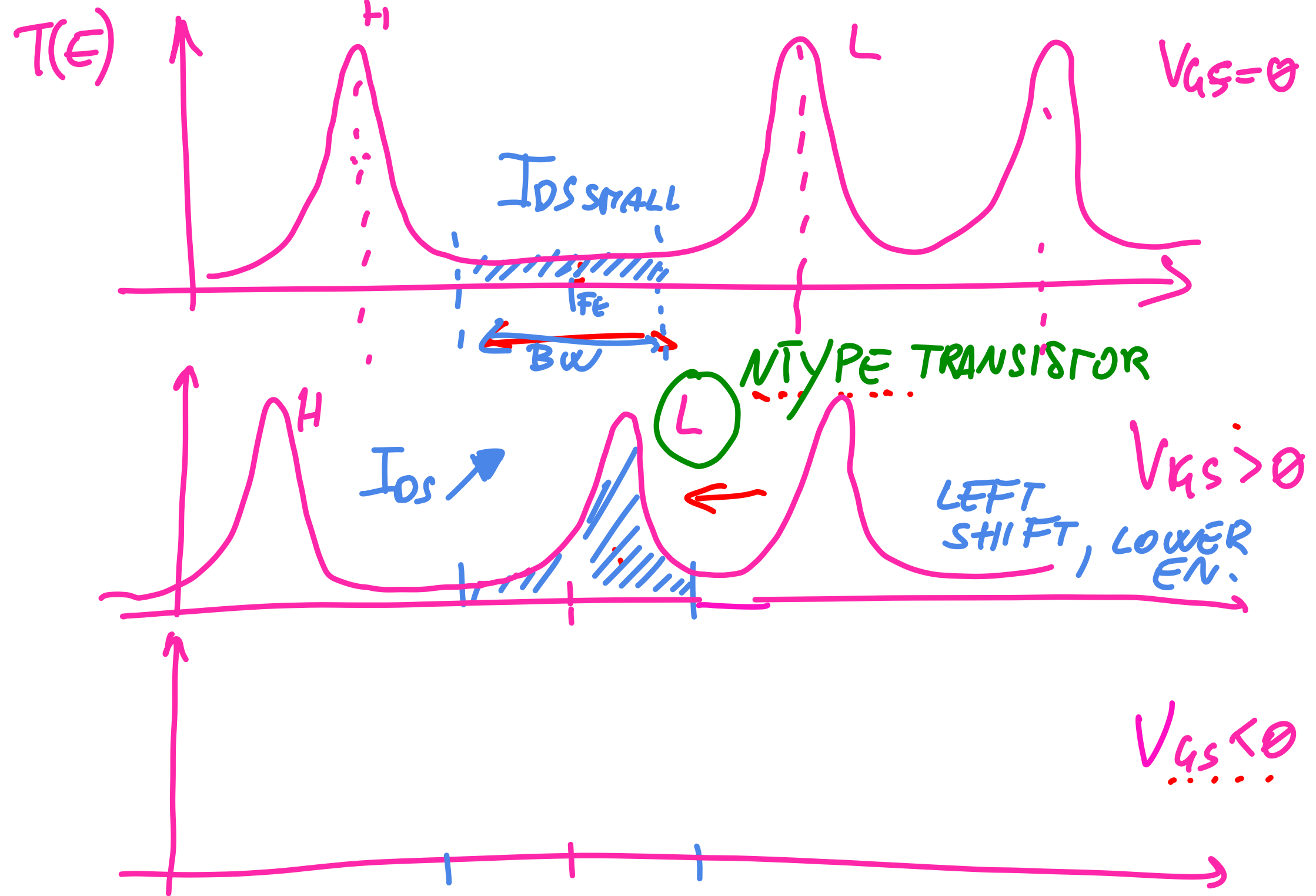


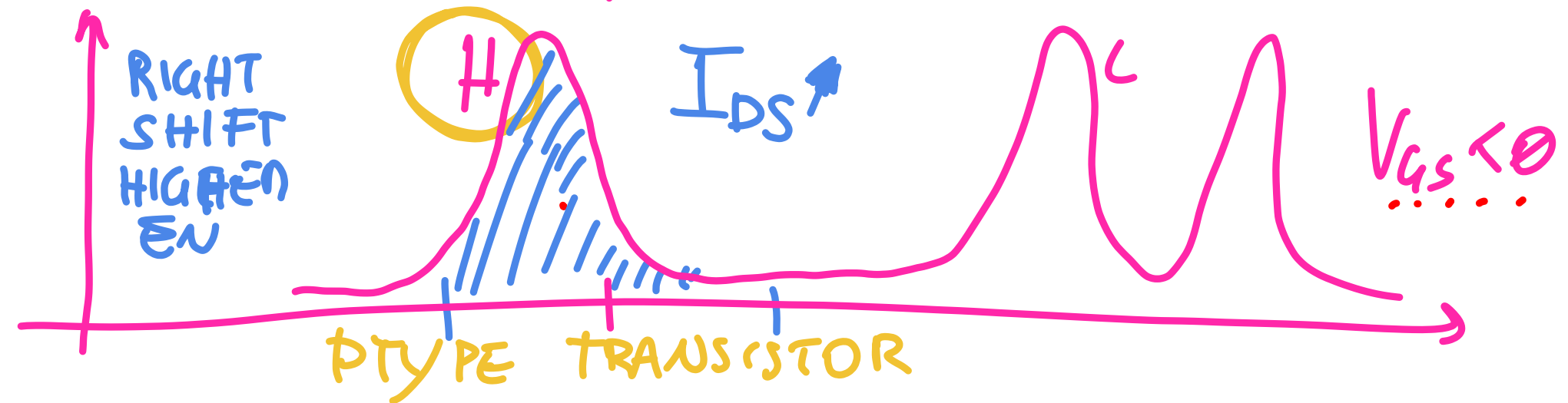
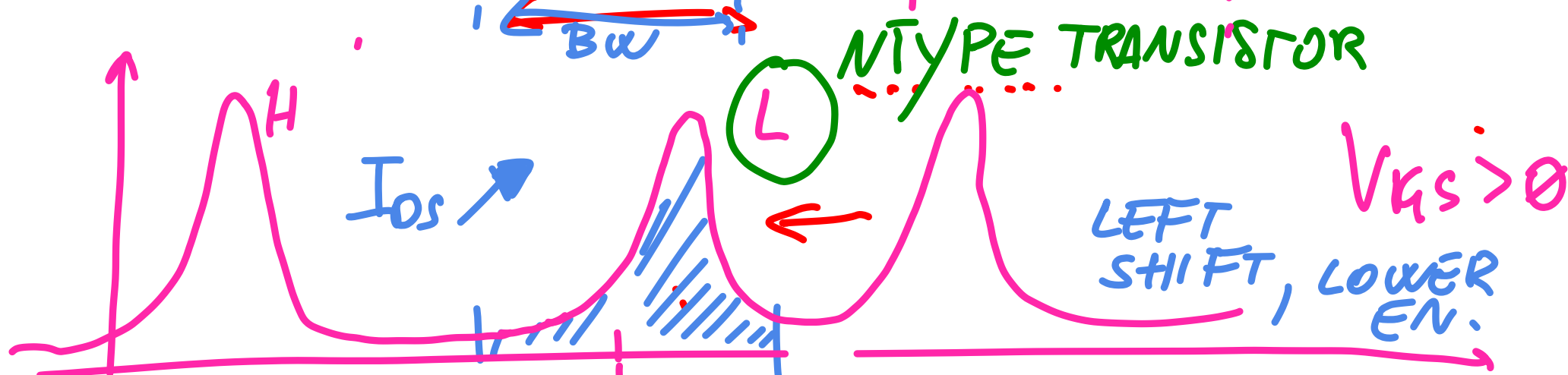
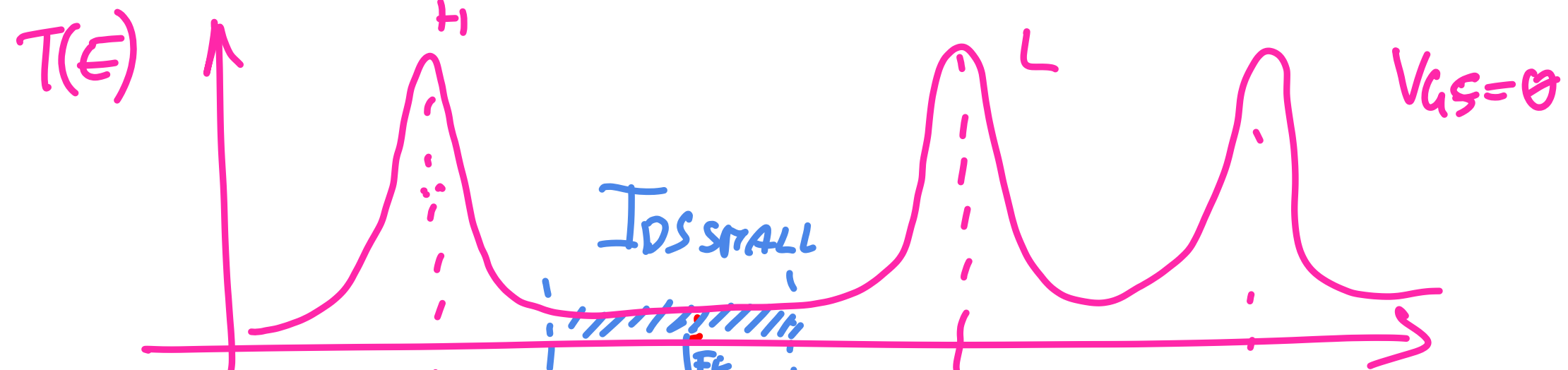
$$V_{GS} < 0$$

POSITIVE CHARGE
 $T_S(E)$ SHIFT
 TOWARD
 HIGHER
 ENERGIES

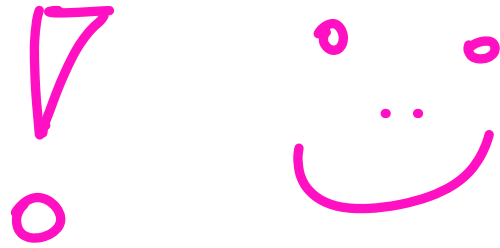








WE CAN USE V_{GS} TO
SWITCH ON/OFF THE TRANSISTOR !



BUT WE ARE ACTUALLY DOING
SOMETHING MORE... WHAT ?

WE ARE USING V_{GS}

→ TO CONTROL

ON/OFF SWITCH

→ TO CONTROL

N TYPE / P TYPE

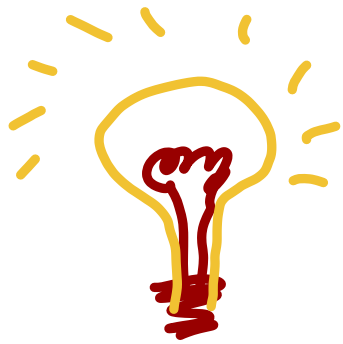
WE ARE USING V_{GS}

→ TO CONTROL

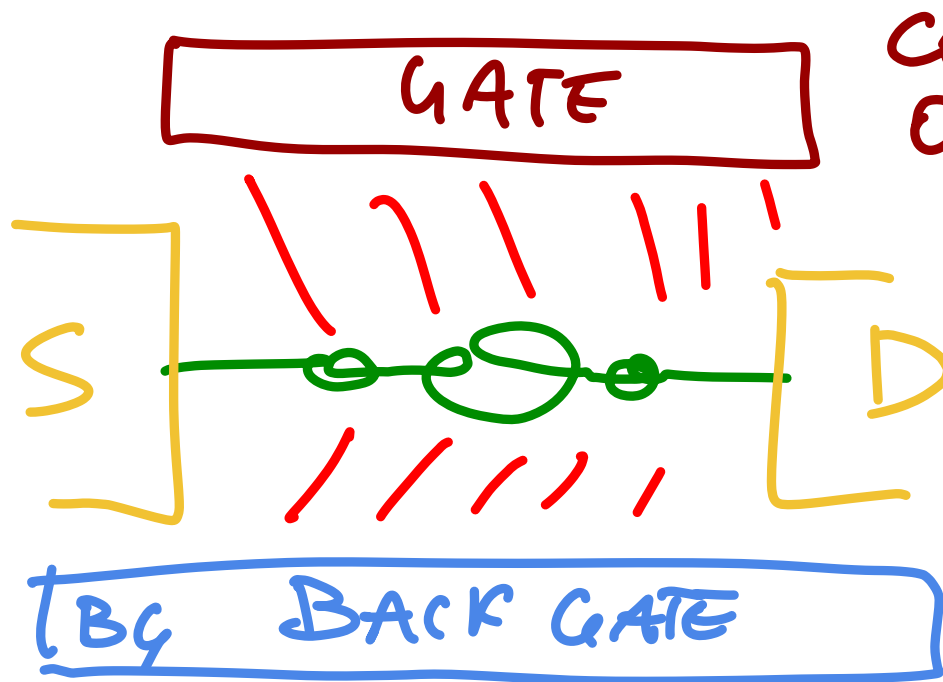
ON/OFF SWITCH

→ TO

CONTROL NTYPE / PTYPE



COMBINE
GATE, BG

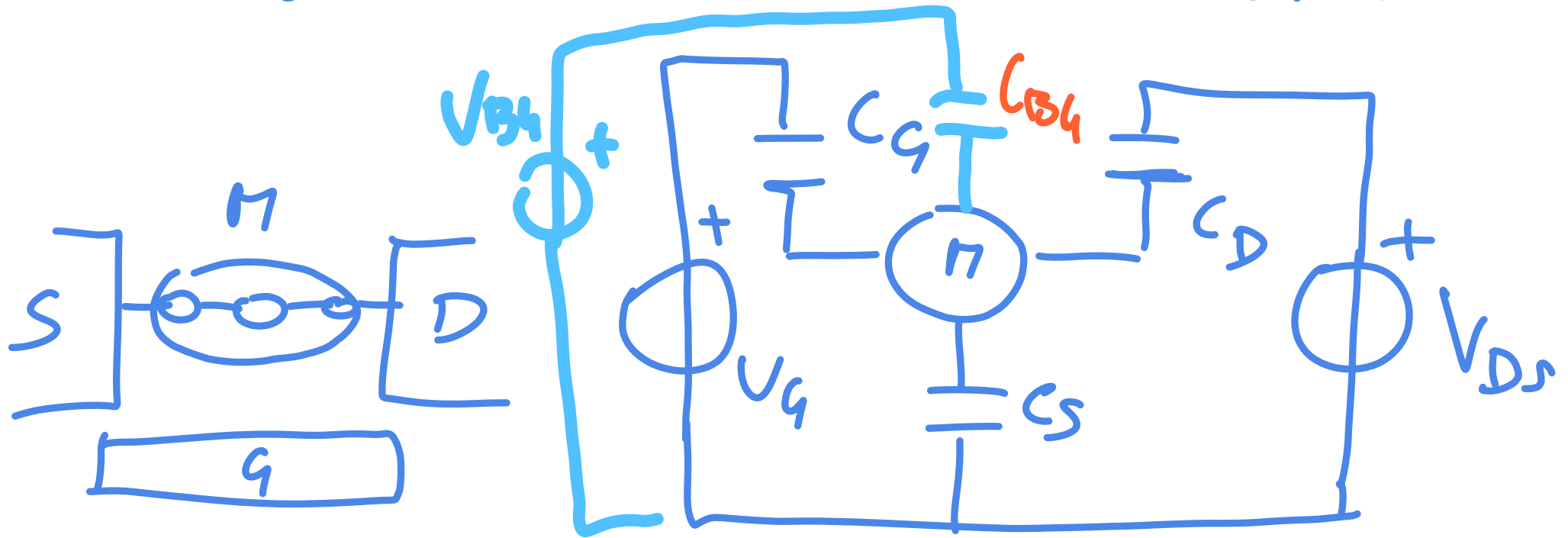


CONTROL
ON/OFF

CONTROLS
NTYPE
OR
PTYPE

WHAT IN THE MODEL ?

CAPACITIVE MODEL OF Q. DOT



$$C_{ES} = C_G + C_S + C_D + C_{BG}$$

$$U = U_{V_{DS}} + U_{C.E} + U_{V_{GS}} + U_{V_{BG}}$$

$$U \left\{ \begin{array}{l} U_{VDS} = -q \frac{C_D}{C_{ES}} V_{DS} \\ U_{VGS} = -q \frac{C_G}{C_{ES}} V_{GS} \\ U_{VBG} = -q \frac{C_G}{C_{ES}} V_{BG} \\ U_{C.E.} = \frac{q^2}{C_{ES}} \cdot \Delta N \end{array} \right.$$

$V_{BG} > 0$
 $V_{GS} > 0$
 $V_{GS} < 0$
 $V_{BG} < 0$

U SHIFTS LOWER ENERGY
 U SHIFTS HIGHER ENERGY

$$\Delta N = N - N_0 \leftarrow \begin{array}{l} \text{No el. } \odot \\ \text{EQ.} \end{array}$$

U_{SCF}

SELF CONSISTING
FIELD

NEXT
SECTION
FOR MODELING

THE GATE & BG

$$V_{Vas} = -q \left(\frac{C_G}{C_{ES}} \right) V_{GS}$$

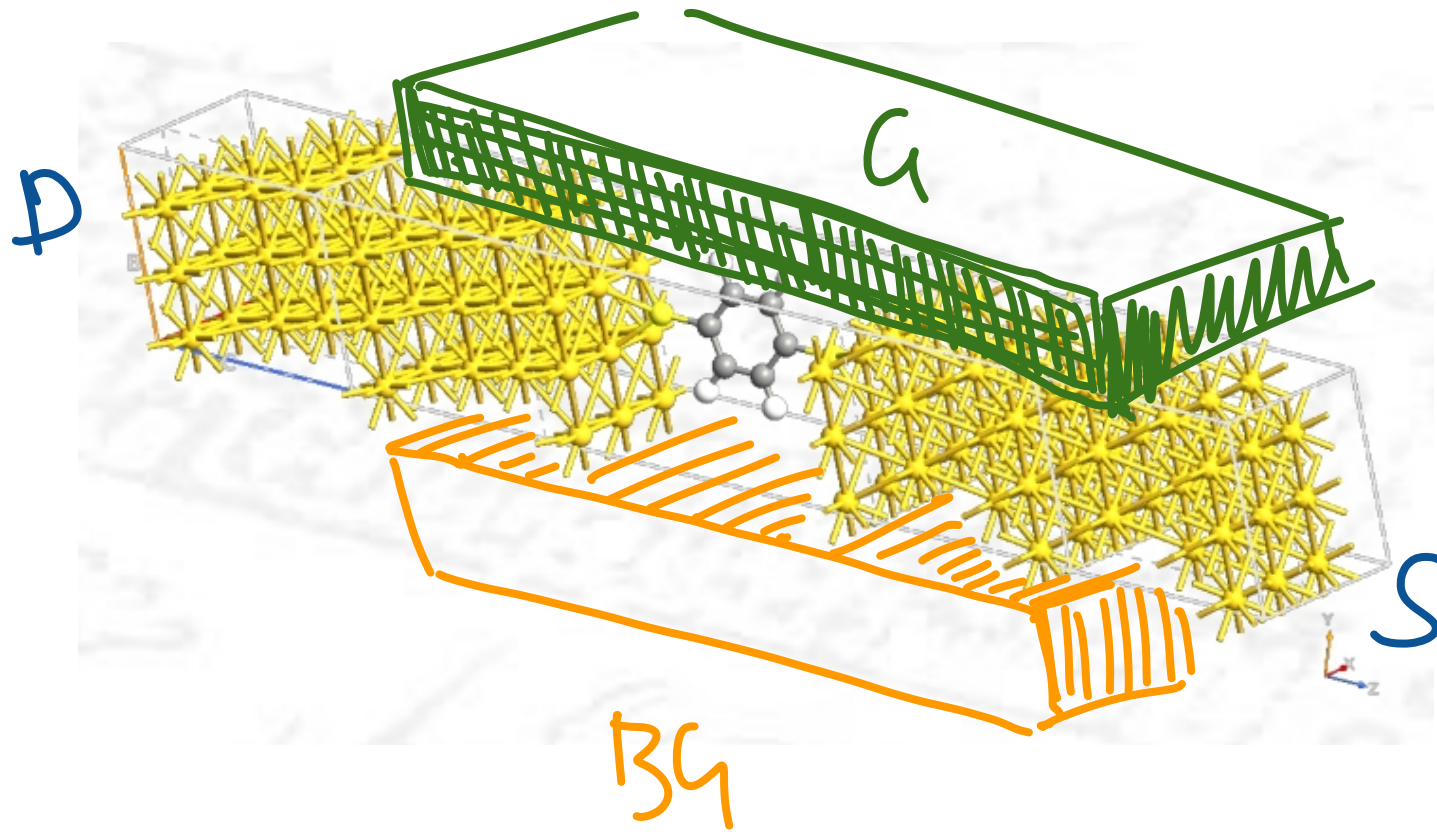
$$V_{V_{BG}} = -q \left(\frac{C_{BG}}{C_{ES}} \right) V_{BG}$$

α
COUPLING
FACTOR

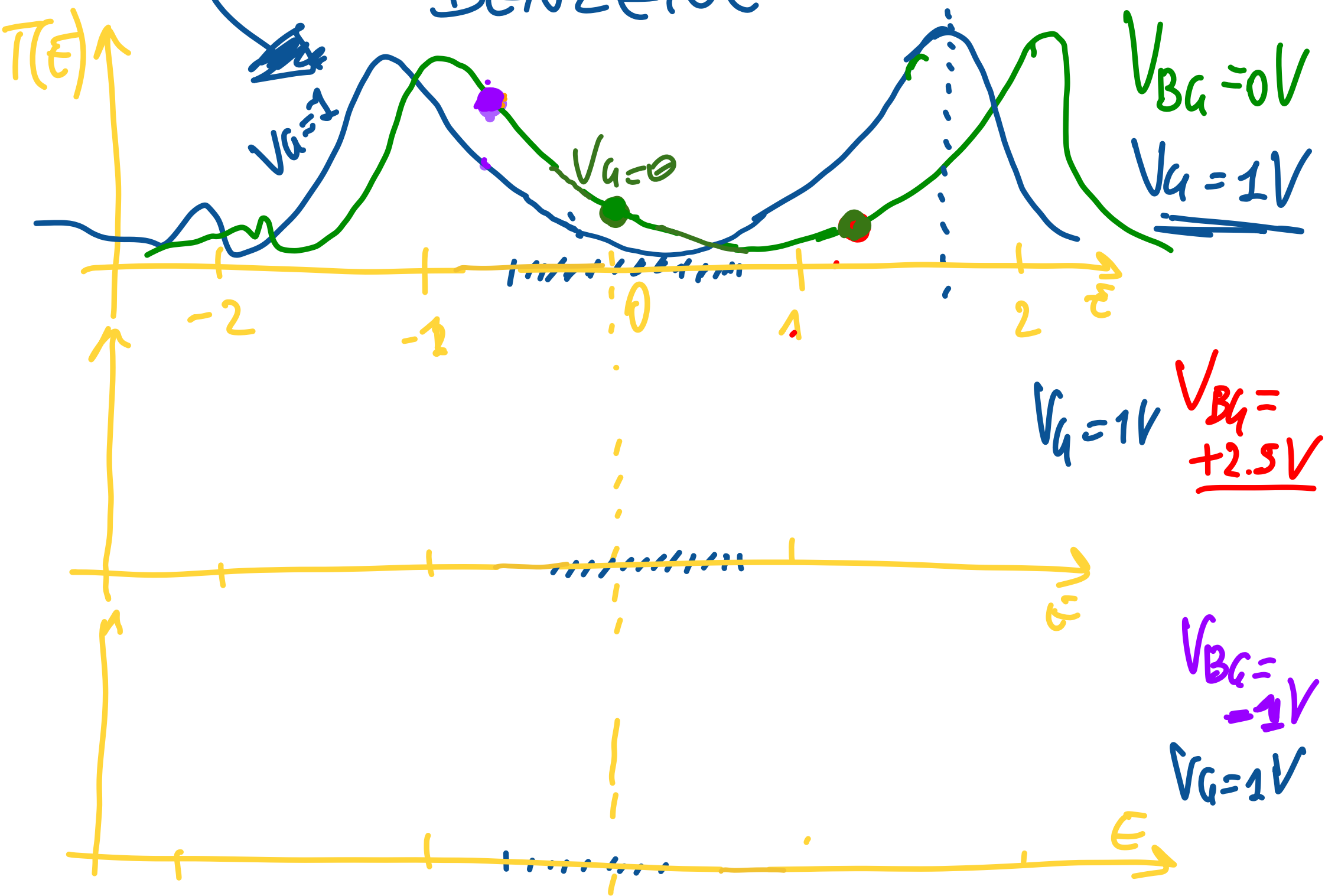
β
BG
COUPLING
FACTOR

NOTE 1. in the example @ THEORY
 $\alpha = 0.5$ but this depends on GATE
ELECTRODE distance and COUPLING MEDIUM
NOTE 2. α and β could be different

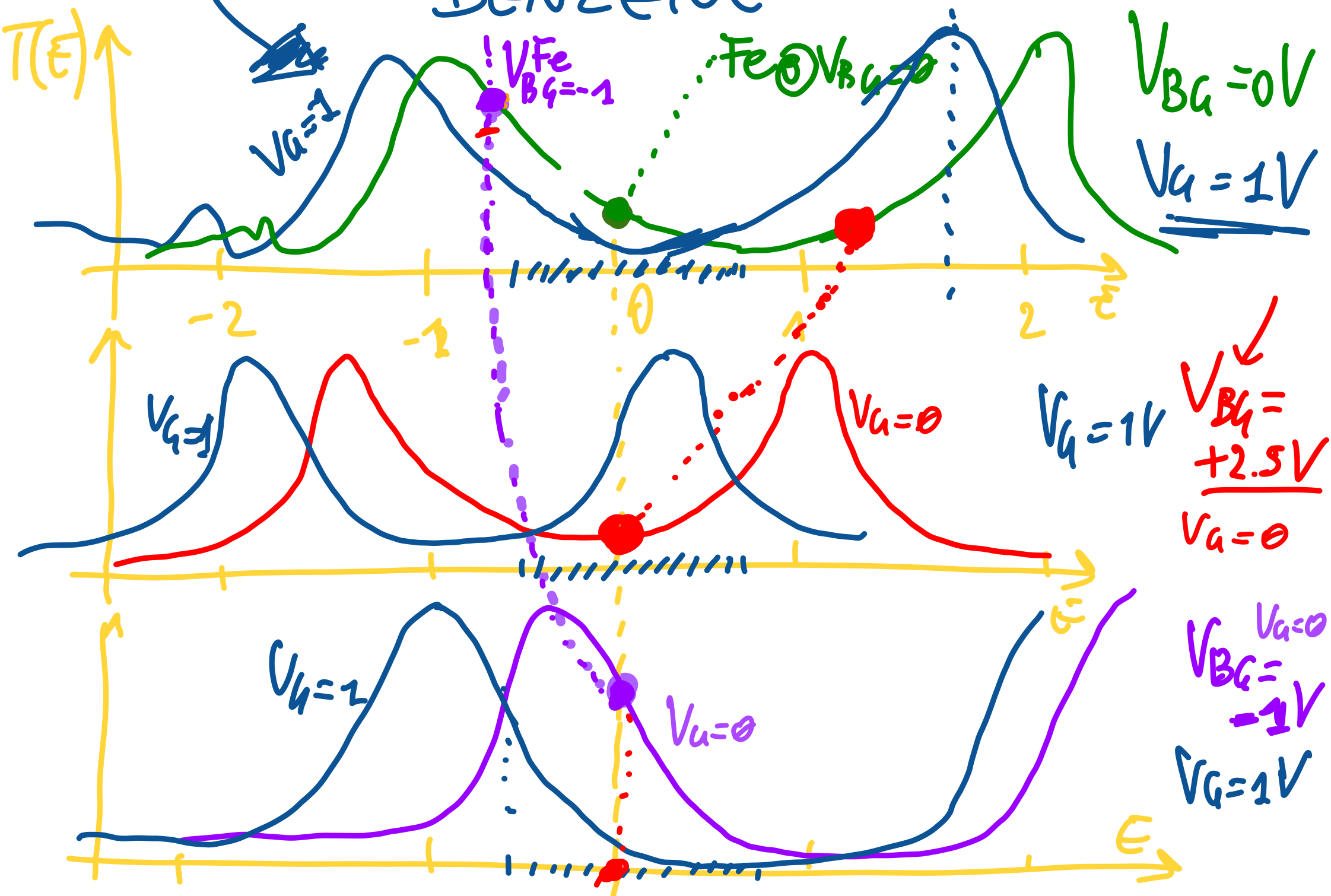
BENZENE

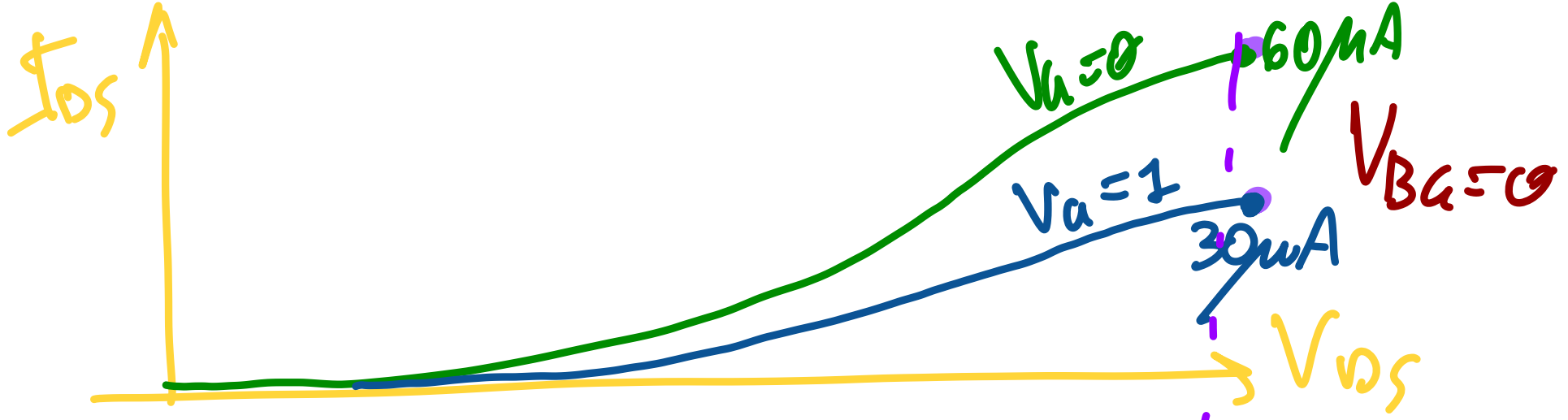


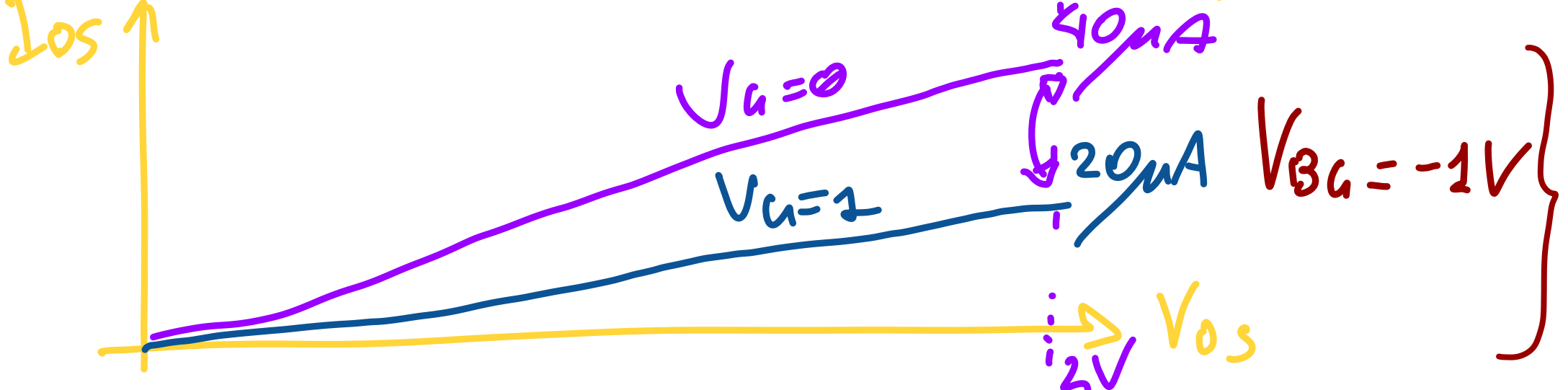
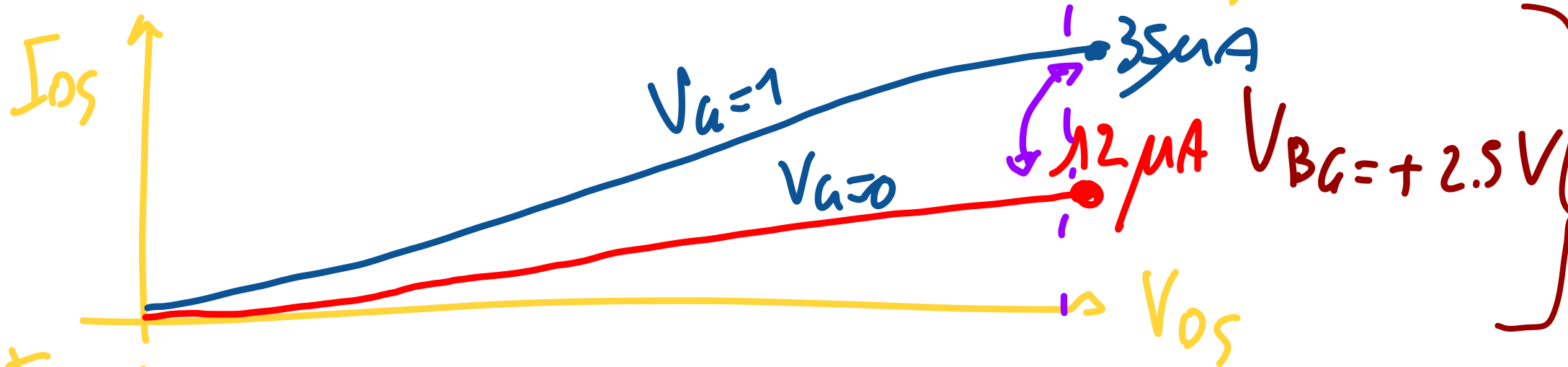
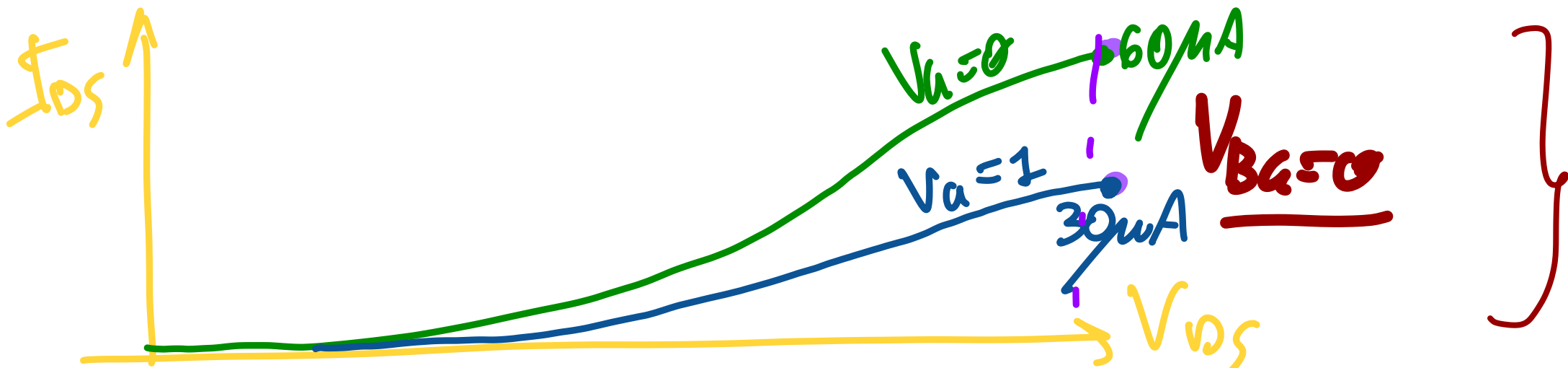
BENZENE



BENZENE





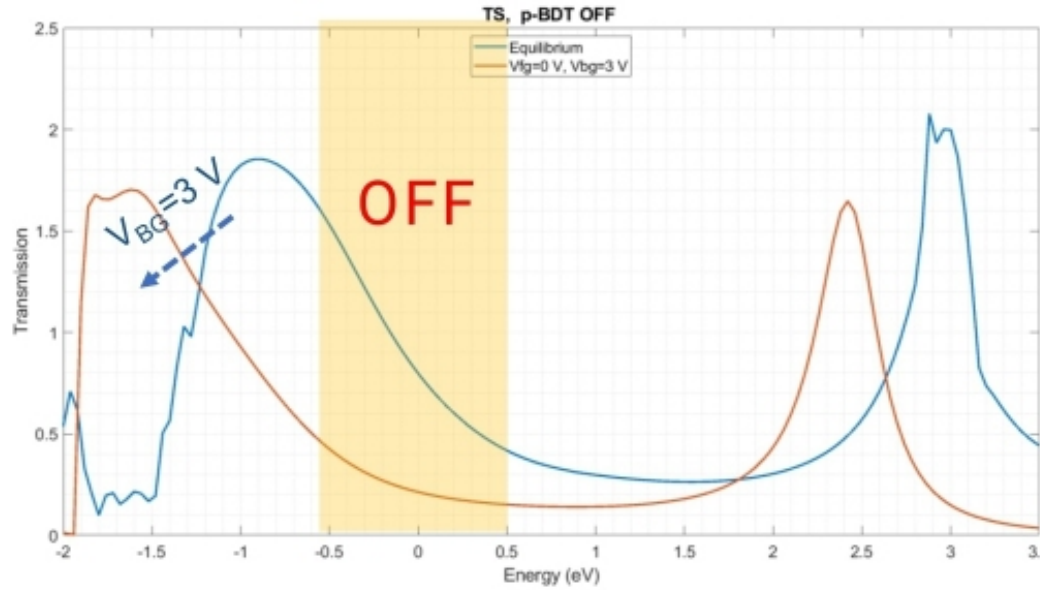


BENZENE, ANOTHER POSSIBLE EXAMPLE OF V_{BG}

V_{FG} = FRONT GATE, I.P. GATE FOR ON/OFF

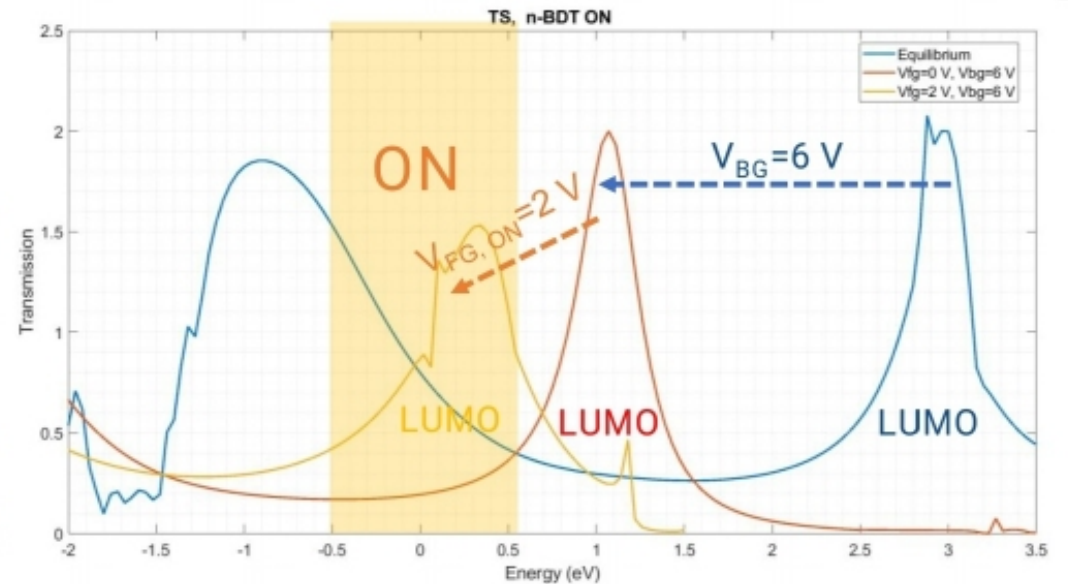
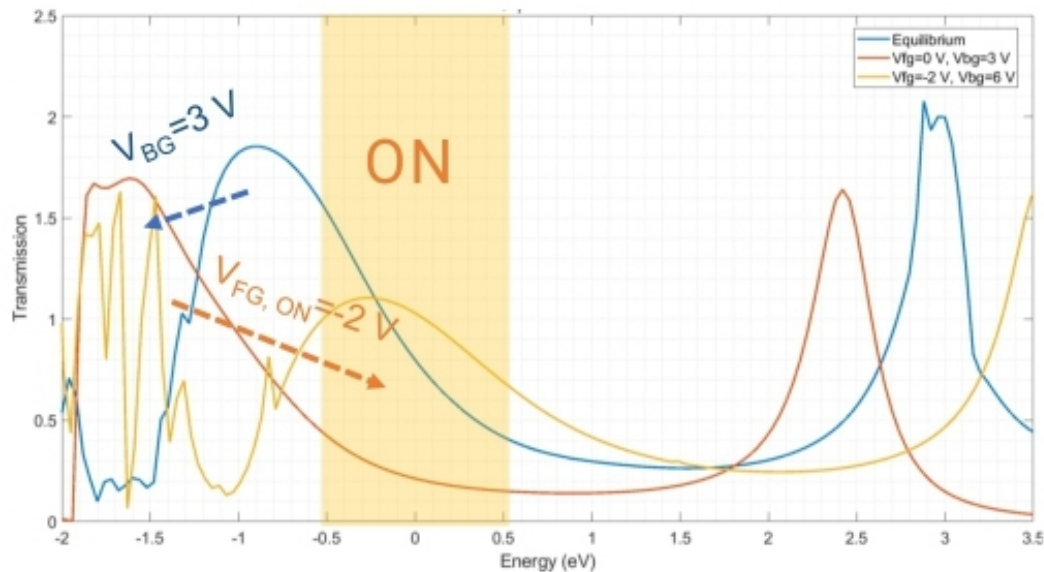
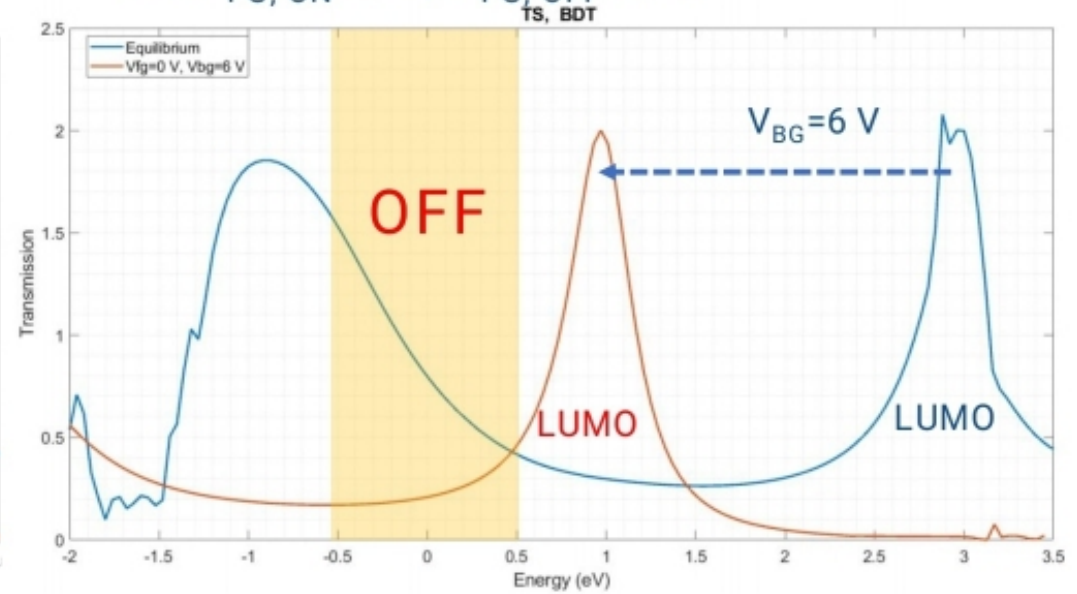
HOMO conduction (p-type)

- $V_{FG, ON} = -2 \text{ V}$, $V_{FG, OFF} = 0 \text{ V}$



LUMO conduction (n-type)

- $V_{FG, ON} = 2 \text{ V}$, $V_{FG, OFF} = 0 \text{ V}$



FOR THE LAST CASE

BENZENE

P TYPE

$$I_{ON} = 40 \mu A$$

$$I_{OFF} = 10 \mu A$$

N TYPE

$$I_{ON} = 40 \mu A$$

$$I_{OFF} = 7 \mu A$$

@ 1V

CONSIDERATION ON BENZENE

- CAN BE USED AS PTYPE & NTYPE
- THERE IS AN APPRECIABLE DIFFERENCE
BW I_{ON} AND I_{OFF}

→ BUT $-I_{OFF}$ IS NOT SMALL
————→ POWER!!

$-\frac{I_{ON}}{I_{OFF}}$ IS POOR FOR DIGITAL
APP

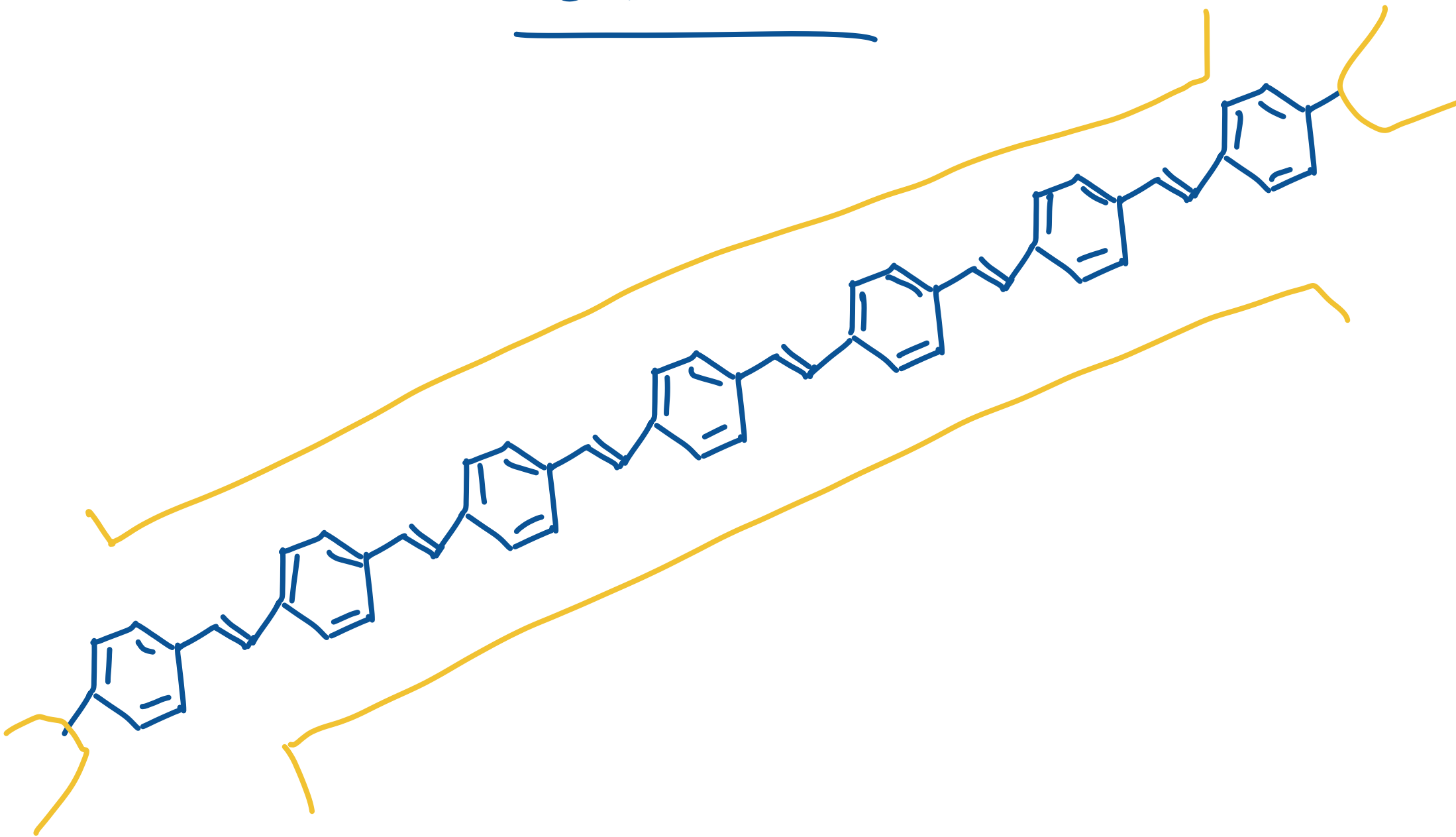
WE CAN DO BETTER!

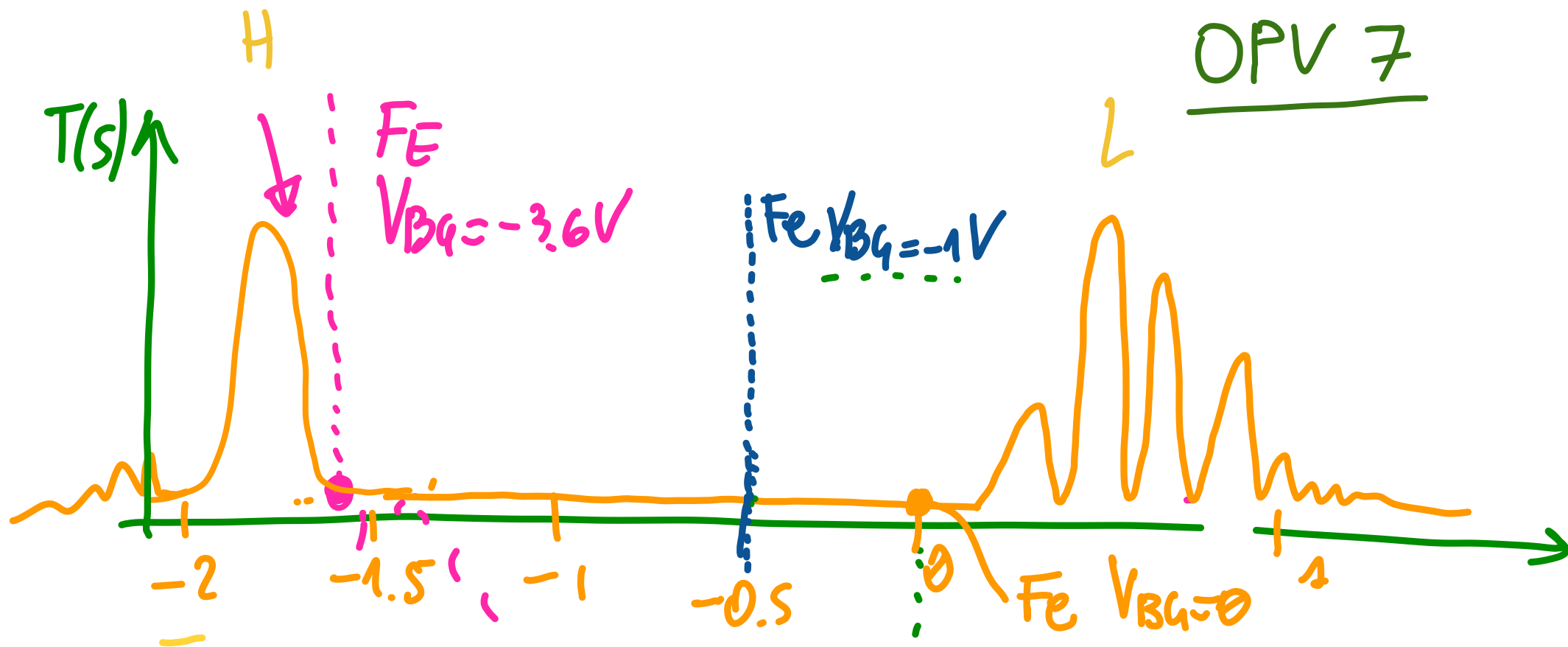
• CASE 1 OPV7

• CASE 2 PCP

NOTE: V_{BQ1} and V_{BQ2} could have different β , either by necessity or by choice; different β complicate tech.

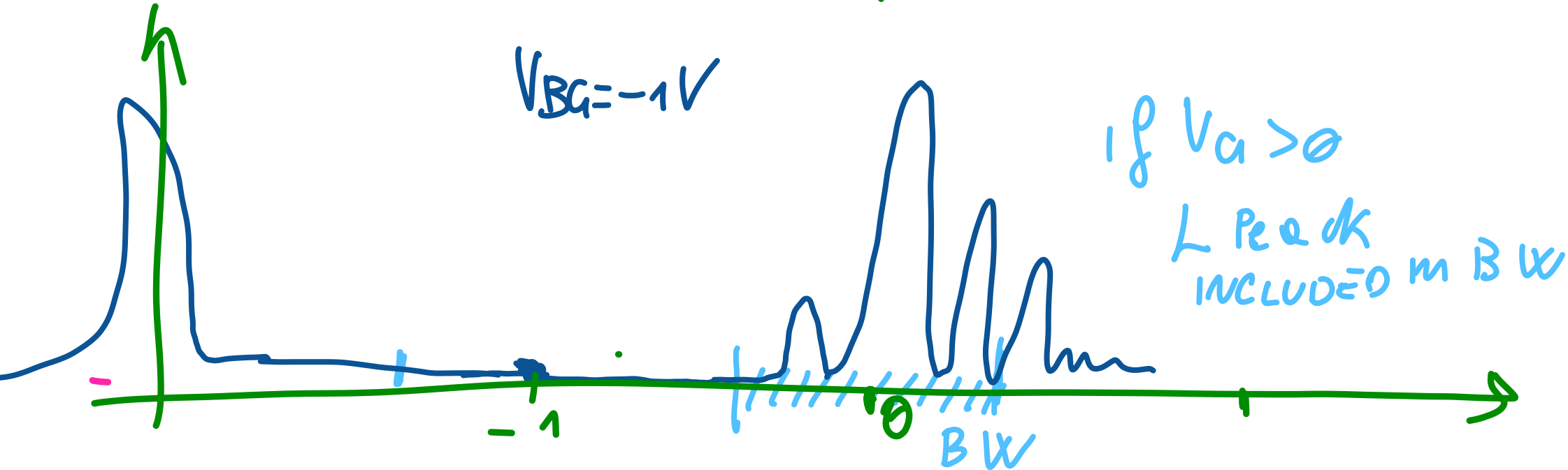
OPV7



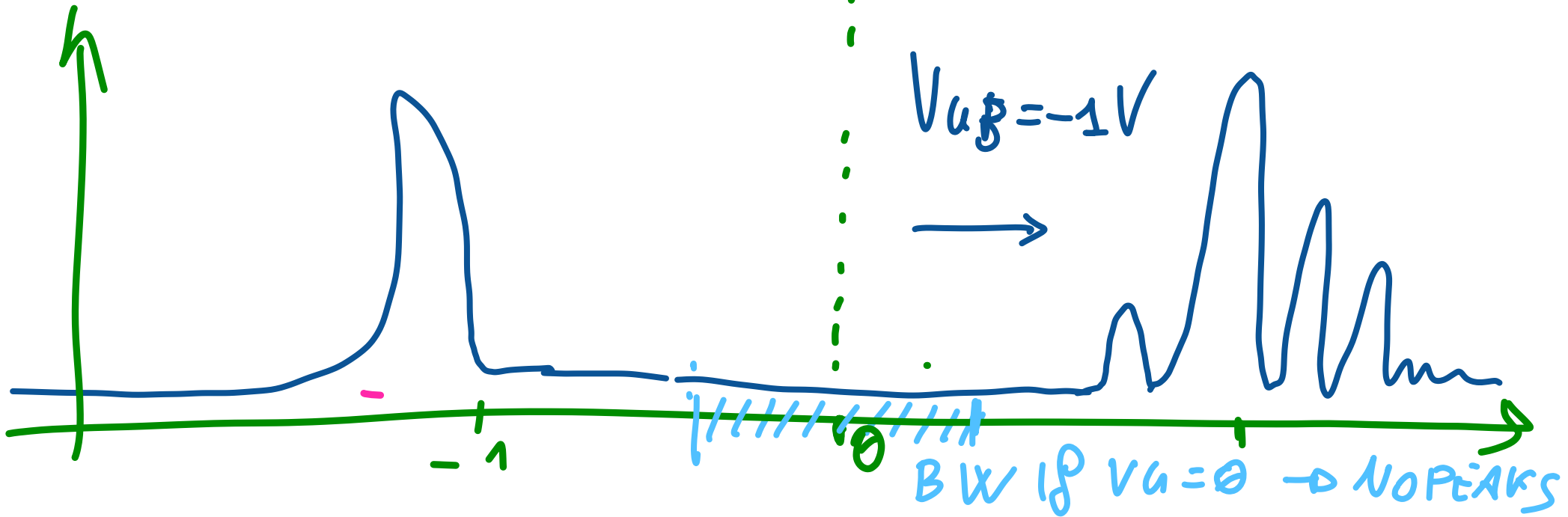
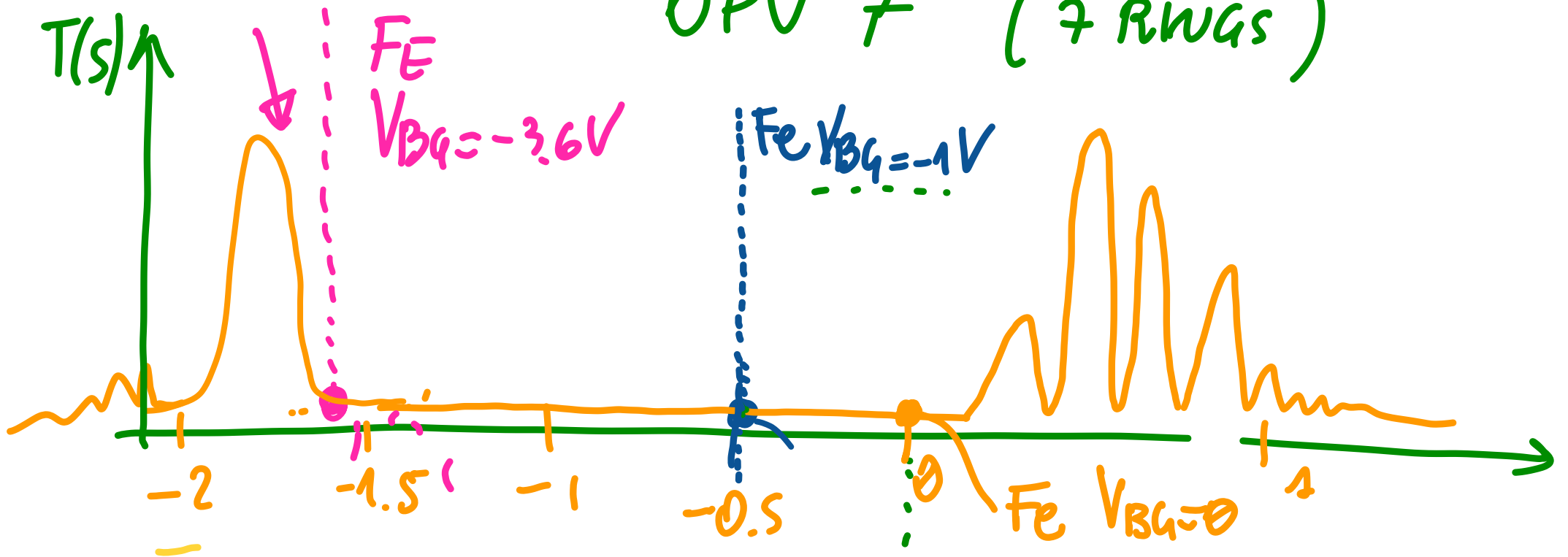


→ USE V_{BQ} TO CHOOSE THE PEAK TO INCLUDE WHEN V_G "ACTIVE"

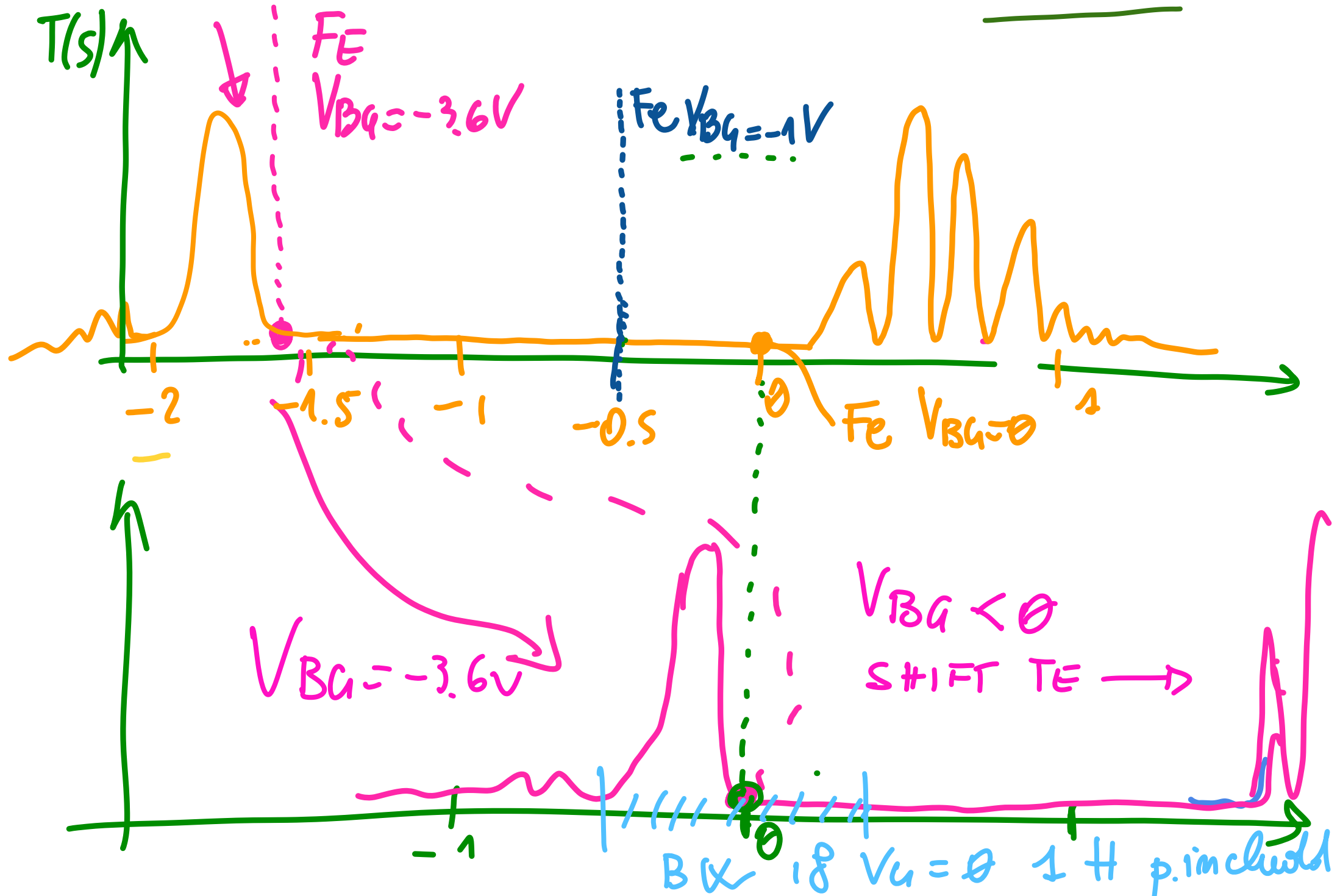
OPV 7 (7 RWGS)



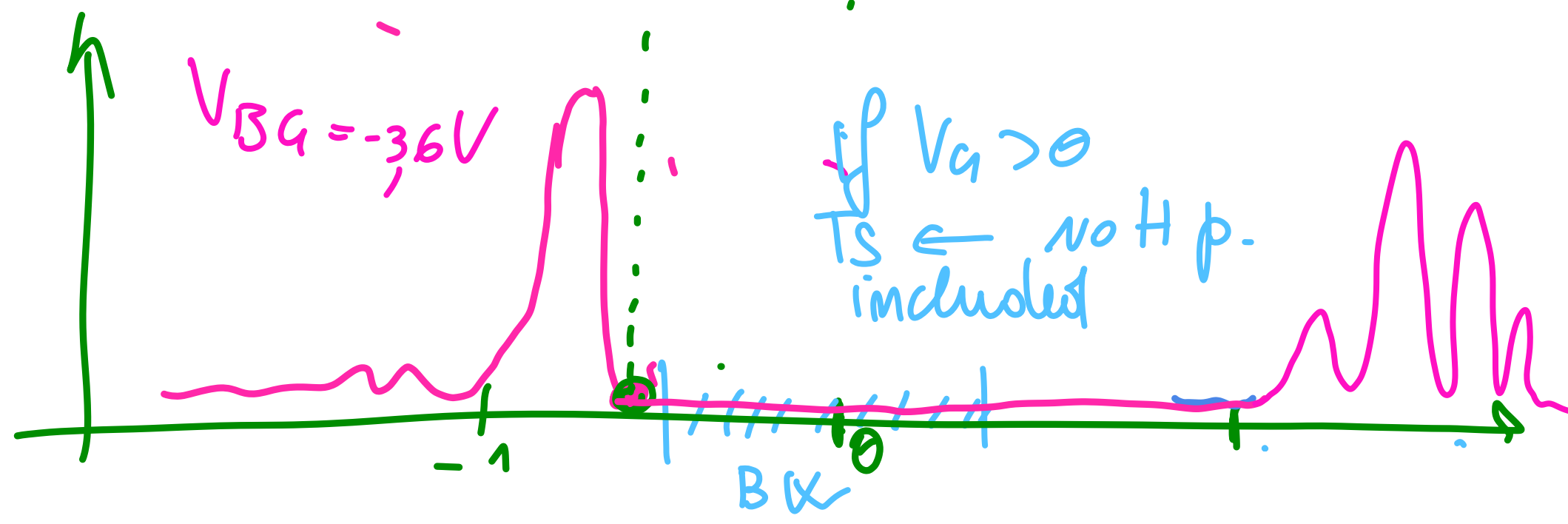
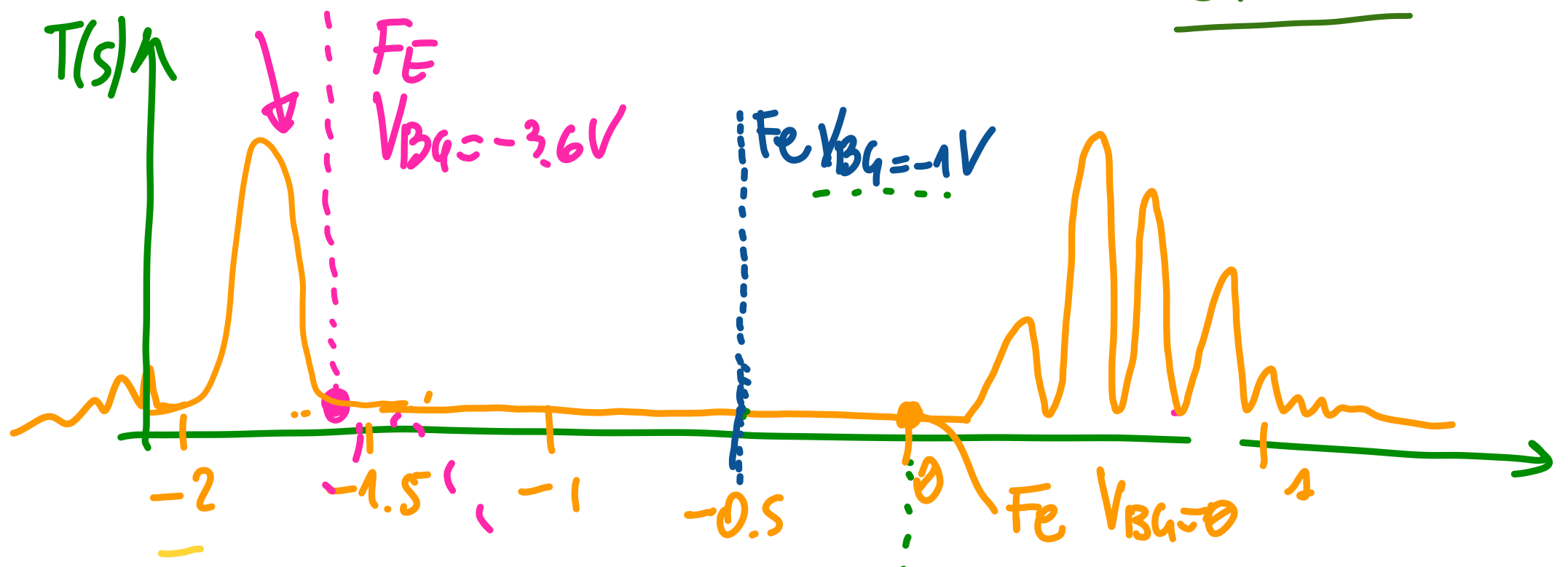
OPV 7 (7 RWGS)



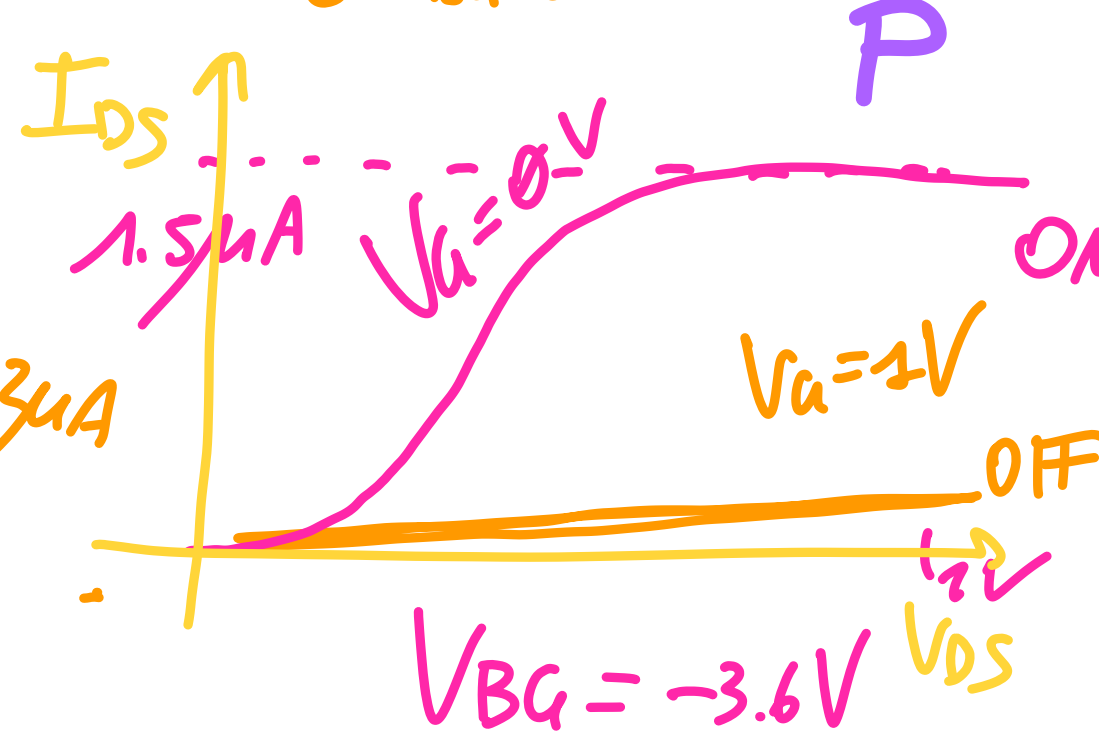
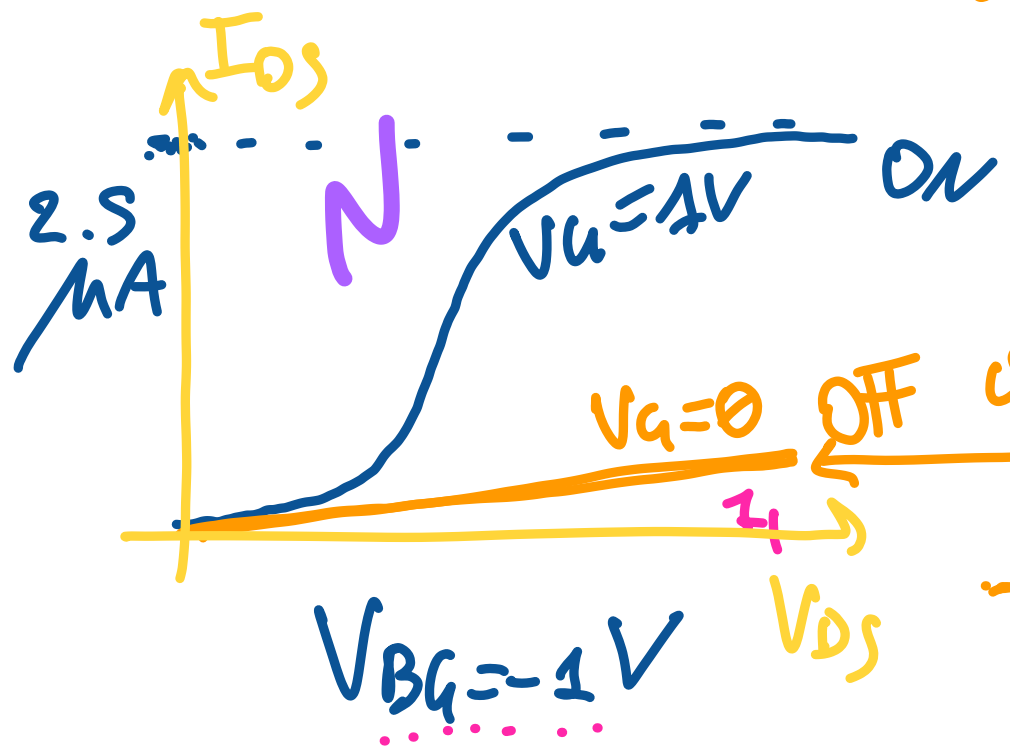
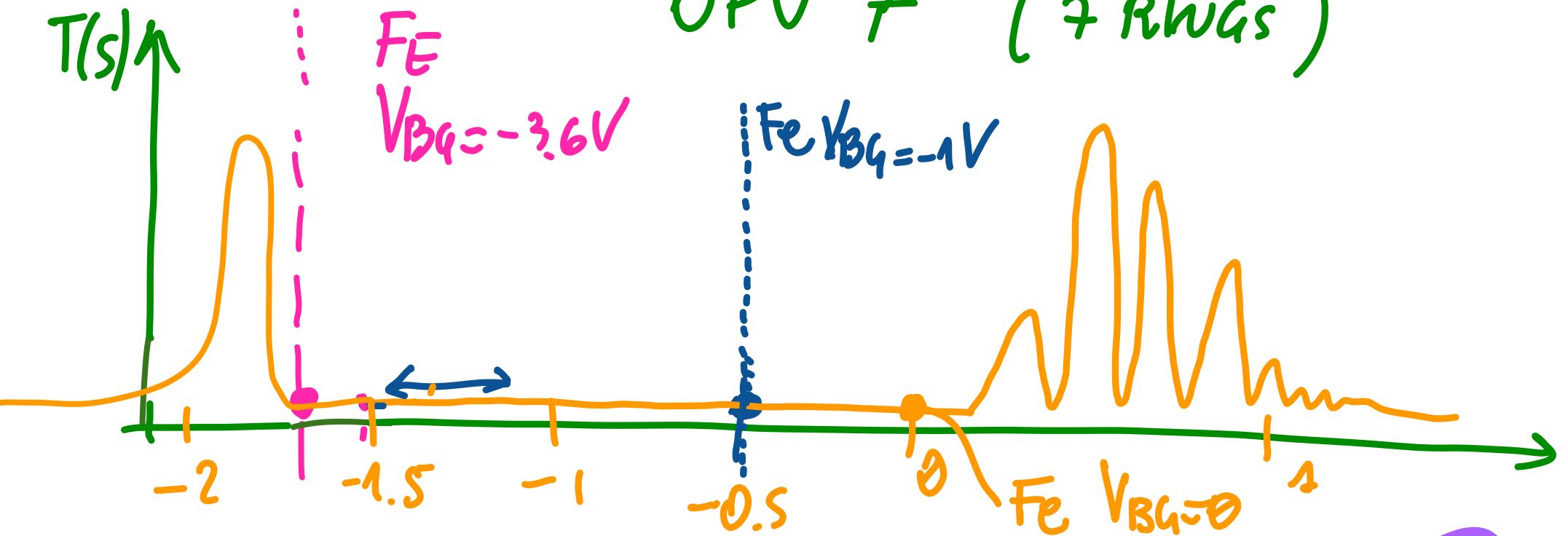
OPV 7



OPV 7



OPV 7 (7 RWGS)



OPV 7

- PTYPE & NTYPE OK ☺

N type	$V_G = 1V$ ON
	$V_G = 0V$ OFF

$V_G = 1V$ OFF	P type
$V_G = 0V$ ON	

COMPLEMENTARY

BEHAVIOUR

- GOOD

$$\frac{I_{ON}}{I_{OFF}}$$

N - $\frac{I_{ON}}{I_{OFF}} \sim 12$..

P - $\frac{I_{ON}}{I_{OFF}} \sim 54$.

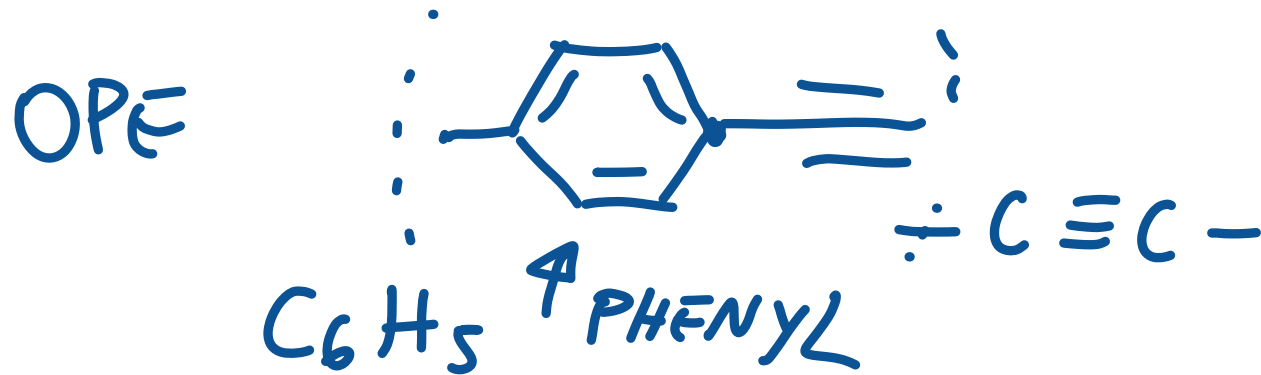
∧ I_{OFF} NOT

SO SMALL

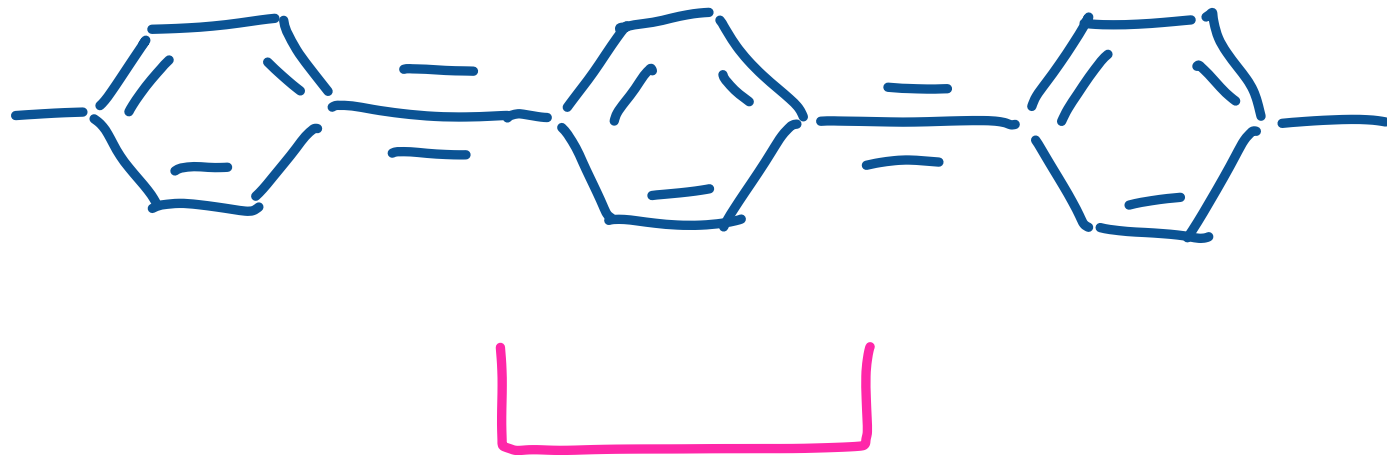
PCP ↓ !!!

BETTER THAN BENZENE

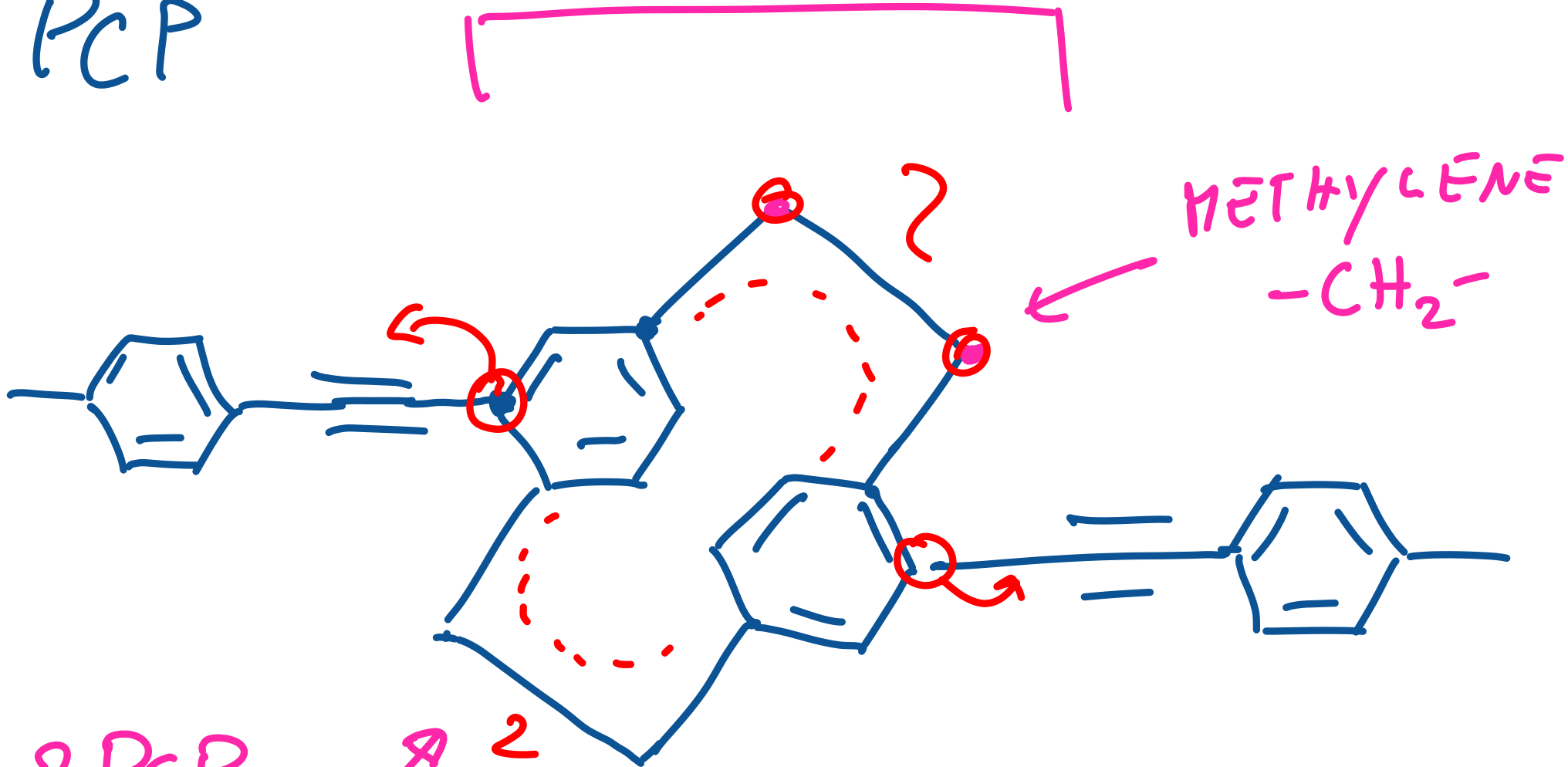
CASE **PCP** PARACYCLOPHANE
BASED ON OPE



OPE3



PCP



2,2 PCP

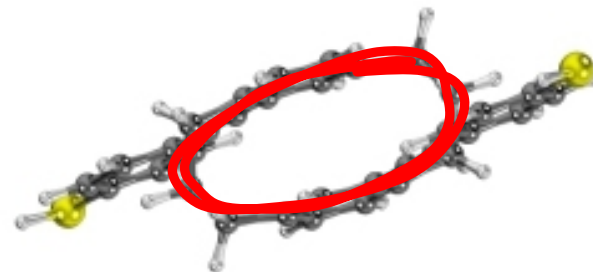
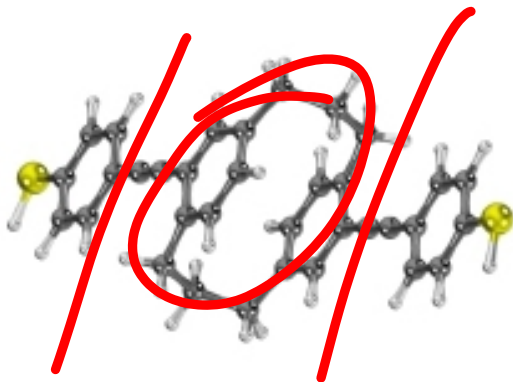
BRIDGE WITH

A CHAIN OF
2 METHYLENE.

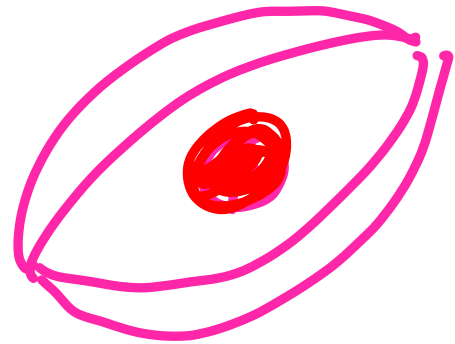
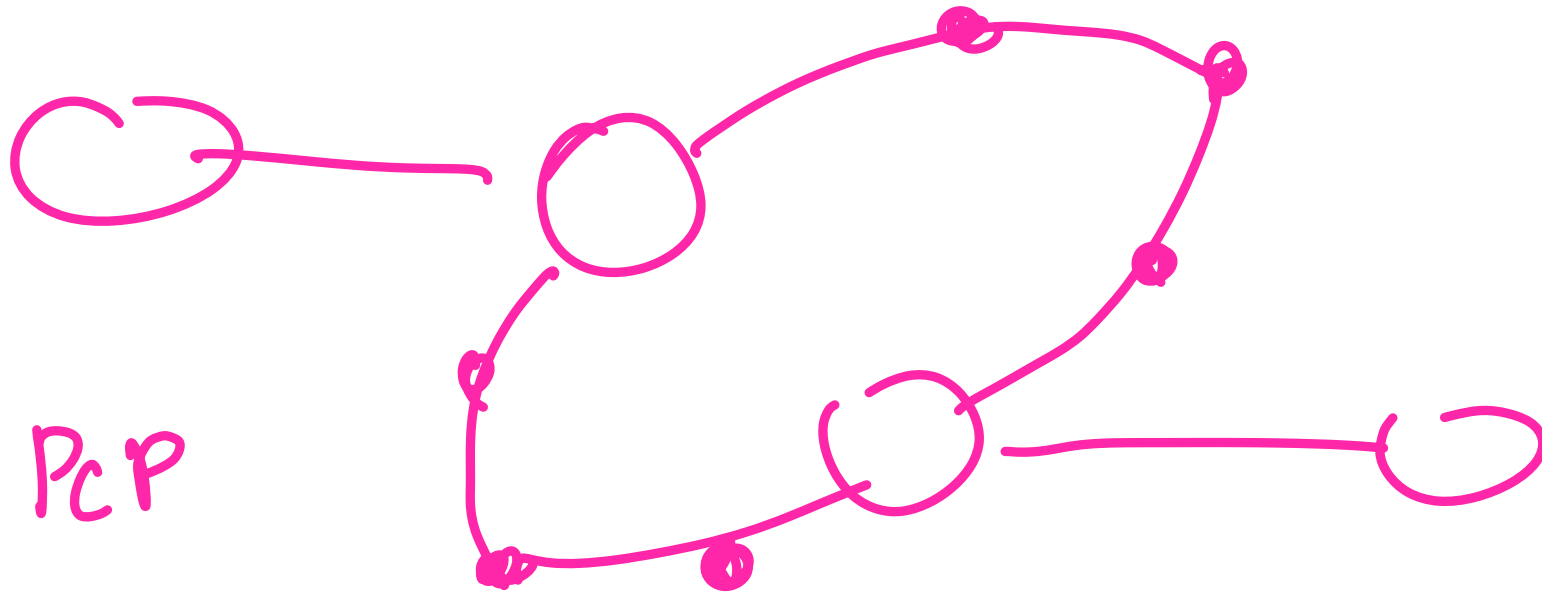
3,3 PCP → BRIDGE

3 METH...

2,2 PCP



3,3 PCP

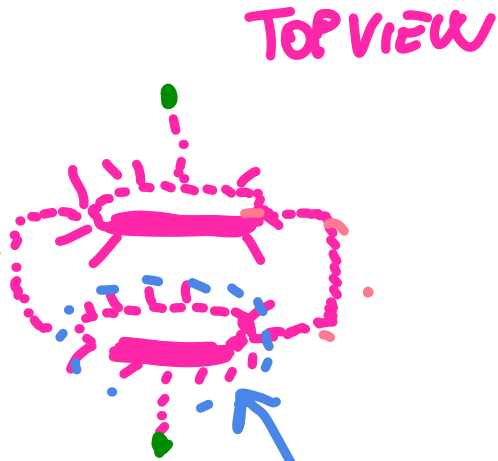


BOAT LIKE
ORG.

CAN BE USED AS A
SENSOR

2,2 PCP

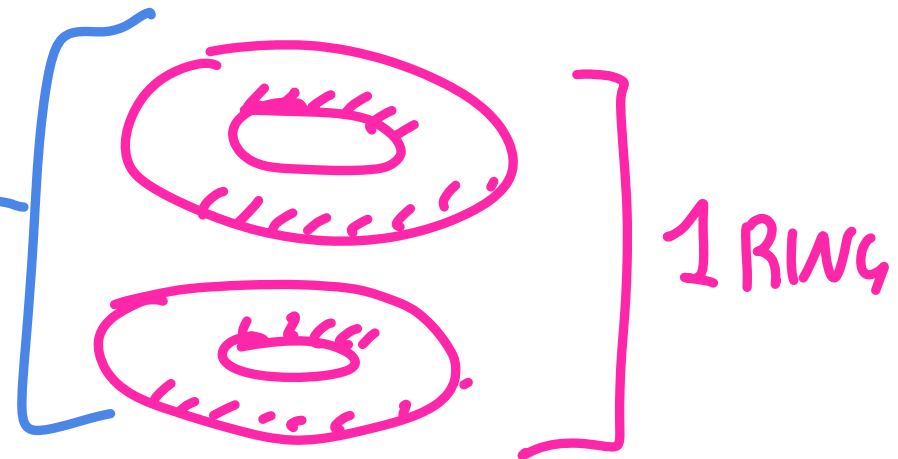
THE BRIDGE
INFLUENCE
THE CONJ. SYSTEM



each
RING
CONJUGATED
SYSTEM
OF π
ORBITALS

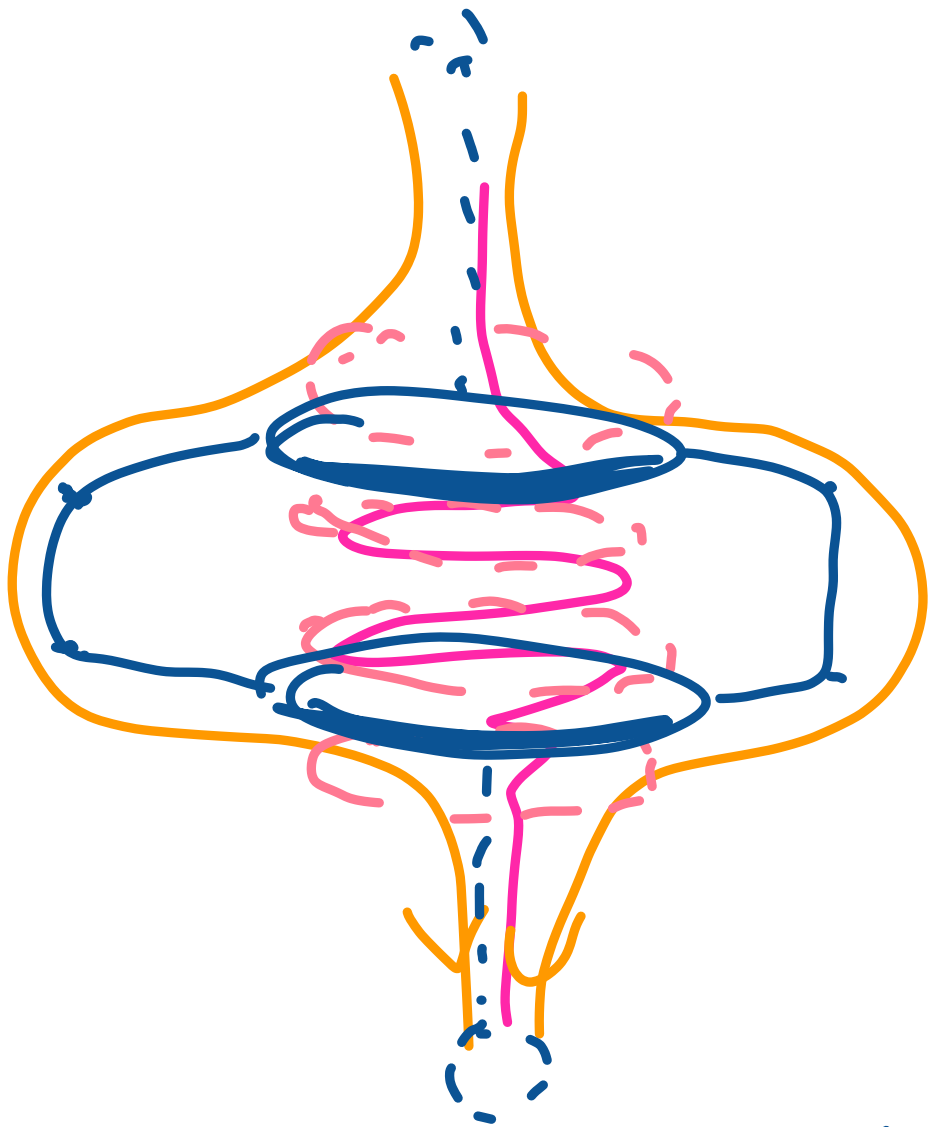
2 RINGS
FACED:

2 FACED
CONJ. SYSTEMS



- 1 CURRENT PATH in π - π OVERLAP BW
THE 2 FACED RINGS I_c
- 1 CURRENT PATH THROUGH THE BRIDGES I_B

\hookrightarrow IN PARALLEL \rightarrow GENERATE
 DESTRUCTIVE QUANTUM
 D QI INTERFERENCE
 $H_{010} \& H_{020-1}$



I_C —
 I_B —

for some
ENERGIES

INTERFERE
DESTRUCTIVELY

D (QI)

$T(E)$

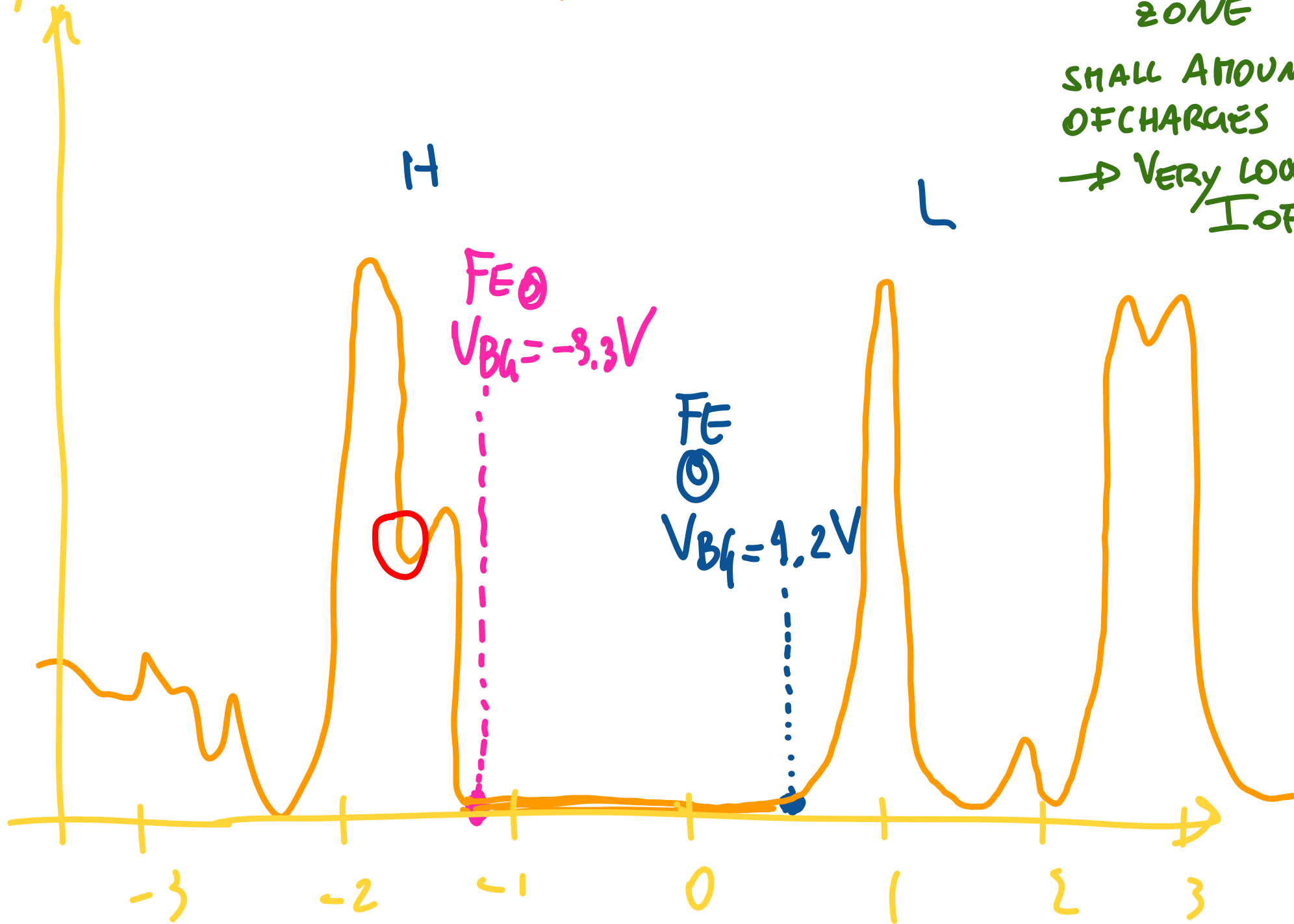
NORMAL SCALE

2,2 PCP + 2 THIOL

IN HLG ZONE

SMALL AMOUNT OF CHARGES

→ VERY LOW I_{OFF}



N TYPE

$$V_{B_{on}} = 1.2V$$

ON	$V_a = 1V$	} ←
OFF	$V_a = -1V$	

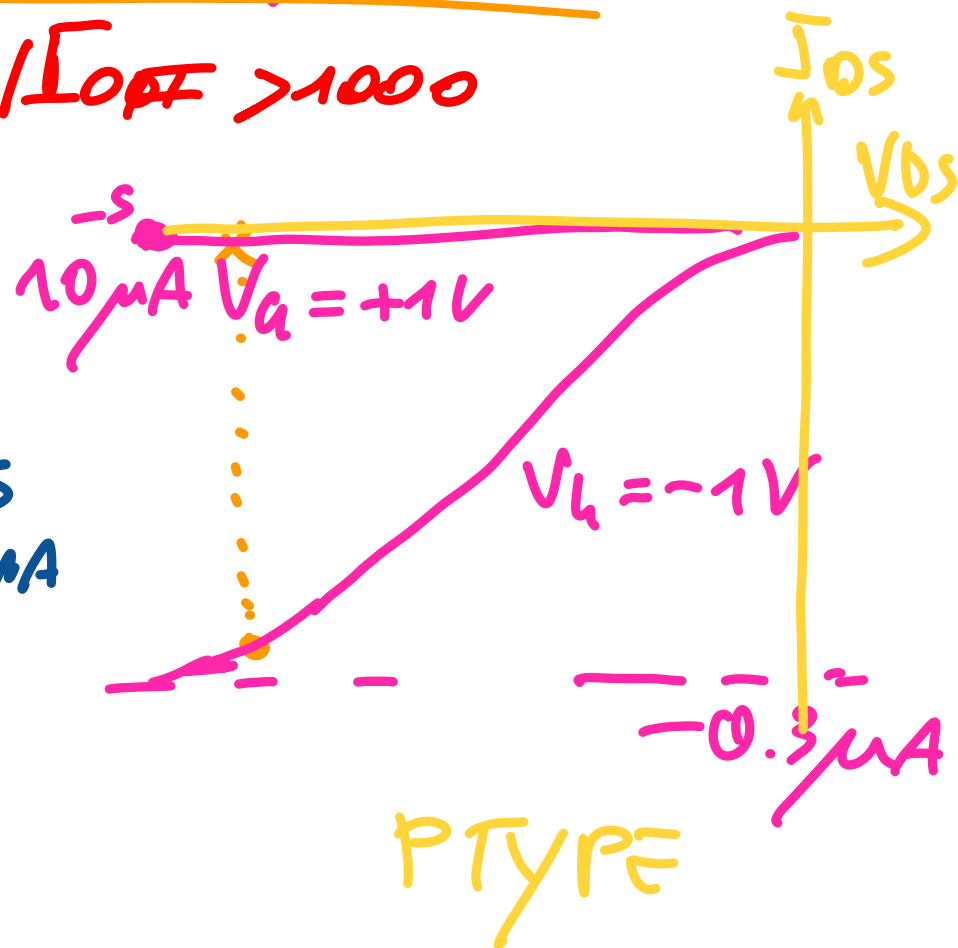
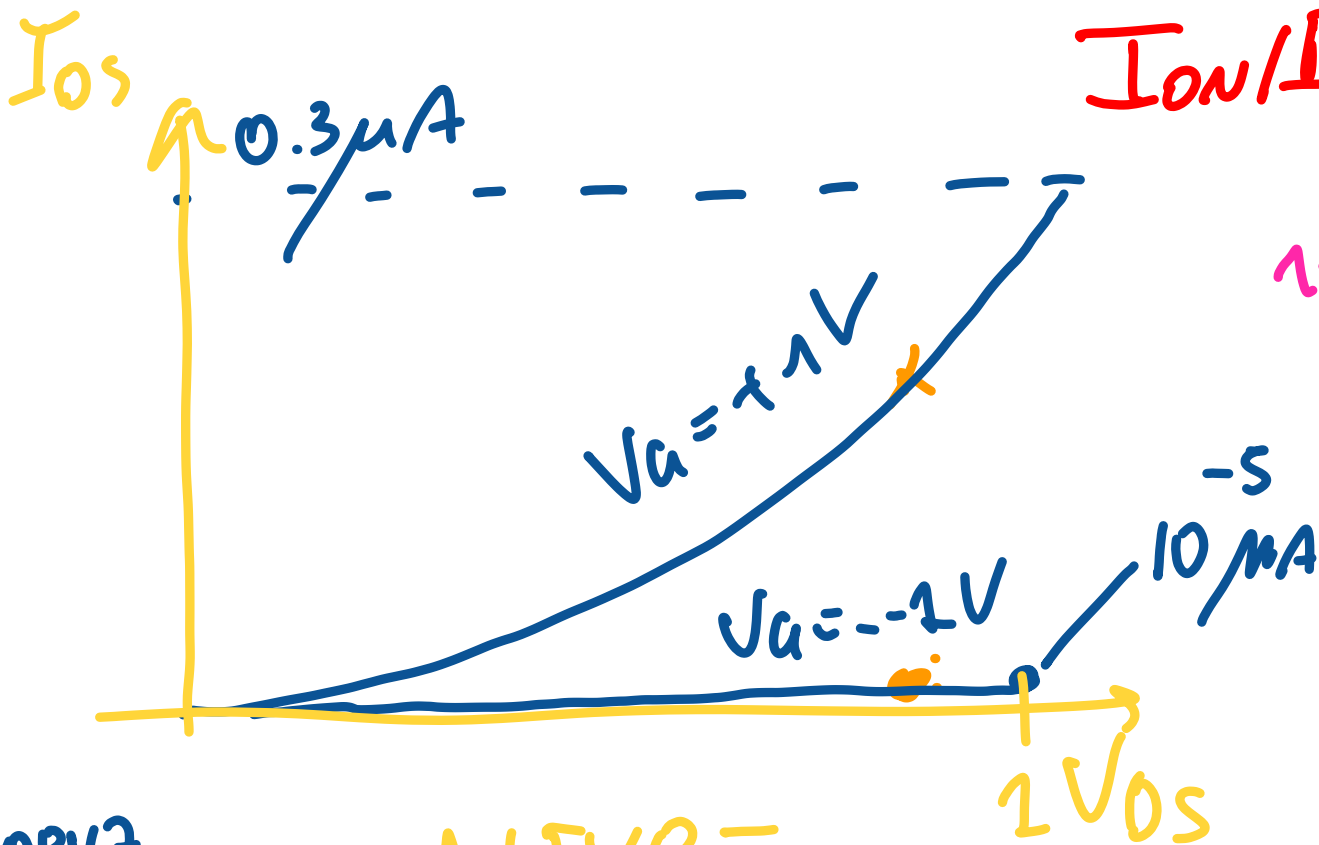
AMBI-POLAR

P TYPE

$$V_{B_{on}} = -3.3V$$

ON	$V_a = -1V$	} ←
OFF	$V_a = 1V$	

$$I_{ON}/I_{OFF} > 1000$$



OPV7
 $I_{OFF} \approx 0.02 \mu A$

N TYPE

P TYPE

→ GOOD AMBIPOLAR BEHAVIOR
(DISTINCT PEAKS, VERY CLEAR)

→ VERY LOW I_{OFF} → GOOD $\frac{I_{ON}}{I_{OFF}}$

→ V.R.T. OPV7

$$I_{OFF} \Big|_{OPV7} \sim 2 \cdot 10^{-2} \mu A$$

$$I_{OFF} \Big|_{PCP_{212}} \sim 10^{-5} \mu A$$

⇒ LOWER POWER ⇒ BETTER BEHAVIOR IN A LOGIC GATE (NEXT SECTION)

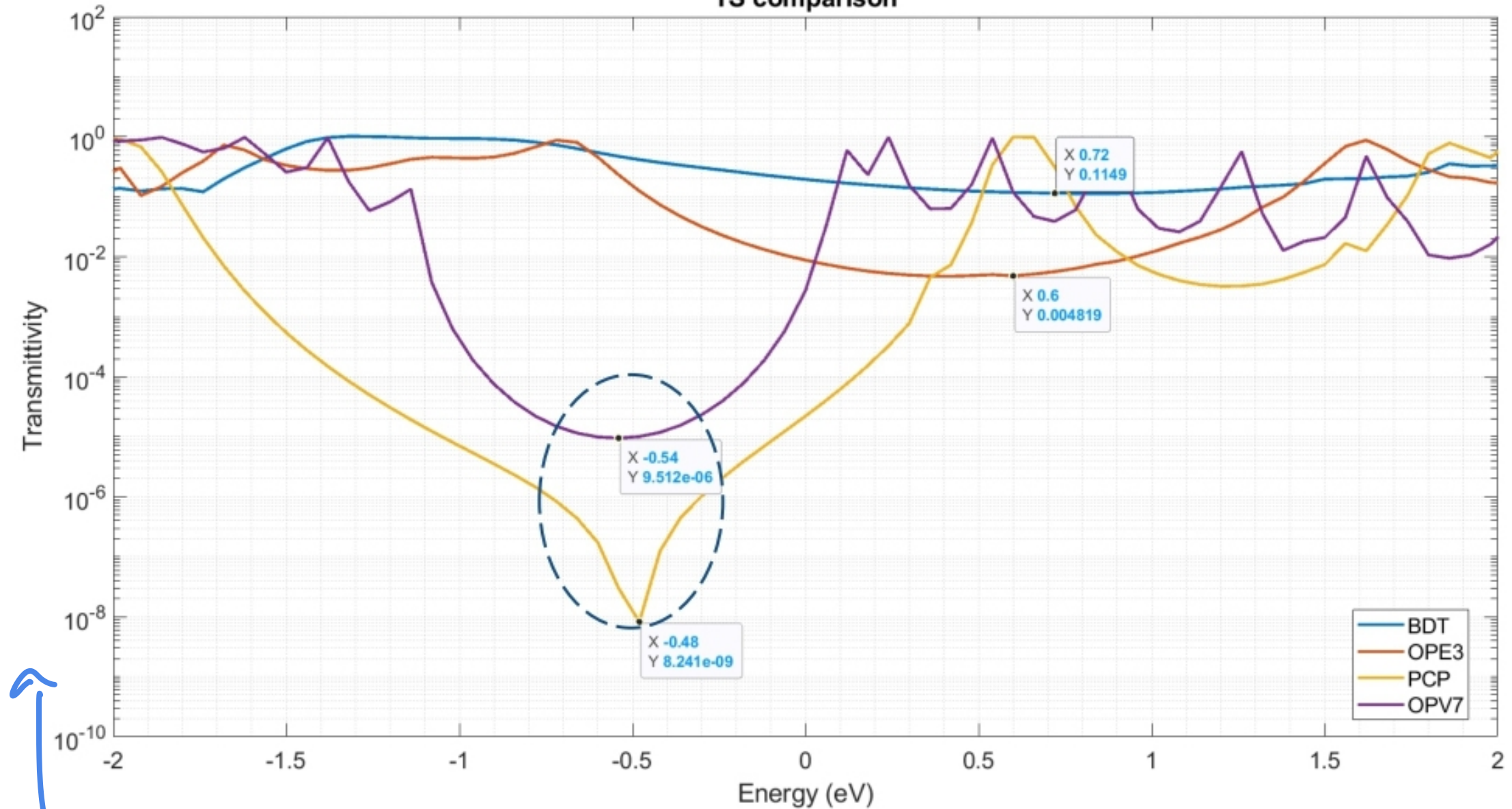
BDT

OPE3

OPV7

PCP

TS comparison



LOG SCALE

QI FUNDAMENTAL FOR 4000 I_{ON}/I_{OFF} RATIO

CRISTALLIZATION

a) CHARGING EFFECT & RESONANCE

b) EFFECT OF GATE

- GENERAL

- BENZENE

- OPV7

- PCP & Q.I.