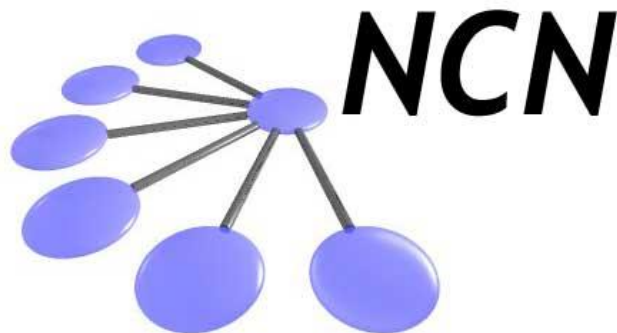


Network for Computational Nanotechnology (NCN)

University of California, Berkeley; University of Illinois at Urbana-Champaign; Norfolk State University; Purdue University; University of Texas at El Paso

First-Time User Guide to MOSCAP*



PURDUE
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*<http://nanohub.org/resources/moscap>

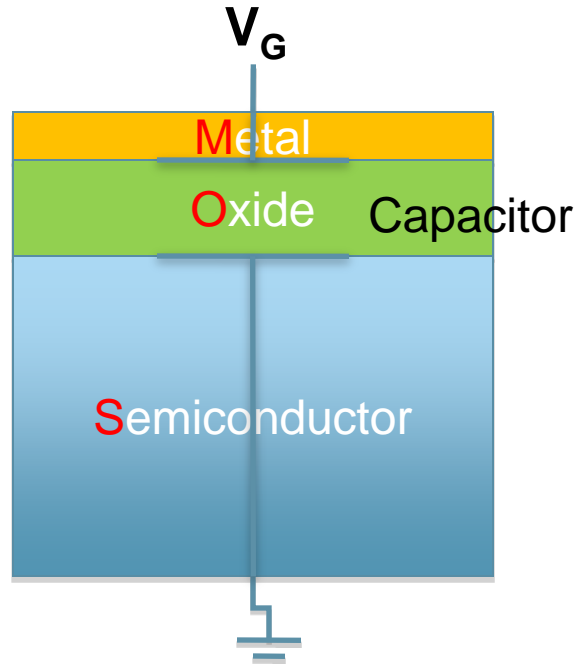
**kim568@purdue.edu

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Introduction: What is a MOS Capacitor?

MOS Capacitor: Metal Oxide Semiconductor^[1] Capacitor



- Metal : metal or poly-silicon material
- Oxide : SiO_2 or high-k dielectric material
- Semiconductor : p-type or n-type semiconductor material
- Importance of MOS Capacitors
 - » Essential for understanding MOSFET*^[2]
 - » Basic structural part of MOSFET

[1] Dragica Vasileska (2008), "MOS Capacitors Description," <http://nanohub.org/resources/5087>

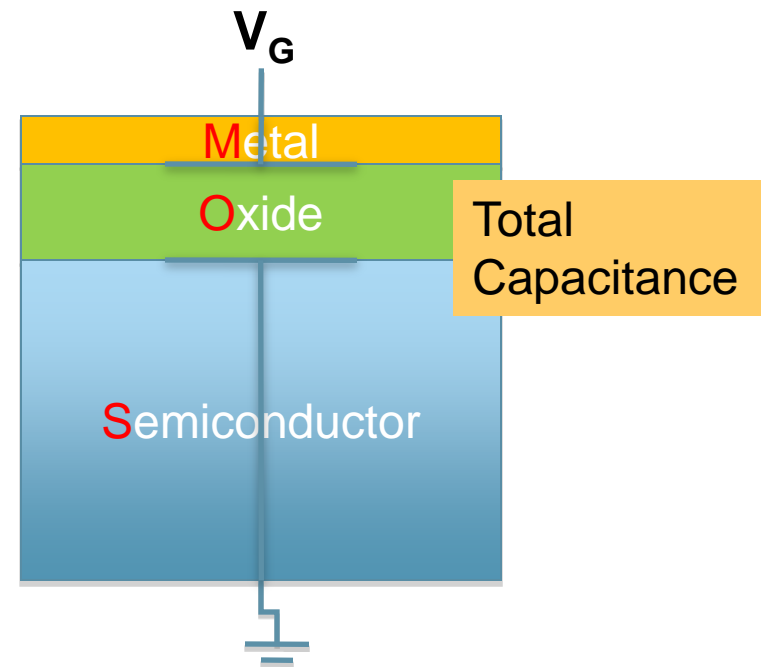
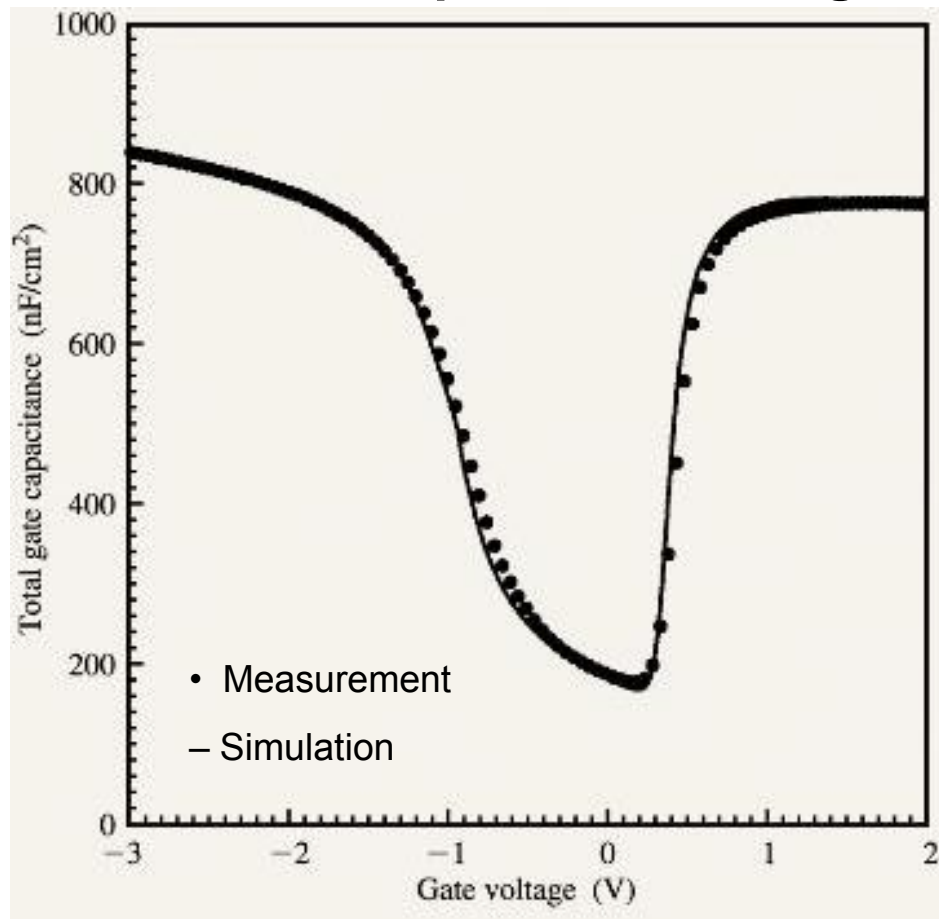
*MOSFET : Metal Oxide Semiconductor Field Effect Transistor has a source and a drain as an additional contacts to the gate contact as shown in the picture above.

For more information, refer to the following reference [2]

[2] Dragica Vasileska (2008), "MOSFET Operation Description," <https://nanohub.org/resources/5085>

Introduction: What Can Be Measured in MOS Capacitors?

Capacitance-Voltage Characteristics^[3]

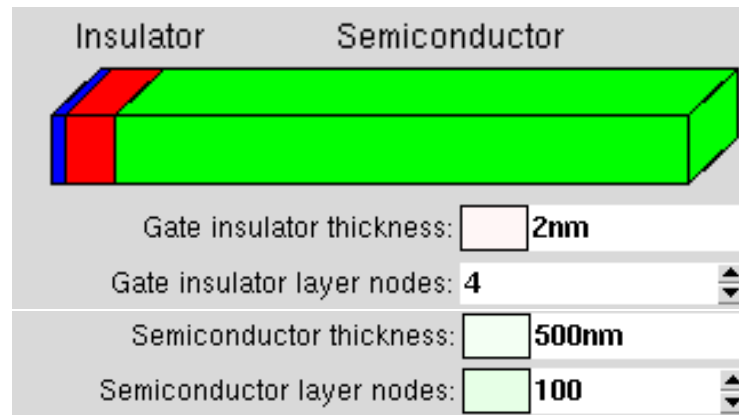


[4] S.-H. Lo, et. al., IBM Journal of Research and Development, volume 43, number 3, 1999

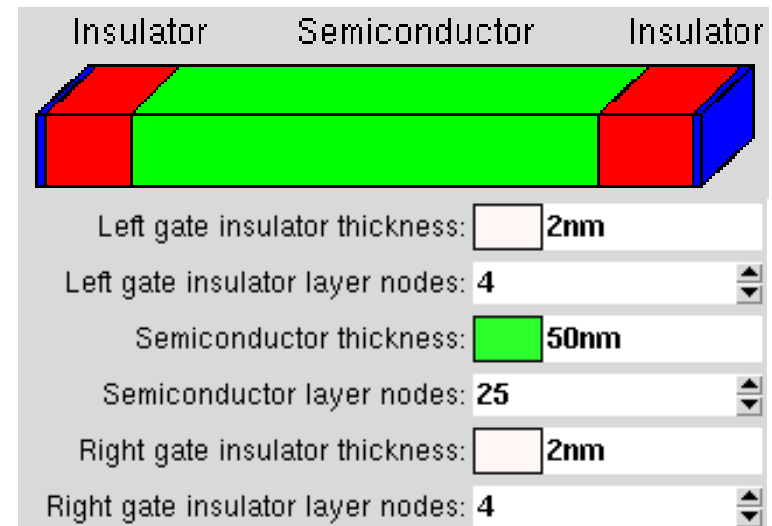
What Can Be Simulated by MOSCAP? (input):

MOS Capacitor with different geometry

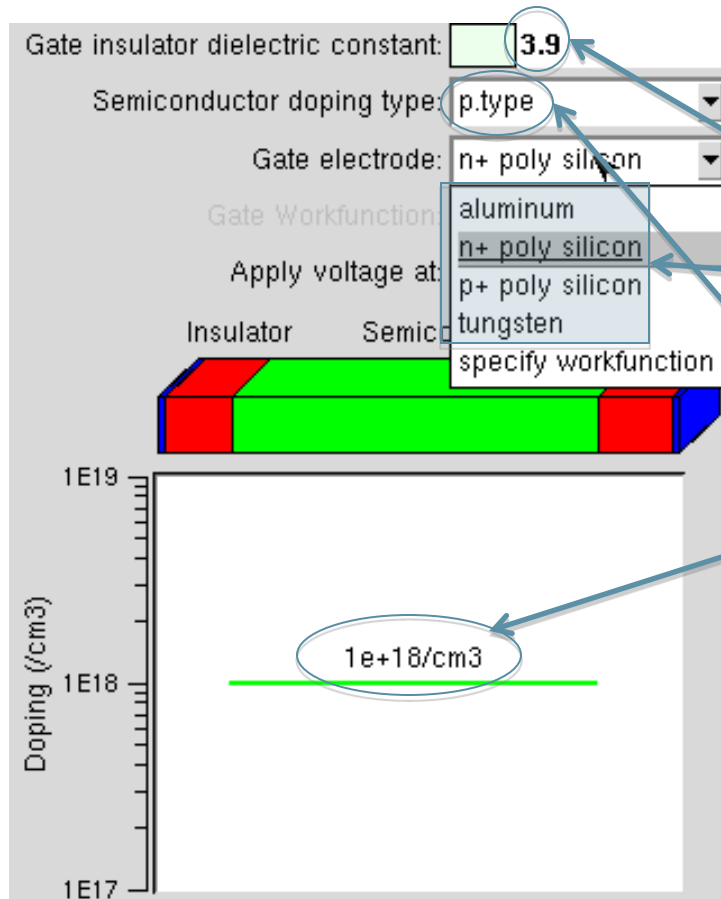
Single Gate



Double Gate



What Can Be Simulated by MOSCAP? (input):



MOS Capacitor with different doping and material

- Insulator dielectric constant
- Gate electrode type
- Gate workfunction* specification
- Semiconductor doping type
- Semiconductor doping density

*Workfunction is the minimum energy that is needed to remove an electron from a solid to a point immediately outside the solid surface
-http://en.wikipedia.org/wiki/Work_function

What Can Be Simulated by MOSCAP? (input):

- Environment parameters
- Special parameters(charge density in the insulator)

The image shows the 'Parameters' tab of a MOSCAP simulation interface. The 'Fixed charge density in gate insulator (/cm3)' and 'Interface trap charge density in gate insulator (/cm2)' are both set to 0. A diagram to the right illustrates the physical structure of a MOS capacitor, showing layers of SiO₂, SiO_x, and Si. It identifies three types of charges: Mobile Ionic Charge (K⁺ and Na⁺ ions), Oxide Trapped Charge (positive charges in the oxide), and Interface Trapped Charge (charges at the Si-SiO_x interface).

Parameters shown in the screenshot:

- Device: Parameters (circled in green)
- Environment: Environment
- Ambient temperature: 300K
- Initial voltage (V): -2
- Final voltage (V): 4
- Number of Voltage Steps: 40
- Frequency for AC analysis: low

Charge density parameters (highlighted in green):

- Fixed charge density in gate insulator (/cm3): 0
- Interface trap charge density in gate insulator (/cm2): 0

Diagram labels:

- MOBILE IONIC CHARGE (K⁺, Na⁺)
- OXIDE TRAPPED CHARGE (+)
- FIXED OXIDE CHARGE (+)
- INTERFACE TRAPPED CHARGE (x)
- Layers: SiO₂, SiO_x, Si

[5]Deal B. E., Electron Devices, IEEE Transactions on, 1980

What If You Just Hit the “Simulate” Button? (output):

Device | Parameters | Environment

Model: Single Gate

Gate insulator thickness: 2nm

Gate insulator layer nodes: 4

Gate insulator dielectric constant: 3.9

Semiconductor thickness: 500nm


Semiconductor layer nodes: 100

Semiconductor doping type: p.type

Gate electrode: n+ poly silicon

Gate Workfunction: 4.05eV

Insulator Semiconductor



Simulate

V1 = -1.7000000E+00 V2 = 0.0000000E+00

Simulation starts with the default input parameters!!

```

Pr
o-i
1 0 7.1775E-01
1 1 3.5102E+00
1 1 2.6705E-02
1 2* 2.3089E-02
1 2* 2.7859E-04
1 3* 2.4722E-04
1
1
1
Solution for bias:
V1 = -1.0000000E+00    V2 = 0.0000000E+00

Previous solution used as initial guess
o-itr i-itr psi-error n-error p-error
1 0 7.9069E-01
1 1 6.6732E+00
1 1 1.3544E-01

```


What If You Just Hit the “Simulate” Button? (output):

C-V characteristic

C/Cox-V characteristic

Energy band diagram (at VG = 0)

Electron density (at VG = 0)

Hole density (at VG = 0)

Net charge density (at VG = 0)

Electrostatic potential (at VG = 0)

Electric field (at VG = 0)

Energy band diagram (at last applied bias)

Electron density (at last applied bias)

Hole density (at last applied bias)

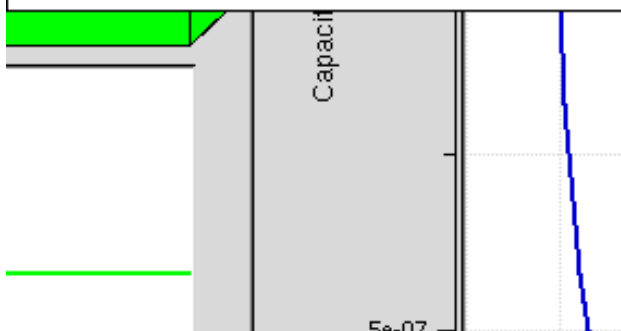
Net charge density (at last applied bias)

Electrostatic potential (at last applied bias)

Electric field (at last applied bias)

Output Log

Download



- **[C-V Characteristics]** : Gate capacitance vs Voltage characteristics
- **[C/Cox-V Characteristics]** : Gate capacitance divided by oxide capacitance vs Voltage characteristics
- **[Energy band diagram]** : Conduction/Valence Band, Fermi level diagram
- **[Electron density]** : electron density in the semiconductor
- **[Hole density]** : hole density in the semiconductor
- **[Net charge density]** : $\rho = q(N_A - N_D + n - p)$
- **[Electrostatic potential]** : $V = \text{const} - q \cdot E_c$
- **[Electric field]** : $E = -\nabla V$

ρ : charge density

N_A/N_D : acceptor/Donor density

n/p : electron/hole density

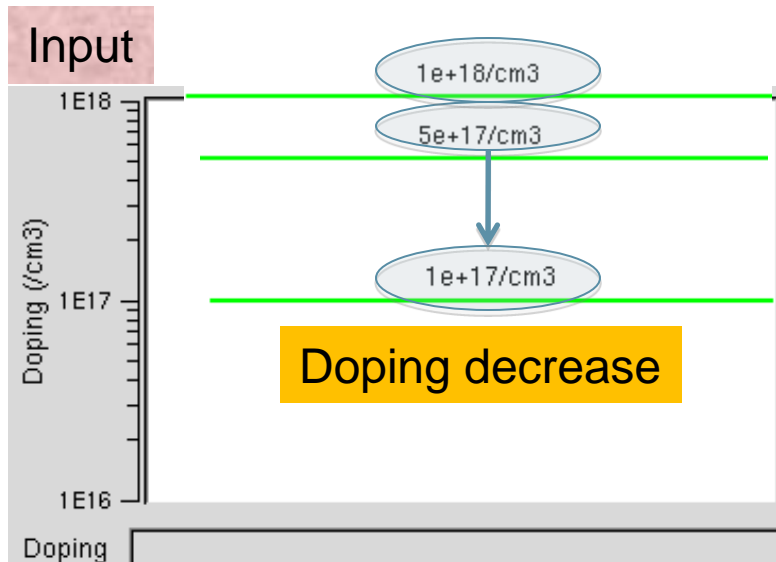
q : electron charge

E_c conduction band edge

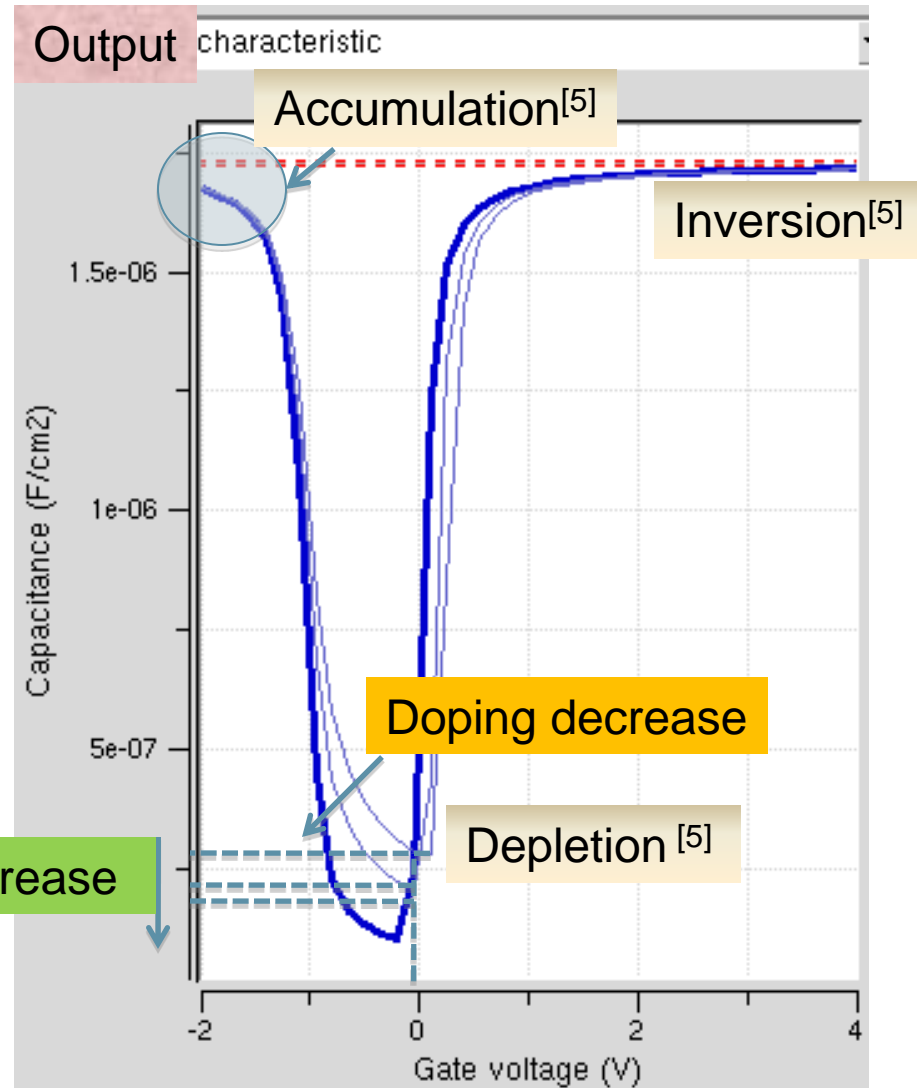
E : electric field

V : electrostatic potential

Example of Simulation (if the doping is changed)



$C_{\text{tot}} |_{V_G=0}$ decrease



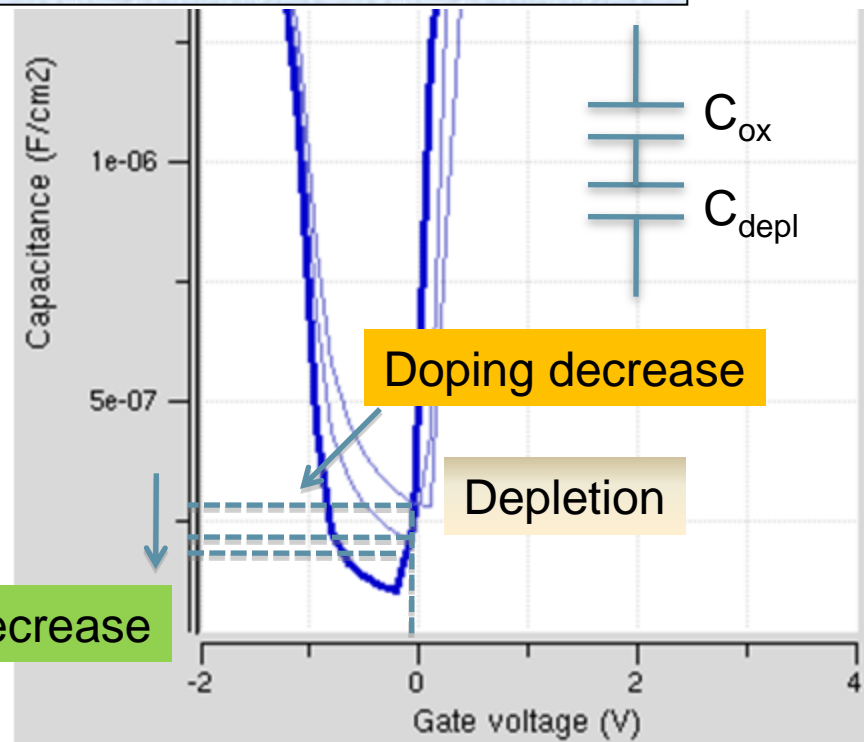
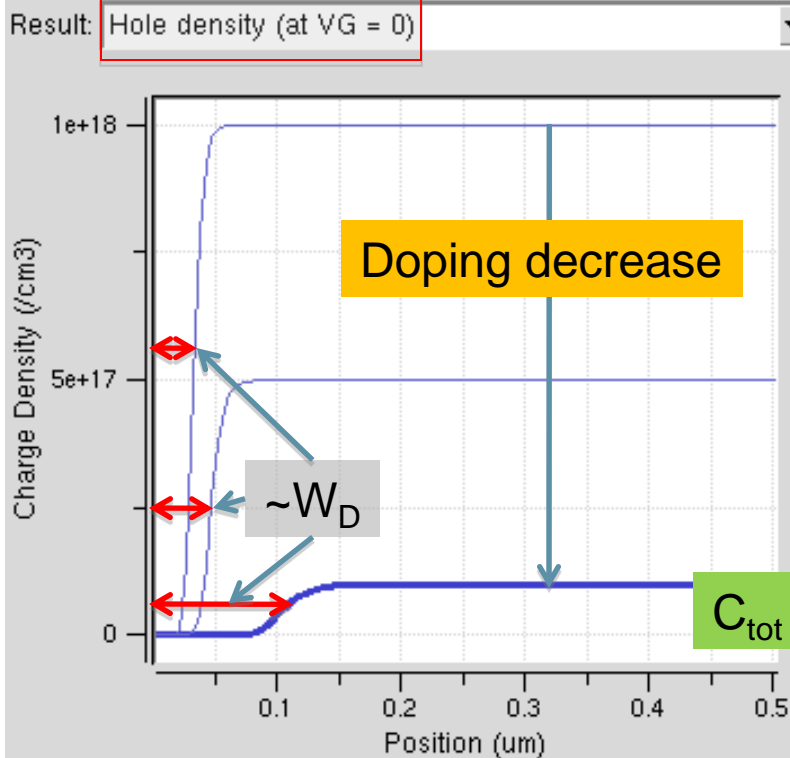
[5] Mark Lundstrom (2008), "ECE 612
Lecture 3: MOS Capacitors,"
<http://nanohub.org/resources/5363>

Example of Simulation (if the doping is changed)

$$C_{tot} = \frac{1}{\frac{1}{C_{ox}} + \frac{1}{C_{Depl}}}, C_{depl} = \frac{\epsilon_S}{W_D}, C_{ox} = \frac{\epsilon_{ox}}{d_{ox}}$$

$C_{tot} / C_{ox} / C_{depl}$: total/oxide/depletion capacitance
 $\epsilon_{ox} / \epsilon_S$: dielectric constant of oxide/semi conductor
 W_D : depletion width, d_{ox} : oxide thickness

W_D increase \rightarrow C_{depl} decrease \rightarrow C_{tot} decrease

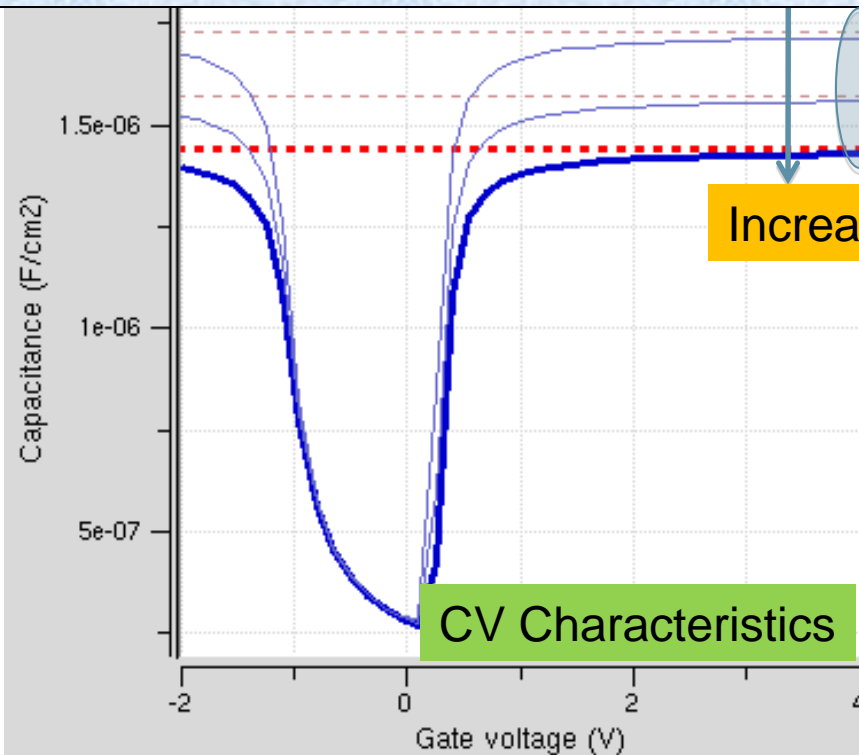


Example of Simulation (if the oxide thickness is changed)

$$C_{tot} = \frac{1}{\frac{1}{C_{ox}} + \frac{1}{C_S}}, C_{ox} = \frac{\epsilon_{ox}}{d_{ox}}$$

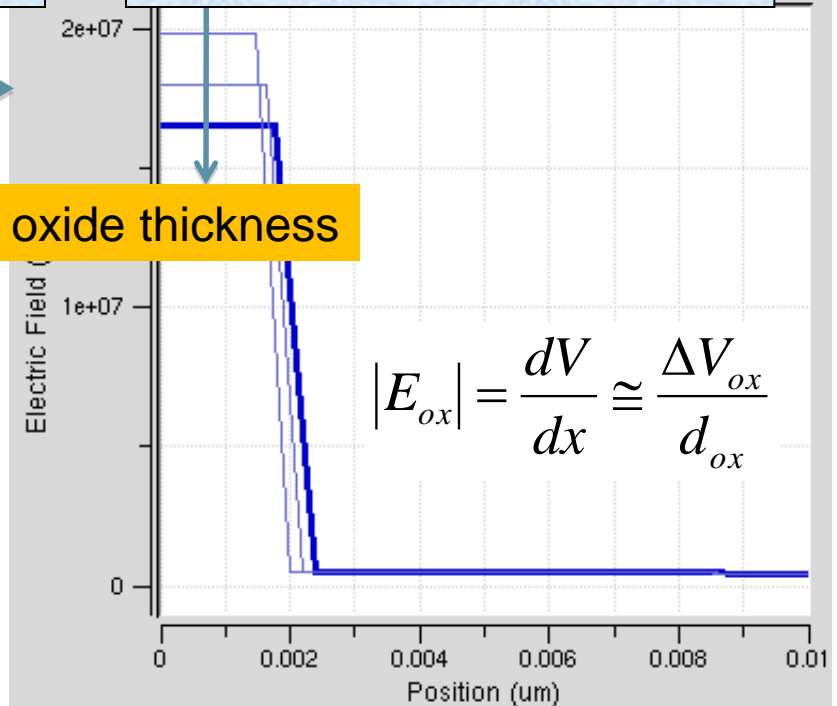
$C_{tot} / C_{ox} / C_S$: total/oxide/semiconductor capacitance
 $\epsilon_{ox} / \epsilon_S$: dielectric constant of oxide/semi conductor
 W_D : depletion width, d_{ox} : oxide thickness

d_{ox} increase \rightarrow C_{ox} decrease \rightarrow C_{tot} decrease



Increasing oxide thickness

d_{ox} increase \rightarrow E_{ox} decrease

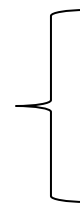


What is PADRE?: the simulator behind MOSCAP

PADRE^[6] : classical drift-diffusion simulator

- PADRE simulates the electrical behavior of devices under steady state, transient conditions or AC small-signal analysis^[7].

Self consistent simulation



Poisson equation

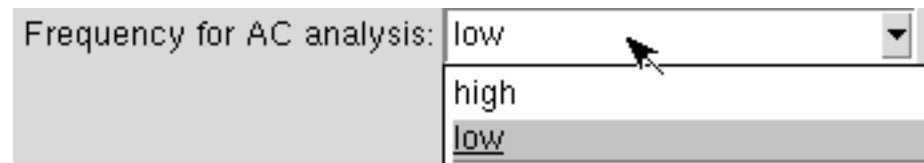
Drift-diffusion equation

[6] Dragica Vasileska; Gerhard Klimeck (2006), "Padre," DOI: 10254/nanohub-r941.3.

[7] http://nanohub.org/resource_files/tools/padre/doc/index.html

Limitations of the MOSCAP Tool

- MOSCAP does not model quantum mechanical effects
 - » Confinement of carriers in the semiconductor
 - » Tunneling from the gate to the semiconductor
- MOSCAP does not have sequence plots – users cannot see what happens in the intermediate steps between $V_G=0$ and the last applied bias*
 - Energy band diagram (at $V_G = 0$)
 - Energy band diagram (at last applied bias)
- MOSCAP does not allow users to choose the frequency for AC analysis*



* These features may be upgraded in the next versions.

References

- MOS Capacitor/MOSFET basic theory

[1] www.eas.asu.edu/~vasilesk, Dragica Vasileska (2008), "MOS Capacitors Description," <http://nanohub.org/resources/5087>

[2] www.eas.asu.edu/~vasilesk, Dragica Vasileska (2008), "MOSFET Operation Description," <https://nanohub.org/resources/5085>

[3] Mark Lundstrom (2008), "ECE 612 Lecture 3: MOS Capacitors," <http://nanohub.org/resources/5363>

- MOS Capacitor C-V measurements

[4] S.-H. Lo, et. al., IBM Journal of Research and Development, volume 43, number 3, 1999

- Charges in oxide

[5] Deal B. E., Electron Devices, IEEE Transactions on, 1980

- PADRE

[6] Dragica Vasileska; Gerhard Klimeck (2006), "Padre," [DOI: 10254/nanohub-r941.3](https://doi.org/10.2554/nanohub-r941.3).

[7] http://nanohub.org/resource_files/tools/padre/doc/index.html