

# **MOS FUNDAMENTALS**

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## **MOS Structure**

## <u>M</u>etal - <u>O</u>xide (SiO<sub>2</sub>) - <u>S</u>emiconductor (Si)





- The most common field plate (gate) materials are heavily doped polycrystalline silicon.
- The silicon-side terminal is called the back or substrate contact.
- The more general designation: metal-insulator-semiconductor (MIS)



Individual energy band diagrams for the metal, insulator, and semiconductor components.



## **Ideal Structure Assumption**

- (1) The metallic gate is sufficiently thick so that it can be considered an equipotential region.
- (2) The oxide is a perfect insulator.
- (3) No charge centers located in the oxide or at the interface.
- (4) Uniformly doped.



- (5) The semiconductor is sufficiently thick so that a field-free region("bulk") is encountered before reaching the back contact.
- (6) An ohmic contact between the semiconductor and the metal on the back side.
- (7) One-dimensional structure.
- (8) No work function difference between metal and semiconductor.





Energy band diagram of an ideal MOS structure with no bias.



# Effect Of An Applied Bias - Qualitative description

- General observations
  - $E_F(metal) E_F(semicondutor) = -qV_G$
  - With  $V_G \neq 0$ , semiconductor Fermi energy is unaffected by the bias and remains invariant as a function of position because of the assumed zero current flow.



- Since the barrier heights are fixed quantities, the movement of the metal Fermi level leads to a band bending.
  - In the metal, no bend-bending.
  - In the oxide and semiconductor, an upward slope when  $V_G > 0$ a downward slope when  $V_G < 0$ .
  - With no oxide charges, the Poisson's equation yields a constant slope in the oxide.
  - Band bending in the semiconductor is somewhat more complex.



• Specific biasing regions - Accumulation ( $V_c < 0$ )



+0

 $V_{G} < 0$ 

-0

• VG < 0 raises EF(metal) and the hole concentration inside the semiconductor,  $p = n_i \exp[-(E_F - E_i)/kT]$ increases as one approaches the oxide-semiconductor interface.

• VG < 0 places negative charges on the gate. To maintain a balance of charge, positively charged holes must be drawn toward the Si-SiO2 interface.





Depletion

• VG > 0 slightly lowers EF(metal) and the hole concentration is decreased (depleted) in the vicinity of the Si-SiO2 interface.

• VG > 0 places positive charges on the gate, which in turn repels holes from the interface and exposes the negatively charged acceptor sites.



#### - Onset of Inversion $(V_{c} = V_{T})$



• As VG is increased positively, the bands at the Si surface will bend down more and the electron concentration at the surface (ns) will increase from less than ni when E i (surface) > EF, to ni when Ei (surface) = EF, to greater than ni when Ei (surface) < EF.



 $V_G = V_T$ 

$$E_{i}(bulk) - E_{i}(surfac) = 2[E_{i}(bulk) - E_{F}]$$

$$n_{s} = n_{i} \exp\left[\frac{E_{F} - E_{i}(surfac)}{kT}\right]$$

$$= n_{i} \exp\left[\frac{E_{i}(bulk) - E_{F}}{kT}\right] = p_{bulk} = N_{A}$$

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#### - Inversion (VG > VT)





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No Bias (V_G = 0)
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Accumulation (V_G < 0)
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Depletion (  $0 < V_G < V_T$  )





**Semiconductor Device Fundamentals** 



# Effect Of An Applied Bias - Quantitative formulation

• Preparatory considerations



For an p-type semiconductor,

<ul> <li>Accumulation</li> </ul>	•	$\phi_s < 0$
– Flat band	•	$\phi_s = 0$
– Depletion	•	$0 < \phi_s < 2\phi_F$
– Onset of inversion :		$\phi_{s}=2\phi_{F}$
– Inversion	•	$\phi_{s} > 2\phi_{F}$



# • Delta-depletion solution

## **Delta-depletion assumption :**

- The functional form of the accumulation charge & the inversion charge :  $\delta$  function.
- Because the depletion width increases only slightly once the semiconductor inverts, it is assumed the  $\delta$  function of charge added in inversion precisely balances the charge added to the gate.
- The actual depletion charge is replaced with a squared-off distribution



- Accumulation :
  - Because of the assumed  $\delta$  function, the electric field and potential are zero for all x > 0.
- Depletion :

For 
$$0 \le x \le W$$
,  

$$\frac{dE}{dx} = \frac{\rho}{K_s \varepsilon_0} \cong -\frac{qN_A}{K_s \varepsilon_0}$$

$$E(x) = \frac{qN_A}{K_s \varepsilon_0} (W - x)$$

$$\phi(x) = \frac{qN_A}{2K_s \varepsilon_0} (W - x)^2$$

with 
$$\phi = \phi_s$$
 at  $x = 0$ ,  
 $W = \left[\frac{2K_s \varepsilon_0}{qN_A}\phi_s\right]^{1/2}$ 

The maximum depletion width,

$$W_{T} = \left[\frac{2K_{S}\varepsilon_{0}}{qN_{A}}(2\phi_{F})\right]^{1/2}$$







- Inversion :
  - The solution is established by merely adding

     a δ function of surface charge to the
     solution existing at the end of depletion.
  - The depletion charge, the x > 0 electric field,
     and the x > 0 potential remain

fixed at their values.

$$\phi_s = 2\phi_F$$



# Exact solution for the charge density and potential assuming

### $\phi_F = 12kT/q$ and T = 300K

(a) Accumulation ( $f_s = -6kT/q$ )



## (b) Middle of depletion ( $f_s = f_F = 12kT/q$ )



(c) Onset of inversion ( $f_s = 2f_F = 24kT/q$ )



#### (d) Deep into inversion





• Gate voltage relationship (delta-depletion solution)

$$V_{G} = \Delta \phi_{Semi} + \Delta \phi_{OX}$$

Because  $\varepsilon_{OX}$  is constant in an ideal oxide with no charges,

$$\Delta \phi_{\mathbf{ox}} = x_o \, \varepsilon_{\mathbf{ox}}$$

Since there is no charges at the interface, (in the depletion region)

$$D_{ox} = D_{Semi} \Big|_{\mathbf{X}=\mathbf{0}}$$
$$E_{ox} = \frac{K_s}{K_o} E_s$$

$$V_G = \phi_S + \frac{K_s}{K_o} \mathbf{x}_o E_s = \phi_S + \frac{K_s}{K_o} \mathbf{x}_o \sqrt{\frac{2qN_A}{K_s\varepsilon_0}} \phi_S \quad (0 \le \phi_S \le 2\phi_F)$$

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At threshold, 
$$\phi_s = 2\phi_F$$
  
$$V_T = 2\phi_F + \frac{K_s x_o}{K_o} \sqrt{\frac{4qN_A}{K_s \varepsilon_0}} \phi_F$$





 $\phi_s$  is a rather rapidly varying function of V<sub>G</sub> when the device is in depletion. However, when it is accumulated or inverted, it takes a large change in V<sub>G</sub> to produce a small change in  $\phi_s$ .

; +++++ delta-depletion solution, —— exact solution.

