Overview of An Advanced Surface-Potential-Based Model (SP)

(Invited paper)

G. Gildenblat and T.-L. Chen

Department of Electrical Engineering
The Pennsylvania State University
University Park, PA 16802, USA

Gildenblat@psu.edu

ABSTRACT

This paper outlines a new surface-potential-based compact MOSFET model (SP) developed at The Pennsylvania State University. The main objective of this work is to find practical engineering solutions of several long-standing problems of surface-potential-based modeling and to use them as a basis for a comprehensive compact MOSFET model. As a result of this approach the physical content of the model is significantly increased while the number of model parameters is reduced without sacrificing the quality of the fit of experimental data.

Keywords: Compact MOSFET model; surface potential.

1 INTRODUCTION

The aggressive reduction of the power supply voltage implies that popular threshold-voltage-based approach to compact modeling is close to its limits and the next generation advanced compact MOSFET models should be surface-potential-based. The challenge is to develop practical and efficient surface-potential-based models which do not suffer from the limitations traditionally associated with this approach. The novel features of SP include

• Analytical (non-iterative) evaluation of the surface potential \( \phi_s \) [1].
• Symmetric bulk charge linearization [2].
• Inclusion of the velocity saturation in a form consistent with Gummel symmetry requirements.
• Bias-dependent lateral field gradient [3].
• Mobility model incorporating Coulomb scattering.
• Analytical approximation for the quantum corrections and polysilicon depletion effect based on the surface potential method [4,5].

The additional features of SP include very simple and physical charge model which is consistent with the drain current model and is free from the unphysical behavior often associated with the more traditional threshold-voltage based models. In what follows we outline the approach taken in the development of SP.

2 SURFACE POTENTIAL

Iterative computations of surface potential represent a traditional weak point of the surface-potential-based models. Analytical approximation for the surface potential that is accurate enough for evaluation of transconductances and transcapacitances for \( \phi_s \geq 3 \phi_i \) \((\phi_i = k_B T / q)\) was presented in [1]. As shown in Fig. 1 this work can be extended to cover all regions of MOSFET operation including accumulation region. The resulting approximation has a maximum absolute error less than 10 nV and is used in SP.

3 LATERAL FIELD GRADIENT

An efficient way to introduce small-geometry effects into a surface-potential-based model is via lateral gradient factor

\[
f = 1 - (\varepsilon_s / q N_{sub}) \frac{\partial^2 \phi_s}{\partial y^2}
\]

Here symbols \( \varepsilon_s \) an \( q \) have their usual meaning and \( y \) is counted from source to drain. The “effective” substrate doping level \( N'_{sub} = f N_{sub} \) is reduced as compared with the actual impurity concentration \( N_{sub} \). This is a physically meaningful alternative to the “charge sharing” approach to the threshold voltage reduction in short-channel devices. In the gradual channel approximation \( f = 1 \), while in more recent work \( f = f(L,W) \) is geometry dependent [6]. SP generalizes this approach to include the bias dependence of \( f \) via a semi-empirical expression

\[
f = f_0 (1 + B_f \phi_f)
\]

where \( f_0 \) is bias dependent, \( B_f \) is a local (i.e. scalable) model parameter and \( \phi_f \) is a surface potential selected at a
suitable point in the channel to assure Gummel symmetry. The resulting increase of $f$ with $V_{gs}$ predicted by (2) is based on a simple physical model and is shown in Fig. 2.

4 MOBILITY MODEL

The effective mobility ($\mu$) model used in SP includes both the universal dependence on the effective vertical field $E_{eff}$ and the deviation from universality associated with Coulomb scattering:

$$\mu = \frac{MU0}{1 + (\mu_E E_{eff})^{\theta_E} + CS [q_b/(q_i + q_b)]^2}$$

(3)

Here $MU0$ and $CS$ are global (non-scalable) model parameters, $\mu_E$ and $\theta_E$ are local parameters and the last term in the denominator introduces Coulomb scattering [7]. The inversion charge $q_i$, the bulk charge $q_b$ and $E_{eff}$ are computed at the potential midpoint in order to assure Gummel symmetry of the model [2]. Here and below “potential midpoint” is defined as a point along the channel where $\phi_s = (1/2)(\phi_{ss} + \phi_{sd}) \Delta \phi_m$. The inclusion of the Coulomb scattering term is required primarily to assure a good fit of low-temperature experimental data shown in Fig. 3 without scarifying the quality of the fits at room and high temperatures (cf. Figs. 4,5).

5 SYMMETRIC LINEARIZATION

Most compact models are developed using bulk and inversion charge linearization. Linearization in terms of $\phi_s - \phi_{ss}$ ($\phi_{ss}$ and $\phi_{sd}$ are surface potentials at the source and drain end of the channel respectively) results in the loss of Gummel symmetry and produces unphysical results of the type “$C_{gs} \neq C_{gd}$ for $V_{ds} = 0$” [8,9]. An alternative approach called “symmetric linearization” was developed in [2] where it was verified using exact results for long-channel MOSFET’s. A small-geometry version of this technique was developed and serves as a basis of SP. In this approach

$$I_d = \beta(q_{im} + \alpha \phi_f) \phi/(r_L + \delta \phi/V_c)$$

(4)

where $\beta = \mu(W/L)C_{ox}$, $q_{im}$ is the normalized inversion charge evaluated at the potential midpoint and $\alpha = -(dq_i/d\phi)_{\phi=\phi_m}$. In (4) $r_L = (1 + \Delta L_{CLM} / L)^{-1}$

where $\Delta L_{CLM}$ is used to introduce channel length

![Fig. 1: Comparison of analytical approximation (solid line) and iterative solution (circles) for the surface potential; $t_{ox} = 20 \text{ Å}, N_{sub} = 10^{18} \text{ cm}^{-3}, V_{fb} = -1 \text{ V and } V_{sb} = 0.$](image1)

![Fig. 2: Typical lateral gradient factor for long-channel (L=20 µm) and short-channel (L=0.2 µm) devices; $t_{ox} = 40 \text{ Å}, N_{sub} = 2.4 \times 10^{17} \text{ cm}^{-3}, V_{ds} = 1 \text{ V and } V_{sb} = 0.$](image2)

![Fig. 3: Measured (circles) and modeled (solid lines) linear transconductance for $T=218 \text{ K}; W/L=10/0.2 \mu m, t_{ox} = 40 \text{ Å}, N_{sub} = 2.4 \times 10^{17} \text{ cm}^{-3}, V_{ds} = 0.1 \text{ V and } V_{sb} = 0 \sim 2 \text{ V.}](image3)
modulation and $V_c = LV_s / \mu$ where $V_s$ denotes saturation velocity. Finally, Grotjohn-Hoefflinger factor $\delta_0 = \delta_0 (\phi)$ serves two purposes: sharpening velocity-field dependence [10, 11] and eliminating the singularity at $V_{ds} = 0$ which is present if one sets $\delta_0 = 1$.

The effective surface potential difference $\phi$ is obtained by softly clamping $\phi_{sat} - \phi_{ss}$ in order to eliminate the negative $G_{ds}$ region which appears when velocity saturation is introduced directly into a charge-sheet model. MOSFET characteristics based on eq. (4) satisfy Gummel symmetry test (cf. Figs. 6, 7) and other benchmark tests.

6 POLYSILICON DEPLETION AND QUANTUM CORRECTION

Polysilicon depletion has been traditionally modeled by changing the effective gate drive and the threshold voltage under an additional assumption of a pinned surface potential. This approach is well suited for the threshold-voltage-based models but produces unphysical results in a subthreshold region [13]. In SP approach surface potential

$$\phi_s = \phi_s^{(0)} + \Delta \phi_s$$

where $\phi_s^{(0)}$ is the surface potential in the absence of polysilicon depletion (cf. section 1 above) and $\Delta \phi_s$ is the perturbation of the surface potential caused by the polysilicon depletion layer. An approximation for $\Delta \phi_s$ developed in [4] is sufficiently accurate for the purposes of compact modeling (Fig. 8). In addition to changing $\phi_s$, the effect of polysilicon depletion is to modify bulk charge linearization coefficient used in (4):

$$\alpha = \alpha^{(0)} + \eta_p - 1$$

where

$$\eta_p = \left[ 1 + (2C_{ox}^2 / q \varepsilon_s N_p) \left[ V_{gs} - V_{fb} - \phi_m^{(0)} \right] \right]^{-1/2}$$

$N_p$ denotes the polysilicon doping level, $\alpha^{(0)}$ and $\phi_m^{(0)}$ represent $\alpha$ and $\phi_m$ respectively in the absence of polysilicon depletion effect.

The quantum effects are considered similarly with the analytical approximation for the perturbation of the surface potential given in [5]. When used in SP the equations developed in [4, 5] are introduced in a manner consistent with the Gummel symmetry requirement.
7 INTRINSIC CHARGE MODEL

The simplicity of the symmetric linearization approach allows one to obtain analytical expressions for the terminal charges in a form consistent with the drain current model. For example,

\[
\frac{Q_c}{WLC_{ox}} = V_{gb} - V_{fb} - \phi_m + \frac{\eta_p \phi}{2} \left( \frac{r_l \phi}{6H} - 1 + r_l \right) \tag{8}
\]

and (using Ward-Dutton partition [14])

\[
\left| \frac{Q_D}{WLC_{ox}} \right| = \frac{q_{im}}{2} - \frac{\alpha \phi (1 - r_l^2)}{4} - \frac{r_l^2 (1 - \eta_p) \phi^2}{24H}
\]

\[
- \frac{r_l \phi (\alpha + 1 - \eta_p)}{12} \left( 1 - \frac{\phi}{2H} - \frac{\phi^2}{20H^2} \right) \tag{9}
\]

where

\[
H = \frac{(q_{im}/\alpha) + \phi_l}{1 + \delta_0 \phi / r_l V_c} \tag{10}
\]

The resulting C-V curves shown in Figs. 9 and 10 are smooth, physical and symmetric. In particular, \( C_{gs} = C_{gd} \) for \( V_{ds} = 0 \).

7 EXAMPLES OF DEVICE CHARACTERISTICS

The verification of SP was accomplished using experimental data for a variety of fabrication processes from several semiconductor companies. Typical results are shown in Figs. 3-5, 11 and 12. Local (fixed \( L/W \)) version of SP requires up to 28 parameters while global version (including scaling) contains 68 parameters of which only 35 are needed for a typical (mature) fabrication process. The relatively small number of the fitting parameters is a consequence of the increased physical content of the model. We note in passing that global version of SP includes reverse short-channel effect via geometry dependences of the doping concentration and of the flat-band voltage.

8 CONCLUSIONS

We have outlined the core SP model consisting of a drain current model and quasi-static model of the terminal charges. The extrinsic SP model will be presented separately. The results of this work indicate that surface-potential-based models are both physical and practical and can form the foundation of the next generation of compact MOSFET models.
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