Simulating Vertical GaN FinFET in Sentaurus

Alexander Simko

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Abstract

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1 Introduction

2 Theory

2.1 Mobility

Mobility describes how easily charge carriers can move through a semiconductor. It is related to the resistivity of the semiconductor by:

$$\rho = \frac{1}{q\mu N} \tag{1}$$

where μ is the mobility, N is the carrier concentration and ρ is the resistivity. The mobility depends on many parameters such as temperature, doping level, electric field and surface scattering. [1]

Multiple models for the mobility can be used and combined using Matthiessen's rule:

$$\frac{1}{\mu} = \frac{1}{\mu_1} + \frac{1}{\mu_2} + \dots \tag{2}$$

By modeling different mobility dependances separately they can then be combined using eq. (2).

2.1.1 Doping Dependance

Experimental data shows that mobility decreases with increasing doping. One simple model for this is the Arora model given by:

$$\mu_{dop} = \mu_{min} + \frac{\mu_d}{1 + \left(\frac{N_A + N_D}{N_0}\right)^A} \tag{3}$$

where N_A and N_D is the doping concentration and μ_{min} , μ_d , N_0 and A are fitting parameters. In order to account for temperature dependance every fitting parameter can be multiplied with a temperature dependant scaling factor, k_i

$$k_i = \left(\frac{T}{300 \,\mathrm{K}}\right)^{\alpha_i} \tag{4}$$

where T is the temperature in Kelvin and α_i is another unique fitting parameter for each parameter mentioned above. Combining eqs. (3) and (4) then yields:

$$\mu_{dop} = \mu_{min} \cdot \left(\frac{T}{300 \,\mathrm{K}}\right)^{\alpha_m} + \frac{\mu_d \cdot \left(\frac{T}{300 \,\mathrm{K}}\right)^{\alpha_d}}{1 + \left(\frac{N_A + N_D}{N_0 \cdot \left(\frac{T}{300 \,\mathrm{K}}\right)^{\alpha_N}}\right)^{A \cdot \left(\frac{T}{300 \,\mathrm{K}}\right)^{\alpha_A}}}$$
(5)

2.1.2 Electric Field Dependance

Velocity no longer scale proportionally to applied field at high field strength, leading to field dependant mobility.

2.1.3 Surface Scattering Dependance

Interfaces between semiconductor and oxide introduce defects. This leads to a mobility degradation close to the surface due to several scattering phenomena. These include surface roughness scattering, acoustic phonon scattering and Coulomb scattering due to charged traps at the interface. Surface roughness and acoustic phonon scattering are both described in the enhanced Lombardi model and depend strongly on the electric field normal to the interface $(E_{\perp}(x))$. The mobility degradation due to acoustic phonon scattering is given by:

$$\mu_{ac}(x) = \frac{B}{E_{\perp}(x)} + \frac{C \left(N_A + N_D\right)^{\lambda}}{E_{\perp}^{\frac{1}{3}}(x) \cdot \left(\frac{T}{200 \text{ K}}\right)^k} \tag{6}$$

and the degradation from surface roughness is given by:

$$\mu_{sr}(x) = \left(\frac{E_{\perp}^2(x)}{\delta} + \frac{E_{\perp}^3(x)}{\eta}\right)^{-1} \tag{7}$$

with B, C, λ , k, δ and η being fitting parameters and x being the distance from the interface. The mobilities from eqs. (6) and (7) can then be combined with any bulk mobility (μ_b) using eq. (2). However as these effects only occur at the surface they need to be weighted with a distance factor $D(x) = \exp(-x/l_{crit})$. l_{crit} is also a fitting parameter. This gives the final mobility:

$$\frac{1}{\mu} = \frac{1}{\mu_b} + \frac{D(x)}{\mu_{ac}(x)} + \frac{D(x)}{\mu_{sr}(x)} \tag{8}$$

The mobility degradation from Coulomb scattering, i.e. scattering from trap charges at the interface can be modeled using the following formula:

$$\mu_c(x) = \frac{\mu_1 \left(\frac{T}{300 \,\mathrm{K}}\right)^k \left[1 + \left(\frac{n}{c_{trans} \cdot \left(\frac{N_{p,i}}{N_0}\right)^{\eta_1}}\right)^{\nu}\right]}{\left(\frac{N_{p,i}}{N_0}\right)^{\eta_2} \cdot D(x) \cdot f(E_{\perp}(x))}$$
(9)

Coulomb scattering only occurs if the field normal to the interface is sufficiently large. Therefore the field dependant function $f(E_{\perp})$ is introduced:

$$f(E_{\perp}(x)) = 1 - \exp\left[-\left(\frac{E_{\perp}(x)}{E_0}\right)^{\gamma}\right] \tag{10}$$

In eqs. (9) and (10) μ_1 , k, ν , c_{trans} , N_0 , η_1 , η_2 , γ and E_0 are fitting parameters and T is the temperature, n is the electron charge density, $N_{p,i}$ is the positive interface charge density and D(x) is the above mentioned distance factor.

2.1.4 Effective Mobility

As many of the above mentioned mobility models depend on the position it can be hard to measure these in physical devices. Therefore the concept of effective mobility is very useful. The effective mobility is given by [1]:

$$\mu_e = \frac{\int_{-x_i}^{x_i} \mu(x) n(x) dx}{\int_{-x_i}^{x_i} n(x) dx}$$
 (11)

This can be interpreted as a weighted average of the mobility, weighted by the position of the actual charge carries, meaning that areas without or with very few mobile charges does not contribute to the overall effective mobility and areas with a high charge carrier concentration have a large contribution.

Another way of calculating the effective channel mobility is by using the definition of mobility. The final formula for the mobility then becomes [2]:

$$\mu_e = \frac{g_d L_g}{L_w Q_n} \tag{12}$$

3 Simulations

Different simulations were created to simulate the different behaviors of the transistor and to calibrate the physics model. This section describes the different simulations.

3.1 Standalone Fin Simulation

The first simulation created was the standalone fin simulation as seen in fig. 1. This was used to calibrate the interface mobility models described in section 2.1. The effective mobility of the fin was calibrated to match the measured effective mobility in [2].

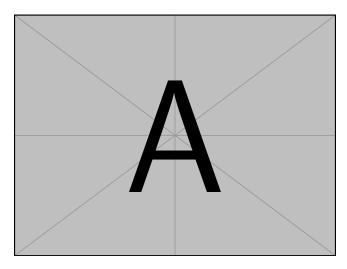


Figure 1: Structure of fin simulation.

3.2 Full Device Simulations

Using the calibrated mobility models the full device, as seen in fig. 2, could be simulated. The first simulation was a V_g sweep. Using this simulation the source contact resistance was calibrated.

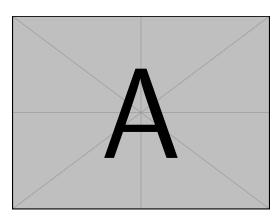


Figure 2: Full device structure

4 Result

5 Discussion

References

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- [2] P Gribisch et al. "Capacitance and Mobility Evaluation for Normally-Off Fully-Vertical GaN FinFETs". In: *IEEE Transactions on Electron Devices* 70.8 (2023), pp. 4101–4107. ISSN: 1557-9646 VO 70. DOI: 10.1109/TED.2023.3287820.