University of Massachusetts Amherst

From the SelectedWorks of Eric Polizzi

July, 2005

A Theoretical Investigation of Surface Roughness Scattering in Silicon Nanowire Transistors

Eric Polizzi, *University of Massachusetts - Amherst* M. Lundstrom S. Datta A. Ghosh J. Wang



Available at: https://works.bepress.com/eric_polizzi/6/



Jing Wang, Eric Polizzi, Avik Ghosh, Supriyo Datta, and Mark Lundstrom

Applied Physics

Letters

Citation: Appl. Phys. Lett. **87**, 043101 (2005); doi: 10.1063/1.2001158 View online: http://dx.doi.org/10.1063/1.2001158 View Table of Contents: http://apl.aip.org/resource/1/APPLAB/v87/i4 Published by the American Institute of Physics.

Additional information on Appl. Phys. Lett.

Journal Homepage: http://apl.aip.org/ Journal Information: http://apl.aip.org/about/about_the_journal Top downloads: http://apl.aip.org/features/most_downloaded Information for Authors: http://apl.aip.org/authors

ADVERTISEMENT



Theoretical investigation of surface roughness scattering in silicon nanowire transistors

Jing Wang

School of Electrical and Computer Engineering, Purdue University, West Lafayette, Indiana 47907

Eric Polizzi

Department of Computer Sciences, Purdue University, West Lafayette, Indiana 47907

Avik Ghosh, Supriyo Datta, and Mark Lundstrom

School of Electrical and Computer Engineering, Purdue University, West Lafayette, Indiana 47907

(Received 8 February 2005; accepted 13 June 2005; published online 19 July 2005)

Using a full three-dimensional (3D), quantum transport simulator, we theoretically investigate the effects of surface roughness scattering (SRS) on the device characteristics of Si nanowire transistors (SNWTs). The microscopic structure of the Si/SiO₂ interface roughness is directly treated by using a 3D finite element technique. The results show that (1) SRS reduces the electron density of states in the channel, which increases the SNWT threshold voltage, and (2) the SRS in SNWTs becomes less effective when fewer propagating modes are occupied, which implies that SRS is less important in small-diameter SNWTs with few modes conducting than in planar metal-oxide-semiconductor field-effect-transistors with many transverse modes occupied. © 2005 American Institute of Physics. [DOI: 10.1063/1.2001158]

The silicon nanowire transistor (SNWT) is attracting broad attention as a promising structure for future electronics.^{1,2} Therefore, understanding carrier transport in Si nanowires becomes increasingly important. Careful studies are needed to experimentally explore transport in SNWTs, but it is also clear that a theoretical understanding is similarly important. In this letter, we present a theoretical exploration of the Si/SiO₂ interface roughness scattering, or surface roughness scattering (SRS),^{3–5} in SNWTs.

It is well-known that scattering due to Si/SiO₂ interface roughness is important in planar silicon metal-oxidesemiconductor field-effect transistors (MOSFETs), and it is expected to be even more important in ultrathin-body silicon-on-insulator (UTBSOI) MOSFETs.3 For bulk MOS-FETs, electrons are confined at the Si/SiO₂ interface by an electrostatic potential well. Under high gate bias, the potential well is thin, electrons are confined very near the interface, SRS increases, and the effective mobility decreases. For UTBSOI MOSFETs, the confining potential is determined by the film thickness, and SRS can be enhanced by the roughness at the two interfaces.³ In a SNWT, the channel is surrounded by the Si/SiO₂ interfaces, so one might expect SRS to dominate transport. We will show, however, that SRS may be less important in SNWTs than in planar devices because of the one-dimensional (1D) nature of the SNWT channel.

In Ref. 6, we developed a full three-dimensional (3D), quantum transport simulator of SNWTs based on the effective-mass approximation. In this work, to investigate the effects of SRS on small-diameter (\sim 3 nm) SNWTs with physically rough Si/SiO₂ interfaces, we make use of this previously developed simulator. The simulated structure is a gate-all-around SNWT with a rectangular cross section and a [100] oriented channel (see Fig. 1). Following previous work on SRS,^{3–5} we assume an abrupt, randomly varying interface between the Si and SiO₂, parametrized by a root mean square (rms) amplitude and an autocovariance function.^{7,8} The statistical nature of the roughness will depend on the nanowire fabrication methods and may differ considerably from that arising during the high temperature oxidation of a planar Si surface. Nevertheless, since our objective is to discuss general insights into the physics of SRS in SNWTs, we will employ the roughness parameters for a planar (100) Si/SiO₂ interface obtained from Ref. 7. Our use of a continuum level description may be questioned, but we believe that it is a

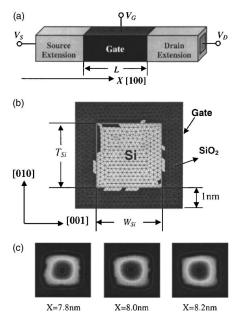


FIG. 1. (a) Schematic diagram of the simulated gate-all-round SNWT. The source/drain doping concentration is 2×10^{20} cm⁻³ and the channel is undoped. There is no source/drain overlap with the channel and the gate length is L=10 nm. V_S , V_D , and V_G are the applied voltage biases on the source, drain, and gate, respectively; (b) cross section of the SNWT with a specific interface roughness pattern for the slice at X=9.0 nm. For the device with smooth Si/SiO₂ interfaces, the Si body thickness is $T_{Si}=3$ nm, the wire width is $W_{Si}=3$ nm, and the oxide thickness is 1 nm; (c) confined wave functions for the slices at X=7.8 nm, X=8.0 nm, and X=8.2 nm, respectively.

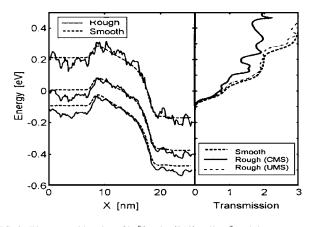


FIG. 2. Electron subband profile [for the (010) valleys] and the corresponding transmission coefficients for the simulated SNWT ($T_{\rm Si}=W_{\rm Si}=3$ nm) with smooth and rough Si/SiO₂ interfaces. The roughness parameters used are $L_m=0.7$ nm and rms=0.14 nm (Ref. 7). The device is at the ON-state ($V_{\rm CS}=V_{\rm DS}=0.4$ V), so the source and drain Fermi levels are equal to 0 eV and -0.4 eV, respectively.

useful first step that gives insight into how the magnitude and spatial coherence of potential fluctuations influence carrier transport. In contrast to previous work,^{3–5} which made use of perturbation theory to compute the surface roughness scattering rate, we treat the physically rough structure directly.

The microscopic structure of the Si/SiO₂ interface roughness is implemented into the 3D simulator according to the following procedure. We first discretize the simulation domain with a 3D finite element mesh;⁶ each element is a triangular prism with a 2 Å height and edge length, comparable to the size of roughness at the (100) Si/SiO₂ interface.⁷ Next, we generate a two-dimensional (2D) random distribution across the whole Si/SiO₂ interface (unfolding the four interfacial planes into a sheet) according to an exponential autocovariance function;⁷

$$C(x) = \Delta_m^2 e^{-\sqrt{2}x/L_m},\tag{1}$$

where L_m is the correlation length, Δ_m is the rms fluctuation of the roughness, and x is the distance between two sampling points at the interface. Based on the 2D random distribution, the types of the elements at the Si/SiO₂ interfaces may be changed from Si to SiO₂ or reversely, to mimic the rough interfaces [see Fig. 1(b)].

After the roughness is implemented, electron transport through the rough SNWT is simulated by using the nonequilibrium Green's function approach.⁹ With a coupled mode space (CMS) representation, $^{\delta,10-12}$ the wave function deformation due to the roughness is treated. (The simulation methodology has been described in detail in Refs. 6 and 10.) To emphasize the role of SRS on electron transport, we do not include any other scattering mechanisms, so coherent transport is assumed inside the device. (Oscillations in the current due to quantum interference might be expected, but the averaging over a thermal distribution of wavelengths that occurs is sufficient to suppress them.) The length of the channel (10 nm) is long enough to ensure that sufficient averaging takes place so that sample specific effects are not observed. The simulated results for the rough SNWT are then compared with those for a device with the same geometrical parameters (e.g., nominal oxide thickness and Si

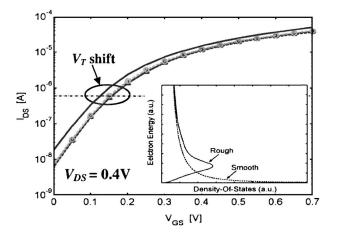


FIG. 3. $I_{\rm DS}$ vs $V_{\rm GS}$ curves for the simulated SNWT ($T_{\rm Si}=W_{\rm Si}=3$ nm) with smooth (solid) and rough (dashed with symbols) Si/SiO₂ interfaces. ($V_{\rm DS}$ =0.4 V). Three samples (triangles, crosses, and circles) of the rough SNWT are generated based on the same roughness parameters (L_m =0.7 nm and rms=0.14 nm) but different random number seeds. The SNWT threshold voltage (V_T) is defined as $I_{\rm DS}$ ($V_{\rm GS}=V_T, V_{\rm DS}=0.4$ V)=2×10⁻⁷ $W_{\rm Si}(A)$, where $W_{\rm Si}$ is in nm. The inset illustrates the reduction of electron DOS at low injection energies caused by SRS.

body thickness) but smooth Si/SiO₂ interfaces. By doing this, the effects of SRS on SNWT device characteristics can be clearly identified.

Figure 2 plots the electron subband profile (left column) at the ON-state ($V_{GS} = V_{DS} = 0.4$ V) in the simulated SNWT with rough and smooth Si/SiO₂ interfaces. The corresponding transmission coefficients (right column) for both the rough and smooth SNWTs are also shown. Note that the modes are coupled in the simulation; we show them separately for illustrative purposes only. It is clearly seen in the energy versus X plot that the presence of the roughness introduces significant fluctuations in the electron subbands, which lead to fluctuating elements in the diagonal terms of the device Hamiltonian [for details, see Eq. (7) in Ref. 6] and act as a scattering potential. At the same time, the shape of the confined wave function also varies from slice to slice along the wire in the rough SNWT [see Fig. 1(c) for an example], which produces deformation and coupling elements in both diagonal and off-diagonal terms of the device Hamiltonian [for details, see Eqs. (7), (8b), and (8c) in Ref. 6], and consequently lowers the transmission. (This effect has been named "wavefunction deformation scattering."^{13–15}) To examine the significance of wave function deformation scattering, we plot an energy versus transmission curve (dotdashed) for the rough SNWT calculated by the uncoupled mode space (UMS) approach,⁶ in which only the variations in the electron subbands are included while the deformation and coupling terms are discarded. The fact that the UMS approach significantly overestimates the transmission for the rough device infers that wave function deformation scattering dominates the transport. This is an important finding because common perturbation theory treatments³⁻⁵ of SRS scattering typically treat the subband energy fluctuations but not the wave function deformation scattering.

From the energy versus transmission plot, we find that the difference between the transmission curve for the rough SNWT and that for the smooth device becomes more and more noticeable as energy increases. This occurs because as energy increases, more subbands (modes) become conductive and the coupling between different modes efficiently re-

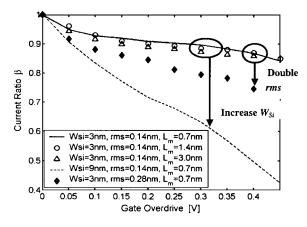


FIG. 4. Current ratio β vs gate overdrive curves for the simulated SNWTs with different wire widths ($W_{\rm Si}$) and roughness parameters (L_m and rms). At all the cases, the Si body thickness is fixed to be $T_{\rm Si}$ =3 nm and the drain bias is $V_{\rm DS}$ =0.4 V.

duces the transmission in the rough SNWT. In other words, SRS becomes more significant as more modes conduct. As we will show later, this effect has an important impact on the role of SRS on SNWT device characteristics.

Figure 3 plots the $I_{\rm DS}$ vs $V_{\rm GS}$ curves in a semilogarithmic scale for both the rough and smooth SNWTs. The results show that there is a distinct threshold voltage (V_T) increase caused by SRS. In the low gate bias region, the lateral displacement of the smooth and rough characteristics implies a V_T increment of ~30 mV for the roughness parameters we used ($L_m = 0.7$ nm and rms = 0.14 nm) and varies little from sample to sample. The increase in V_T due to SRS was unexpected and the reason for it is as follows. Due to SRS, injections at low energies are blocked in the rough SNWT, which reduces the density-of-states (DOS) near the band edge (see the Fig. 3 inset). The lowered DOS near the band edge reduces the charge density in the subthreshold regime, and consequently increases V_T in the rough SNWT. This effect would be modest in a conventional MOSFET with an energy-independent DOS above the band edge, but it becomes pronounced in a 1D wire with a singularity in the DOS at the band edge.

Finally, we explore the effects of SRS on the SNWT drain current above threshold. To do this, we compute a current ratio $\beta = I_{DS}^{rough}/I_{DS}^{smooth}$ at the *same* gate overdrive, $V_{GS}-V_T$, for both rough and smooth SNWTs. By comparing currents (rough versus smooth) at the same gate overdrive, the effect of the V_T increasing induced by SRS is removed. This allows us to examine whether the roughness can cause a significant reduction of SNWT ON-current by backscattering. Figure 4 shows the β vs gate overdrive curves for the SNWTs with different wire widths and roughness parameters. Several interesting phenomena are observed. First, all the simulated structures display a decreasing β with an increasing gate overdrive. This occurs because more modes become conductive under higher gate bias, which, as described earlier, enhances SRS in the SNWTs.

Second, based on the roughness parameters we used, $L_m=0.7 \text{ nm}$ and rms=0.14 nm, which are typical of an oxidized Si/SiO₂ interface,⁷ the SNWT with $W_{\text{Si}}=3 \text{ nm}$ (solid) achieves a surprisingly high $\beta \approx 0.9$ at a typical ON-state condition (gate overdrive =0.3 V for a 0.4 V supplied voltage). The same amount of surface roughness scattering severely degrades the mobility of a planar MOSFET under a high gate bias.¹³ To explore the effects of L_m , two additional values (1.4 nm for circles and 3.0 nm for triangles) were examined. The results show that β is insensitive to L_m , as expected from the averaging over a distribution of wavelengths that occurs at room temperature and high drain bias ($V_{\rm DS}$ =0.4 V). In contrast, doubling the rms (diamonds) clearly degrades β at the same gate overdrive, indicating the importance of maintaining relatively smooth Si/SiO₂ interfaces for the high performance application of SNWTs.

Third, increasing the wire width reduces the strength of quantum confinement and thus increases the number of conducting modes in the SNWT. Our results (solid versus dashed) clearly show that with a larger number of conducting modes in the wider ($W_{Si}=9$ nm) SNWT, SRS is much stronger than in the narrower ($W_{Si}=3$ nm) device. This observation also suggests that SRS is more serious in a planar MOS-FET, which can be viewed as a SNWT with a very large wire width.

In summary, we theoretically investigated SRS in SN-WTs by using a full 3D, self-consistent, quantum mechanical simulator. The microscopic structure of the Si/SiO₂ interface roughness was implemented into the simulator using the 3D finite element method. We found that (1) SRS reduces the electron density of states in the channel, which increases the SNWT threshold voltage, and (2) SRS in SNWTs becomes less serious when fewer propagating modes conduct, implying that SRS will be less important in small-diameter SN-WTs than in planar MOSFETs with many transverse modes occupied. This work provides important insights into the nature of SRS in SNWTs and suggests that SRS may not be as important in small-diameter nanowires as it is in conventional, planar MOSFETs.

This work was supported by the Semiconductor Research Corporation (SRC), and the Microelectronics Advanced Research Corporation (MARCO) Focus Center on Materials, Structures and Devices (MSD). The authors thank Jin Zhang, Min Chin Chai, Joseph Taylor, and Professor M. Alam at Purdue University for their sincere help.

- ¹Y. Cui, Z. Zhong, D. Wang, W. U. Wang, and C. M. Lieber, Nano Lett. **3**, 149 (2003).
- ²F. -L. Yang, D. -H. Lee, H. -Y. Chen, C. -Y. Chang, S. -D. Liu, C. -C. Huang, *et al.*, IEEE Symp. VLSI Tech., Digest of Tech. 2004, 196.
- ³F. Gamiz, J. B. Roldan, J. A. Lopez-Villanueva, P. Cartujo-Cassinello, and J. E. Carceller, J. Appl. Phys. **86**, 6854 (1999).
- ⁴M. V. Fischetti, Z. Ren, P. M. Solomon, M. Yang, and K. Rim, J. Appl. Phys. **94**, 1079 (2003).
- ⁵C. -Y. Mou and T. -M. Hong, Phys. Rev. B **61**, 12612 (2000).
- ⁶J. Wang, E. Polizzi, and M. Lundstrom, J. Appl. Phys. 96, 2192 (2004).
- [']S. M. Goodnick, D. K. Ferry, C. W. Wilmsen, Z. Liliental, D. Fathy, and O. L. Krivanck, Phys. Rev. B **32**, 8171 (1985).
- ⁸T. Yoshinobu, A. Iwamoto, and H. Iwasaki, Jpn. J. Appl. Phys., Part 1 26, 1447 (1994).
- ⁹S. Datta, Superlattices Microstruct. 28, 253 (2000).
- ¹⁰J. Wang, E. Polizzi, A. Ghosh, S. Datta, and M. Lundstrom, J. Comp. Electron. (in press).
- ¹¹E. Polizzi and N. B. Abdallah, Phys. Rev. B 66, 245301 (2002).
- ¹²E. Polizzi and N. B. Abdallah, J. Comput. Phys. 202, 150 (2004).
- ¹³K. Uchida, H. Watanabe, A. Kinoshita, J. Koga, T. Numata, and S. Takagi, Int. Electron. Dev. Meet. Tech. Digest 2003, 47.
- ¹⁴H. Sakaki, T. Noda, K. Hirakawa, M. Tanaka, and T. Matsusue, Appl. Phys. Lett. **51**, 1934 (1987).
- ¹⁵A. Gold, Phys. Rev. B **35**, 723 (1987).