

Crosslight Simulation of Effects of Bending in Thinfilm Transistors



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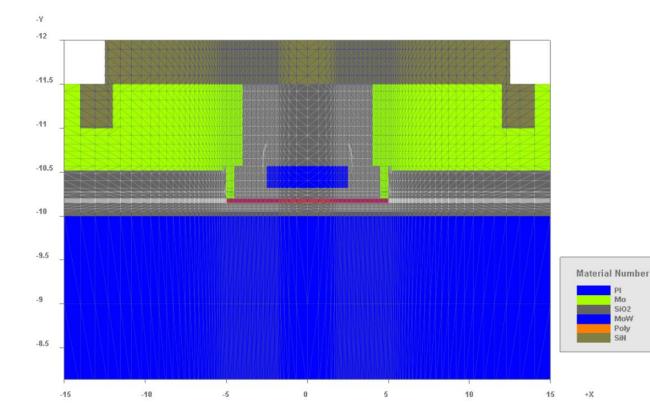
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- Structure and experiments
 - Stress and material models
 - Device model: Stressed metal
 - Device model: Stressed poly
 - Summary



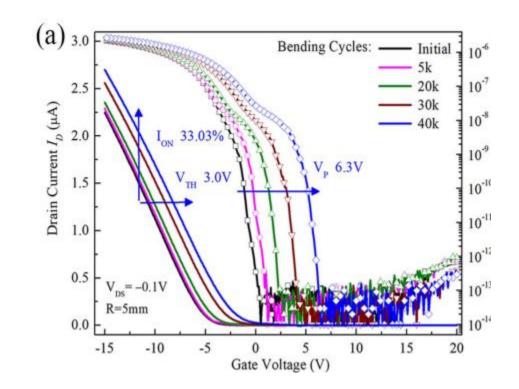
Structure



IEEE TRANSACTIONS ON ELECTRON DEVICES, VOL. 66, NO. 5, MAY 2019 p2214

Spontaneous Degradation of Flexible Poly-Si TFTs Subject to Dynamic Bending Stress

Wei Jiang[®], Mingxiang Wang[®], *Senior Member, IEEE*, Huaisheng Wang[®], and Dongli Zhang[®]





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Model

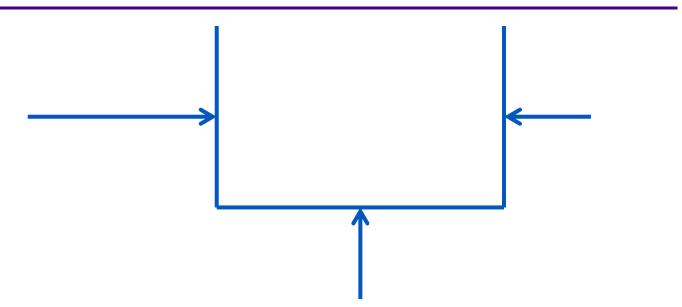
$$c_{11} * \frac{\partial^2 V_x}{\partial x^2} + c_{44} * \frac{\partial^2 V_x}{\partial y^2} + c_{12} * \frac{\partial^2 V_y}{\partial x \partial y} + c_{44} * \frac{\partial^2 V_y}{\partial y \partial x} = -bx$$

$$c_{12} * \frac{\partial^2 V_x}{\partial y \partial x} + c_{44} * \frac{\partial^2 V_x}{\partial x \partial y} + c_{44} * \frac{\partial^2 V_y}{\partial x^2} + c_{11} * \frac{\partial^2 V_y}{\partial y^2} = -by$$
V-displacement vector
bx/by=initial
stress/boundary stress
$$bx = \frac{\partial \sigma_{0,xx}}{\partial x} + \frac{\partial \sigma_{0,xy}}{\partial y}$$

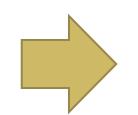
$$by = \frac{\partial \sigma_{0,yy}}{\partial y} + \frac{\partial \sigma_{0,xy}}{\partial x}$$



Strain boundary (default)



By default, boundaries in CSUPREM model are such that all displacement at left/right/bottom are restricted



Not suitable for bending



Revised Strain Boundary

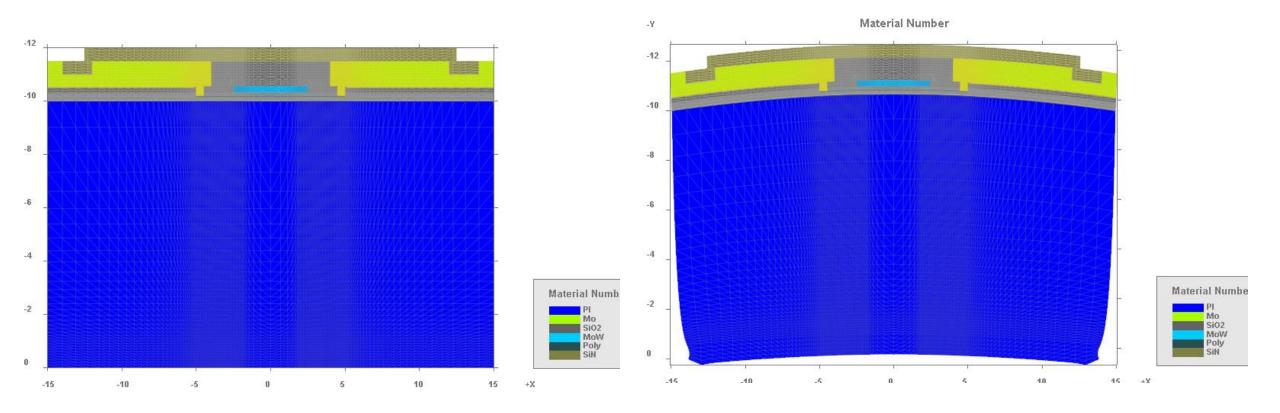
y.free.boundary=bottom Revised bottom boundary

x.free.boundary=left/right Revised left/right boundary

y.fixed.boundary=left/right Fix the y-direction so that the object would not move in y-direction after bending

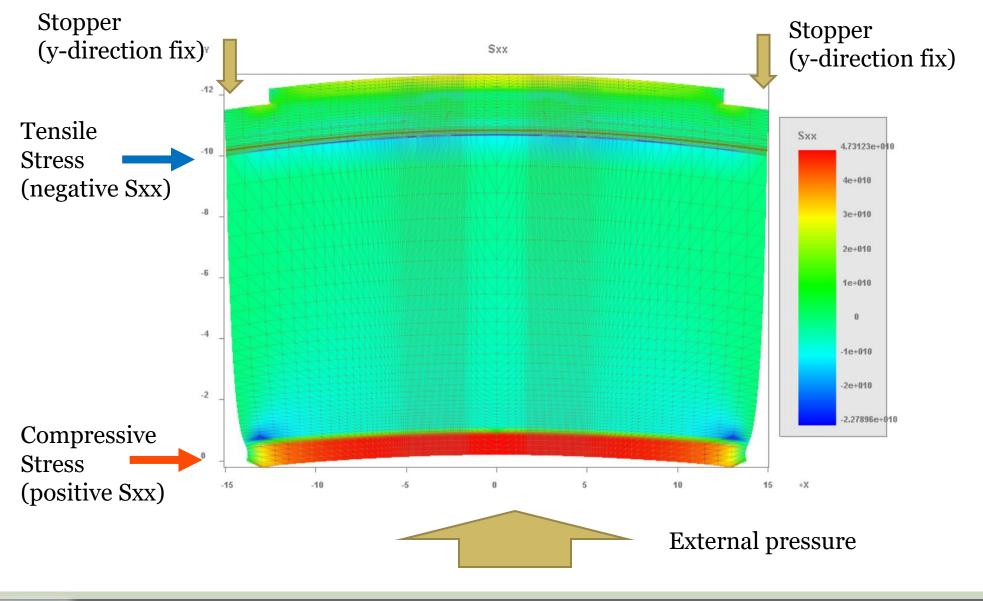


With and Without Bending





Stress Distribution





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Files and Commands

Ltps_TFT_30um_nostrain for nostrain

- (1) Run Csuprem on 2d.in
- (2) Run APSYS on tft2_vt.sol, tft2_iv.sol

(3) Customized MoW, Pi material macro defined as LTPS.mac Ltps_TFT_30um_strain for with strain

Let work function of MoW be stress dependent:

affinity variation=function function(stress_xx) 4.9+abs(stress_xx) end_function



Files and Commands

#Key commands in 2d.in

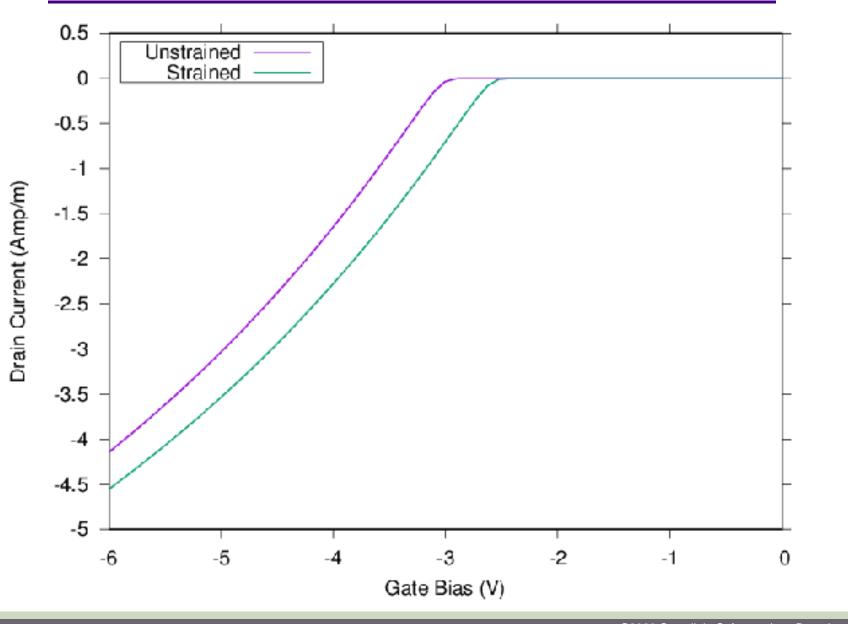
external_pressure xrange_from=-15.0 xrange_to=15.0 depth=0.5 sigma=-1.0e11 bottom stress temp1=25 temp2=25 y.free.boundary=bottom x.free.boundary=left/right y.fixed.boundary=left/right

struct outf=08_final_disp.str add.disp=t

External pressure acting from the bottom within a range. Depth affects mesh region where external pressure acts on. The results shall not be sensitive to choice of depth of external initial pressure.



Shift of Threshold Voltage





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\$ Main reference:

\$ (for basic band structure)

\$ [1] M.V. Fischetti and S.E. Laux, "Band structure, deformation

\$ potentials, and carrier mobility in strained Si,

\$ Ge, and SiGe alloys, J. Appl. Phys., vol. 80, pp. 2234-2252, 1996.

\$

\$ [2] "Electronic-band parameters in strained Si(1-x)Ge(x)
alloys on

\$ Si(1-y)Ge(y) substrates," M. M. Rieger and P. Vogl,

\$ Phys. Rev. B. Vol. 48, No. 19, 15 Nov. 1993, pp. 14276-287
\$ To convert uniaxial strain to biaxial strain in the orthorgonal
\$ direction, use eu=-D*eo where eo is the strain in a direction
\$ orthorgonal to the uniaxial direction.

\$ (for shear component of hole mobility enhancement)
\$ [3] WANG et al., IEEE TRANSACTIONS ON ELECTRON DEVICES, VOL. 53, NO. 8,
\$ AUGUST 2006, p. 1840

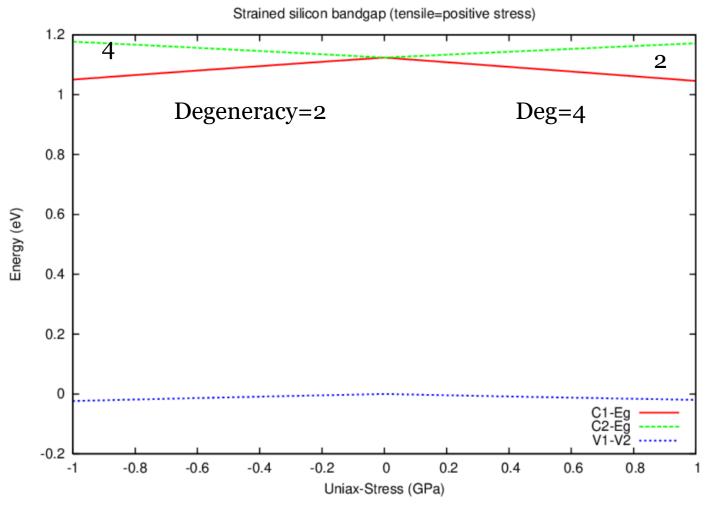
\$ (for shear component of electron mobility enhancement)\$ [4] CHIDAMBARAM et al.,

\$ IEEE TRANSACTIONS ON ELECTRON DEVICES, VOL. 53, NO. 5, MAY 2006, p. 944

\$

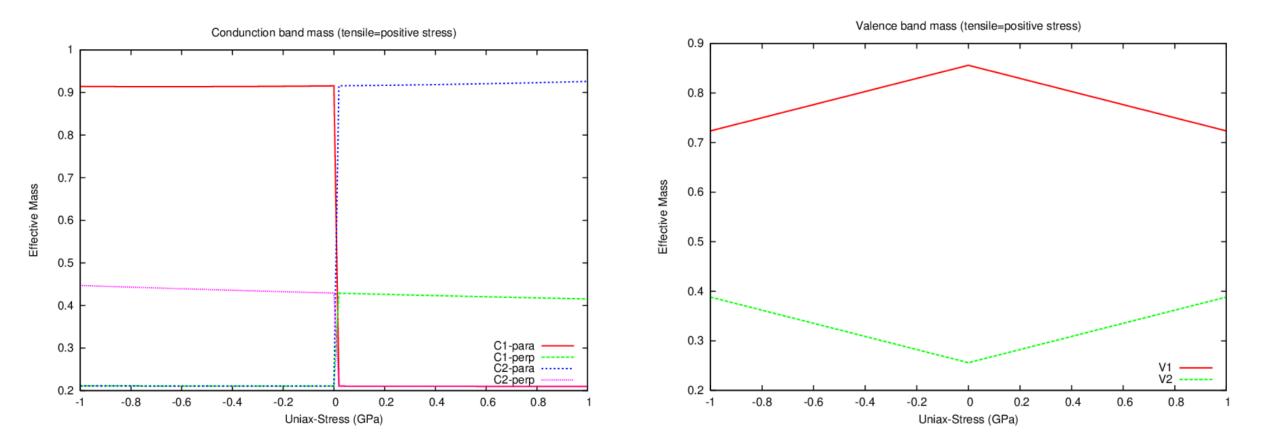
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Stress dependent effective masses





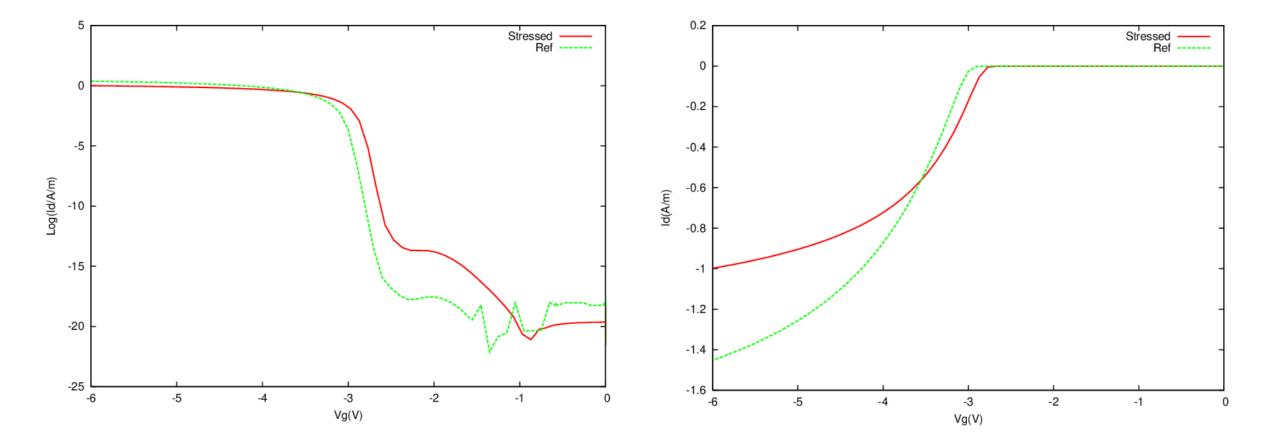
Stress induced mobility change is fitted to theory for single crystal silicon. In material macro presentation, these are stress dependent mobility factors \$ this is electron enhancement factor, stress unit = GPa
\$ tensile stress enhances electron mobility
intern_func1 variation=table
table(stress_xx)
-1. 1.
0. 1.
0.5 1.4
1. 1.7
end_table

\$ this is hole enhancement factor, stress unit = GPa
\$ compressive stress enhances electron mobility
intern_func2 variation=table
table(stress_xx)

-1. 2.
-0.5 1.5
0. 1.
1. 1.
end_table



Vt shift has the same trend as experiment Using strained silicon theory based on material model fitted to kp theory





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Crosslight provides convenient TCAD tools for simulating stress effects in TFT

While stress computation is relatively straightforward, the exact mechanism for Vt shift may be more complicated

Stress induced gate work function shift, strained microcrystal-Si, charged traps generation, doping deactivation can all be factor(s) and Crosslight TCAD can be a useful tool for analysis and design

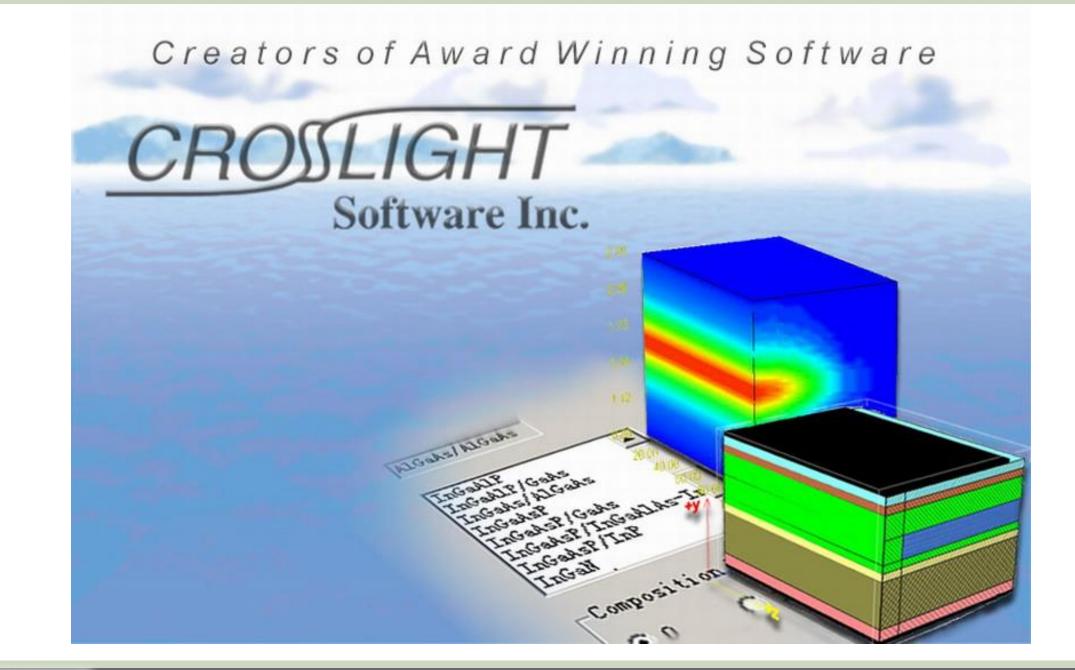


Thanks for your attention!



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