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Physics-Based Compact Model of Independent Dual-Gate BEOL-Transistors for Reliable Capacitorless Memory

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ABSTRACT Capacitorless DRAM architectures based on Back-End-of-Line (BEOL)-transistors are promising for long-retention, high-density and low-power 3D DRAM solutions due to its low leakage, operational flexibility, and monolithic integration capability. Different from classical silicon-based devices, in-depth studies on the performances of nanoscale multi-gate transistors (e.g., a-InGaZnO-FET) are still barely conducted for physical description, due to the complicated multi-gating principle, finite-size effects on transport, increased variation sources and enlarged parasitic effect. Hence, high-performance multinanoscale (down to ~ 50 nm) dual-gate a-IGZO transistors are fabricated, and a physical compact model is developed based on the surface potential for dual-gated coupling and the disordered transport with finite-size-correction. The short channel behaviors on sub-threshold swing, mobility and threshold voltage are investigated, and contact effects are validated by the transfer-line method (TLM). Regarding the specific challenge of dual-gate alignment, possible misalignment and parasitic effects on multi-device fluctuations are important of large-scale circuit design and analyzed by TCAD simulations. Besides, the bias-temperature instability (BTI) has been comprehensively investigated. In awareness of the above effects, this model bridges fabrication-based material properties and structural parameters, assisting in a threshold fluctuation-resistant operation scheme for capacitorless multi-bit memory, showing a great potential in future monolithic integration circuit design using BEOL-transistor.

INDEX TERMS BTI, compact model, contact effects, DRAM, independent dual gate a-IGZO-FET, disorder semiconductor, hopping, transfer-line method.

I. INTRODUCTION

Amorphous In-Ga-Zn-O (a-IGZO) exhibits promising attributes for Monolithic 3D integration (M3D), primarily owing to its compatibility with the Back-End-of-Line (BEOL) process [1]. The a-IGZO field-electric-transistors (FETs) take advantage of moderate mobility and ultra-low leakage current due to the disordered transport mechanism and ultra-wide energy band [2], [3], [4]. This presents a promising solution to circumvent the limitations associated with memory capacitance and cell coupling prevalent

in conventional 1T1C dynamic random access memory (DRAM) technology [5]. By utilizing the gate dielectric layer of the transistor as a storage node capacitor, it offers higher storage density and longer data retention time in capacitor-less memory design [6]. Due to the increasing demand for dense memory design, the scaling down of BEOL-transistors is highly required. However, the sub-threshold property and reliability issues of a-IGZO-FET arise with the scaling-down due to the inferior gate controllability [7], [8], [9]. Variations in material composition and fabrication process



FIGURE 1. (a) Schematic illustration of the IDG a-IGZO FETs with a thickness of ~5 nm. (b) Agreement between analytical and numerical results of back gate surface potentials at different V_{DS} with errors in the inset. $V_{TG} \& V_{BG}$ denotes DG-synchronized-sweep with the same voltage.

contribute to the changes in density of states, and further induce reliability issues in sub-threshold region. It hindered advanced large-scale applications, especially in multi-bit data processing, which is sensitive to device instability [10].

Recently, independent dual-gate (IDG) technology has exhibited the potential to overcome the limitations [11], [12], which are attributable to operational flexibility, adaptive threshold and excellent gate controllability. Nevertheless, when considering complicated effects of dual-gate coupling, disordered transport and sub-threshold degradation at the nanoscale, the compact modeling of BEOL-devices is still incomplete to satisfy future reliable M3D circuit design [13]. To tackle with this issue, the high-performance dualgated a-IGZO-FET is fabricated and modeled to enable cross-layer co-design. On the basis of Newton correction method and dual-gate coupling effects, an accurate surface potential (SP)-based physically compact model is developed to support three operation modes. Particularly, the percolation-theory-based transport mechanisms have been corrected considering the finite-size effects on the amorphous channel [14]. With this model, the critical effects of short channel effects, top gate misalignment, multi-device fabrication variability, bias-temperature instability (BTI), and parasitic effects are investigated. Based on the above effects, this model enables compensation for the attenuation of voltage at the charge storage node (V_{SN}) resulting from fabrication process variations, thereby facilitating accurate and tolerant multi-bit programming. Thus, robust design metrics of monolithic integration circuits can be explored by considering the fabrication-related material and structural parameters.

II. DEVICE FABRICATION AND MODELING A. DEVICE FABRICATION

Fig. 1(a) shows the schematic illustration of the fabricated IDG-IGZO FET with a channel width of 200 nm and a channel length of 50 nm. Firstly, layers of 20 nm W (bottom gate), 4.5 nm HfO₂, and 5 nm a-IGZO were deposited and patterned. The S/D patterns were lithographed by electron beam lithography (EBL), then 10/30 nm Ti/Au was deposited by the E-beam evaporator and lifted off. Next, dual stacks of 2/5 nm Al₂O₃/HfO₂ were deposited as the passivation layer and top-gate dielectric. Finally, the top gate was patterned and lifted off. Fabrication process details can be found in our previous publication [11].

B. NANOSCALE SURFACE POTENTIAL CALCULATION

Due to the conduction band-edge fluctuations in the a-IGZO material [15], disorders (e.g., localized states, etc.) are generally present near or below the conduction bandedge. Therefore, the amorphous channel carrier density is dominated by the density of localized and extended states, closely related to changes in IGZO composition [16]. Based on the Fermi-Dirac distribution, the Poisson equation is given as Eq. (1), shown at the bottom of the page where $\Phi_{TA} = k_0 T/q$, $\Phi_{TA} = k_0 T_A/q$. And φ is the potential in the channel layer, $\varepsilon_{\rm S}$ is the dielectric constant of semiconductor, q is the elementary charge, $N_{\rm T}$ is the total concentration of localized states, τ_0 is the lifetime of carriers, ν_0 is the attempt-to-escape frequency, k_0 is the Boltzmann's constant, and T_A is the characteristic temperature of the exponential DOS. The first and second terms in the right of Eq. (1) represent the localized state and the extended state carrier concentrations, respectively. The defectrelated carrier concentration dominates the subthreshold region [17].

Referring to the mathematic manipulation in [17], based on boundary conditions of bottom/top channel surfaces (Eqs. (2), (3)), the original transcendental equation for IDG-FETs can be obtained in Eq. (4). shown at the bottom of the page

$$Q_{BG} = C_{OXB}(V_{BG} - V_{BF} - \varphi_{BG}) = - \varepsilon_s \frac{\partial \varphi}{\partial y}\Big|_{y=0}$$
(2)

$$Q_{TG} = C_{OXT}(V_{TG} - V_{TF} - \varphi_{TG}) = \varepsilon_s \frac{\partial \varphi}{\partial y}\Big|_{y=t_s}$$
(3)

where $q_1 = Q_{\text{TG}}/(C_{\text{OXT}} \cdot \Phi_{\text{TA}}), q_2 = Q_{\text{BG}}/(C_{\text{OXB}} \cdot \Phi_{\text{TA}}),$ $A_0 = 2 qN_{\text{T}}\varepsilon_{\text{S}} \cdot ((\pi T/T_{\text{A}})/sin(\pi T/T_{\text{A}})) \Phi_{\text{TA}},$

$$\frac{\partial^2 \varphi(x, y)}{\partial x^2} = \frac{q}{\varepsilon_S} \left[\left(\frac{\frac{\pi T}{T_A}}{\sin\left(\frac{\pi T}{T_A}\right)} \right) \cdot N_T e^{\frac{(\varphi - V)}{\Phi_{TA}}} + N_T v_0 \tau_0 e^{\frac{(\varphi - V)}{\Phi_T}} \right]$$
(1)

$$FF(\varphi_{S1}) = \left[k_1q_1 + \alpha \cdot \coth\left(\frac{\alpha}{2} + t_s \cdot \delta\right)\right](k_1q_1 + k_2q_2) - A_0exp\left(\frac{V_{G1} - \varphi_{S1}}{\Phi_T}\right) - B_0exp\left(\frac{V_{G1} - \varphi_{S1}}{\Phi_{TA}}\right)$$
(4)

 $B_0 = 2 q N_{\rm T} \varepsilon_{\rm S} v_0 \tau_0 / \Phi_{\rm TA}$, $t_{\rm s}$ is the active layer thickness, $Q_{\rm BG}$ and $Q_{\rm TG}$ are charge densities, $C_{\rm OXT}$ and $C_{\rm OXB}$ are gate capacitances, $V_{\rm BF}$ and $V_{\rm TF}$ correspond to the flat-band voltages of the top and bottom gates, respectively, and α is closely related to the interfacial coupled electric field.

The segmentation method is employed in the sub-threshold and above-threshold regions to simplify the *Coth/Csch* functions and decouple transcendental equations for the analytical solution of surface potential at the nanoscale. In the subthreshold region, the free charge density is approximately zero, according to this equivalent capacitance model, and φ_{ss} can be obtained by

$$k_{eq} = 1/(1+1/k_1+1/k_2) \tag{5}$$

$$\varphi_{ss} = \left(k_2(k_1V_{G1} + V_{G2}) + k_{eq}(V_{G1} - V_{G2})\right)/k_2(1+k_1)$$
(6)

where $k_1 = C_{\text{OXT}}/C_{\text{S}}$, $k_2 = C_{\text{OXB}}/C_{\text{S}}$, $V_{\text{G1}} = V_{\text{TG}}-V_{\text{TF}}$, and $V_{\text{G2}} = V_{\text{BG}}-V_{\text{BF}}$. In the above-threshold voltage region, *Coth/Csch* functions are approximated by zero, the transcendence equation is transformed into Eq. (7), shown at the bottom of the page and φ_{sn} is easily obtained as Eq. (8), shown at the bottom of the page here, q_2 can be expressed as a function of φ_{S1} , $q_2 = (V_{\text{G2}}-\varphi_{\text{S1}})/\Phi_{\text{TA}} + 2ln(k_1q_1)$. By connecting φ_{sn} and φ_{ss} , an initial analytic solution is given by a smoothing function and its accuracy is further improved by the 1st-order Newton's correction series in

$$\begin{cases} \Delta F = -FF/FF^{(1)}\\ \varphi_{S1} = smothf\{\varphi_{SS}, \varphi_{Sn}\} + \Delta F \end{cases}$$
(9)

where *smothf*(*x*) is the smoothing function that connects piecewise surface potentials. Finally, the analytical solutions of the channel surface potential φ_{S1} at different V_{DS} agree with numerical results, as described in Fig. 1(b), and the error is less than 22.5 mV. And the numerical results were obtained by iterative calculations of Eq. (1) using Wolfram Mathematica numerical computation software.

C. MOBILITY WITH CONTACT EFFECTS AND CURRENT MODEL

The fluctuation of the mobility edge in the amorphous state of a-IGZO leads to localized states in the band gap and potential barriers in the extended state [3]. As for the intrinsic transistor properties, carrier density and T-dependent transport in an amorphous channel have been reported in [18]. As the size of TFTs scales down to the nanoscale, the effects of the extended-state potential barrier, localizedstate characteristic temperature, and gate coupling become increasingly significant and can no longer be ignored. On the classical percolation and hopping mechanisms, the mobility model (Eq. (12), (13), (14)) includes finite-size effects while also considering the coupled carrier density for dual-gate modulation [12]. Due to the obvious contact resistance (R_C) effect in the ultra-scaled device [19], [20], the applied voltage drop in the contact area is much larger than that in the channel, which causes mobility degradation. The mobility considering R_C is incorporated into Eq. (13).

$$\mu_P = \mu_b A_P (V_{G1} - V_{P1} + V_{G2} - V_{P2})^{4(D - W/D)} (10)$$

$$\mu_V = \mu_0 A_V (V_{G1} + V_{G2})^a \tag{11}$$

$$1/\mu_{EF} = 1/\mu_P + 1/\mu_V + 1/\mu_{coup} \tag{12}$$

$$\mu_{EF_{CON}} = \mu_{EF} / (1 + \theta \mu_{EF} R_C (V_{GS} - V_T))$$
(13)

where μ_b and μ_0 are constant mobility terms for percolation and hopping, W/D is the spatial coherence ratio of the potential barriers, $V_{P1,2}$ is the transition voltage, A_P and A_V are related to N_T and T_A , and θ is related to the electrode contact length. Moreover, based on Pierret and Shields' theory [21], combined with μ_{EFCON} , the Pao-Sah current integral equation is solved to obtain

$$I_{D} = \frac{\mu_{EFCON}}{L/W} \begin{cases} C_{OX1} \left[V_{Gf}(\varphi_{D,1} - \varphi_{S,1}) + \frac{1}{2} \left(\varphi_{D,1}^{2} - \varphi_{S,1}^{2} \right) \right] + \\ C_{OX2} \left[V_{Gf}(\varphi_{D,2} - \varphi_{S,2}) + \frac{1}{2} \left(\varphi_{D,2}^{2} - \varphi_{S,2}^{2} \right) \right] + \\ 2\Phi_{TA} \left[C_{OX1}(\varphi_{D,1} - \varphi_{S,1}) + C_{OX2}(\varphi_{D,2} - \varphi_{S,2}) \\ + AF \cdot N_{T} t_{s} e^{\Phi_{TA} V_{Geff}} \right] \end{cases}$$
(14)

III. RESULTS AND DISCUSSION A. MODEL VALIDATION

Transfer curves on a logarithmic and linear scale at different V_{TG} are simulated in Fig. 2(a), which shows the great potential of threshold modulation. For IDG devices, under the influence of the bottom gate, electrons accumulate at the bottom interface of the a-IGZO and form the main channel [22]. Due to the dual-gate coupling effect, the top gate can flexibly control the energy band and current density distribution of the channel, as shown in Fig. 2(b). The output and transfer curves for DG-synchronized-sweep are shown in Figs. 2(c, d). An excellent agreement between the model and experimental data has been demonstrated. Due to the dual-gate coupling effect, the carrier scattering induced by the interface defects can be mitigated, resulting in a weakened interfacial effect and better *SS* performance.

In order to explore the *L*-dependent device characteristics at nanometer size, the model reproduces the IV characteristics of IDG-FETs at different *L* (e.g., Fig. 3(a)). As shown in Fig. 3(c), as *L* decreases (300 nm \sim 50 nm), the *SS* decays to 92.49 mV/Dec, the mobility decreases by 70%,

$$FF(\varphi_{S1}) = k_1 q_1 (k_1 q_1 + k_2 q_2) - A_0 exp\left(\frac{V_{G1} - \varphi_{S1}}{\Phi_T}\right) - B_0 exp\left(\frac{V_{G1} - \varphi_{S1}}{\Phi_{TA}}\right)$$
(7)

$$\varphi_{sn} = \frac{k_1 V_{G1} + k_2 V_{G2}}{k_1 + k_2} - \Phi_T \cdot Lam \left(\frac{B_0 \Phi_{TA} t}{k_1 V_{G1} (k_1 + k_2)} exp \left(\frac{k_1 (V - V_{Gf}) + k_2 (V - V_{Gb})}{\Phi_T (k_1 + k_2)} \right) \right)$$
(8)



FIGURE 2. (a) Comparison of model and experimental transfer curves and trans-conductance at different V_{TG} . (b) Energy band diagrams and carrier concentration distributions under different top-voltage conditions. Agreement between our model and experimental data of the (c) transfer curves, (d) output curves under the DG&BG synchronized sweep.



FIGURE 3. (a)Agreement between model and measured results for transfer curves for *L* from 300 to 50nm. (b) Comparison of model (w/o contact effects) and experimental transfer curves in linear scale at L = 50 nm. (c) *SS* increases and u_{eff} decreases with *L*. (d) Mobility with and w/o contact effects. (e) Total device resistance vs. *L* measured by transfer-line method under different V_{G} conditions.

and the V_{TH} is significantly shifted. The DIBL effect is also not significant for the DG-FET as *L* decreases to 50 nm compared to the single-gate device [23], as shown in Fig. 3(b). An extreme point of the channel potential may shift and indicate short channel effects at small *L* and large V_{DS} . The *SS* performance of the short-channel device is defined as $\Phi_{\text{T}}(\ln 10)(1/\delta + C_{\text{D}}/C_{\text{OX}})$, where δ is related to the minimum value of the channel potential [12]. The *SS* degradation leads to an increase in the leakage current, which in turn affects the DRAM retention. There are non-ideal



FIGURE 4. I V curves of BG single-sweep and TG&BG synchronized-sweep of BG stress measured under (a) PBTI for L_{CH} =100nm IDG-FET. Degradation of ΔV_{TH} at (b) PBTI was investigated using various sweeping methods ΔV_{TH} was subjected to fitting using a semi-classical model. (c) Schematic of top gate misalignment and the potential distribution diagram. (d) Band diagrams and carrier conc. corresponding to the surface un- covered and covered by the top gate. (e) TCAD simulated ΔV_{TH} matrix for different V_{TG} and misalignment distances, with 30 nm-channel length and 5 nm-thickness. (f) Distribution characteristics of 12 groups of devices under different V_{TH} and insert the variations of V_{TH} . (g) Normalized mobility properties extracted from I-V curves and average μ is inserted.

ohmic contacts in short-channel devices, the effect of contact resistance increases as the channel length decreases, and the channel resistance ($R_{\rm CH}$) divider decreases, resulting in lower mobility and lower current. Furthermore, a comparison of different mobility models (classical and contact-limited) and experimental data is depicted in Fig. 3(d), where the contact resistance leads to a significant decrease in mobility and current when L=50 nm. Based on the TLM, the total resistance ($R_{\rm TOT}$) was calculated for different L corresponding to different $V_{\rm GS}$ and the $2R_{\rm C}$ was extracted from the intercept of the linear fit to the $R_{\rm TOT}$ as shown in Fig. 3(e) with $2R_{\rm C} = 860\Omega \cdot \mu$ m.

B. RELIABILITY ASSESSMENT

The continuous scaling improves metrics of efficiency, density, and area in circuits and architectures. However, continued miniaturization leads to increasingly serious BTI effects and non-uniformity of the manufacturing process.

Since a-IGZO FETs are inherently n-channel devices, the reliability of positive bias temperature instability (PBTI)



FIGURE 5. (a) Description of write operation and read mode in a 2TOC cell. (b) V_{SN} compensation values corresponding to different V_{RBG} . (c) Simulated V_{SN} development over time for different storage states and different V_{WBG} , where (1111)₂, (0111)₂ and (0011)₂ are different binary-based input values. (d) 2TOC cell multi-bit operation waveforms, and Input1 to Input7 correspond to I_{WBL} input values of 0.1 μ A to 0.7 μ A in steps of 0.1 μ A, respectively. WWL: write word line, WBL: write bit line, WBG: write back gate.

is very important. Here, Fig. 4(a) shows the PBTI with measured IV curves (V_{DS} =100 mV) for BG-sweep and DGsynchronized-sweep. PBTI leads to a positive V_{TH} shift, and during synchronized-sweep, the BTI effect is mitigated due to the dual-gate coupling effect, which leads to the bulk accumulation of carriers away from the interface [24], thus exhibiting more reliable threshold characteristics. PBTI in IGZO is mainly due to charge trapping or relaxation kinetics, which have different time kinetics, voltage acceleration factors and activation energies. As reported in [7], [25], a physics-based model is employed to reproduce stress and relaxation traces recorded under various test conditions. Here, we simplify the complex physics-based model to a unified empirical BTI degradation model that can be expressed as

$$\Delta V_{TH} = A_E t^\beta V^a_{\rho\nu} e^{-\frac{E_a}{\Phi_T}} \tag{15}$$

where, ΔV_{TH} is the threshold voltage offset due to BTI, *t* is the stress time, *V*ov is the stress voltage, E_a is the activation energy, A_{E} , α and β are degradation fitting parameters. The degradation fitting parameters are extracted by fitting the ΔV_{TH} measured from the experiment, as shown in Fig. 4(b). Then, the above ΔV_{TH} is embedded into the compact model via the effective gate voltage $V_{\text{GEFF}}=V_{\text{GS}}$ - ΔV_{TH} , further capturing DRAM cell retention degradation induced by stressing.

The source/drain asymmetry effect due to top gate misalignment in an advanced process node is investigated using TCAD simulation. For poorly aligned IDG-FET, the offset area induces additional series resistance and weakens controllability for the channel potential, and the increased overlapping area aggravates the effects of overlap capacitance and gate leakage, as shown in Fig. 4(c), which results in degradation of the data integrity of the 2T0C DRAM cell. The threshold adjustment feature based on IDG-FETs compensates the data integrity by adjusting the $V_{\rm TH}$ to alleviate leakage current through the top-gate voltage. It facilitates accurate and tolerant multi-bit programming. When applied a negative top-gate voltage, conduction band decreases at surface, which is not covered by top-gate. Correspondingly, carrier concentration will increase, as shown in Fig. 4(d). Moreover, Fig. 4(e) exhibits a relationship between top gate offset distance and ΔV_{TH} at different V_{TG} .

We further investigated the device-to-device variations of the nanoscale IDG-FETs, Fig. 4(f) shows the variation of measured IV curves for 12 groups of devices using the same process steps. The average V_{TH} decreases as V_{TG} increases, accompanied by an increase in V_{TH} variation. Additionally, the intrinsic parameters of the transistor are extracted from transfer characteristics, demonstrating mobility fluctuations among different devices, as shown in Fig. 4(g). These fluctuations primarily originate from variations in material composition and fabrication process of a-IGZO, which lead to changes in density of states, thereby affecting the extended-state potential barrier and localized-state characteristic temperature. It will result in fluctuations of retention time among cells, challenging multibit computing.

C. NOVEL 2TOC CELL SIMULATION

This model is written in Verilog-A and compiled to commercial circuit design simulator as new active components. For the design of a high-density storage DRAM architecture, a multi-bit storage current read/write operation scheme based on IDG-2T0C was verified [26], and the novel capacitorless circuit unit is shown in Fig. 5(a). This is mainly gate voltage modulation that alters the DG coupling effect to change the channel potential and carrier concentration, which in turn stabilizes the threshold voltage of the read/write transistor in the read, write and hold cases of the 2T0C cell. During write operation, the write bit line (WBL) is shorted to read bit line (RBL), the write transistor (T_W) and the read transistor (T_R) forms a diode connection to charge the storage node (SN) through WBL, the V_{SN} can be adaptively changed to compensate for the $V_{\rm TH}$ variations of $T_{\rm W}$. During read operation, different current values are obtained at RBL depending on the $V_{\rm SN}$ nodes, as shown in Fig. 5(a). Based on the $T_{\rm R}$ threshold regulation capability, the $V_{\rm RBG}$ is adjusted so that the $V_{\rm SN}$ rises to a high level after the write word line (WWL) is turned on and stabilizes at the target value after the WWL is turned off, as shown

in Fig. 5(b). This method effectively mitigates the current degradation as well as compensates for the degradation of the $V_{\rm SN}$ due to the coupling effect of the WWL turned-off device. For the $T_{\rm W}$, adjusting $V_{\rm WBG}$ increases the $V_{\rm TH}$ value, thereby reducing its leakage current and improving data retention, as shown in Fig. 5(c). Flexible operation based on the IDG-2TOC can be used to balance device fluctuations due to short-channel effects. Fig. 5(d) demonstrates the timing diagram for the multi-bit operation of the 2TOC memory cell, where $I_{\rm WBL}$ is the write current, $I_{\rm RBL}$ is the read current.

IV. CONCLUSION

In summary, a surface potential-based compact physical model of ultra-scaled independent dual gate a-IGZO-FET was successfully developed, which can be applied to three modes of operation: single gate sweep, dual gate synchronized sweep, and independent gate modulation. This model considers gate coupling and finite-size disordered transport from device physics, and characterizes the short channel effect and contact resistance. The reliability issues due to variations in material composition and manufacturing processes have been addressed. It serves as a novel methodology for evaluating device-to-circuit reliability, thus accelerating the design-technology co-optimization flow in ultra-large-scale integrated circuits using **BEOL**-transistors.

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