# Ferroelectric Thickness Dependent Domain Interactions in FEFETs for Memory and Logic: A Phase-field Model based Analysis

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**Abstract**— We present a phase-field simulation framework for ferroelectric (FE)-FET which captures multi-domain effects by self-consistently solving 2D time-dependent Ginzburg-Landau (TDGL), Poisson's, and semiconductor charge/transport equations. Using our phase-field model and experiments, we analyze electrostatics-driven multi-domain formation and voltage-induced polarization (P) switching for different FE thickness ( $T_{FE}$ ). We show that for  $T_{FE} = 5 \text{nm} - 10 \text{ nm}$ , FEFETs exhibit multi-level memory functionality; while for  $T_{FE} = 1.5$ nm – 3 nm, FEFETs can serve as non-hysteretic switches with enhanced gate control. Our results signify that as  $T_{FE}$  is reduced from 10nm to 5nm, denser domain patterns emerge in FE, and the dominant P-switching mechanism changes from nucleation to domain-wall motion based leading to a decreased memory window with  $T_{FE}$ scaling. Moreover, as  $T_{FE}$  is scaled further from 3nm to 1.5nm, effective permittivity of the gate stack increases due to multidomain electrostatic interactions.

#### I. Introduction

By virtue of its CMOS process compatibility, Hafnium Zirconium Oxide (HZO) based FEFET is emerging as one of the most promising candidates for future electronics. FEFET have been demonstrated to offer multi-level memory/synaptic functionalities for high FE thickness ( $T_{FE}\sim10$ nm) [1-2] (Fig. 1(b)) and non-hysteretic switch behavior with enhanced gate control for low  $T_{FE}$  (<3nm) (Fig.1(c)) [3-4]. However, the multi-domain effects in FE and its correlation with  $T_{FE}$ , as well as the enhanced permittivity  $(\epsilon_r)$  behavior is yet to be understood. To that end, in this work, we analyze multi-domain polarization (P) switching and the origin of enhanced- $\epsilon_r$  in FEFETs and their dependence on  $T_{FE}$ . Our analysis is based on self-consistent phase-field model of FEFET, validated with our experimental characteristics of metalferroelectric-insulator-metal (MFIM) and metal-ferroelectricinsulator-semiconductor (MFIS) stack for different  $T_{FE}$ . Unlike previous FEFET models [5-8], which assume a certain number of FE domains [5], in our model, the multi-domain formation in the FE is self-consistently determined by electrostatics, which allows us to comprehensively analyze the microscopic domain interactions in FE and its influence on the underlying transistor channel. Using our model, we provide several insights into the  $T_{FE}$ dependent behavior of FEFETs.

#### II. PHASE-FIELD MODELING OF FEFET

In our phase-field simulation, we self-consistently solve the TDGL equation [9-10] for *P*, Poisson's equation for potential and semiconductor charge equations for charge density profile. The considered FEFET structure, simulation flow and parameters are shown in Fig. 3. We consider HZO (Hf<sub>0.5</sub>Zr<sub>0.5</sub>O<sub>2</sub>) as FE, where the *P* direction is assumed to be along the film thickness, which is analogous to the c-axis of its orthorhombic crystal phase. In addition, the Landau coefficients used in simulation are assumed to be strain normalized based on the assumption of a stress free

interface. Note, unlike a previous FEFET model [6] based on 1D Landau-Khalatnikov equation, we use 2D TDGL equation to capture *P* variation along the FE thickness. Other advancements of our model are summarized in Fig. 2.

#### III. VALIDATION OF PHASE-FIELD MODEL

To validate our phase-field model, we fabricate and measure the P-V characteristics of MFIM (HZO-Al<sub>2</sub>O<sub>3</sub>) stack. Our simulation results (Fig. 4(a-c)) signify good agreements with the measured P-V characteristics for different  $T_{FE}$  (5, 7 and 10nm). The P profile of the FE at 0V in Fig. 4(d) suggests the formation of multi-domain (MD) state. Formation of such MD state occurs to suppress the depolarization electric(E)-field in FE by forming inplane E-field (stray field) in the FE-DE interface (Fig. 5). In addition to the electrostatic energy associated with stray field, there is a gradient energy ( $\sim$ (dP/dx)<sup>2</sup>) associated with the P variation near the domain-wall (DW). With the decrease in  $T_{FE}$ , the span of DW (along the z-axis) decreases, which reduces the gradient energy of DW. This leads to the formation of larger number of domains with decreasing  $T_{FE}$  (Fig. 4(d)), which suppresses the depolarization Efield in FE more effectively. As the number of domains in the FE layer depends on electrostatic interaction between the FE and DE layer, the DE layer thickness  $(T_{DE})$  as well as its permittivity impacts the number of domains. The simulated and measured major/minor P-V characteristics of 10nm HZO with T<sub>DE</sub>=3nm and 5nm are shown in Fig. 6(a) and Fig. 7(a), respectively, showing good agreement. For  $T_{DE}$ =3nm, the domain pattern is less dense compared to  $T_{DE}$ =5nm. Therefore, the depolarization-field is more significant in the former case. As a result, the applied voltage driven P-switching in FE for TDE=3nm takes place as a combination of domain nucleation and DW motion (Fig. 6(b)). However, for  $T_{DE}$ =5nm, the domain pattern is relatively denser and hence, the P-switching takes place only through DW motion (Fig. **7(b)**). The number of domains for different  $T_{FE}$  are shown in Fig. 8 suggesting denser domains with decreasing  $T_{FE}$ . Therefore,  $T_{FE}$ scaling should prefer DW motion-based P-switching over domain nucleation. Based on this understanding, let us now discuss the FEFET characteristics and their dependence on  $T_{FE}$  scaling.

#### IV. MULTI-LEVEL MEMORY FUNCTIONALITY

We first analyze the FEFETs for  $T_{FE}$  scaled from 10nm to 5nm. The simulated gate charge (Q) versus gate voltage  $(V_{GS})$  for  $T_{FE}$ =10nm (Fig. 9) shows hysteretic characteristics with a two-step increase (decrease) in Q during the forward (reverse) sweep. The P profile of FE layer (Fig. 10) suggests that, with the increase in  $V_G$ , P-switching first happens through nucleation of new +P domain at the source (S) and drain (D) side of FE giving rise to a sharp increase in Q- $V_G$  characteristics. With the further increase in  $V_G$ , the -P domain between the newly nucleated +P and the previous +P domains completely switches to +P through DW motion. This gives rise to second step in the Q- $V_G$  characteristics. Note that the reason for dominant P-switching at the S/D side is the higher E-

field (originating from the depletion regions in S/D) compared to the center of the FE. Considering the SET and RESET condition (after applying  $V_G$ =+5V and  $V_G$ =-5V, respectively), the potential profile of the FEFET is shown in Fig. 11 for  $V_G$ =0V, illustrating highly non-uniform potential distribution in the channel due to MD state of FE. For RESET state, domains near the S/D side are in -P state which induce a large negative potential across the channel-S/D junction. In contrast, for SET state, due to the dominant +Pdomain near the S/D side, the potential across the channel-S/D junction becomes positive. As a result of P-induced change in channel potential, we observe a large  $V_T$  (= $V_{GS}$  at  $I_D$ =1 $\mu$ A/ $\mu$ m) shift in the I<sub>D</sub>-V<sub>GS</sub> characteristics of FEFET between RESET and SET states (Fig. 12). Moreover, due to the step wise P-switching in the FE layer, different  $V_T$  shift in the  $I_D$ - $V_{GS}$  characteristics of FEFET are obtained for different  $V_{SET}$  value as shown in Fig. 12, leading to multi-level memory behavior.

Similar to  $T_{FE}$ =10nm, Q- $V_G$  characteristics for  $T_{FE}$ =5nm (Fig. 13) signifies hysteretic characteristics with smaller step-wise change in Q. This is because for lower  $T_{FE}$  (5nm), the domain density is more (compared to  $T_{FE}$ =10nm). Hence,  $V_G$ -induced Pswitching takes place via DW motion (Fig. 14) leading to a smaller step-wise change in Q- $V_G$  characteristics. Similar to  $T_{FE}$ =10nm, the *P*-switching for  $T_{FE}$ =5nm is more dominant near S/D (Fig. 14). The FEFET potential profile for RESET and SET states at  $V_G$ =0V (Fig. 15) show a larger channel potential and smaller source barrier for the SET state due to the P-switching in FE. Due to P-switching with smaller steps, a smaller change in  $V_T$  is observed in the  $I_D$ - $V_{GS}$ characteristics (Fig. 16) at lower  $T_{FE}$  (~5nm). Note that the  $V_T$  shift between  $V_{SET}$ =3.4V and 3.6V is insignificant as P-switching takes away from the S/D side and thus has less impact on source barrier. Moreover, we observe an increased memory window (MW= $\Delta V_G$ at  $I_D$ =1 $\mu$ A/ $\mu$ m) (Fig. 17) for higher  $T_{FE}$  (=10nm) due to nucleation dominated P-switching. The MW decreases, and  $V_T$  shift becomes smaller for lower  $T_{FE}$  (=5nm) due to denser domain pattern and DW motion based P-switching, however, leads to multi-level memory operation.

 $I_D$ - $V_{DS}$  characteristics for FEFET ( $T_{EE}$ =5nm) with SET state illustrates hysteretic characteristics (Fig. **18(a)**) similar to experimental results in [11]. In the forward path,  $I_D$  increases with the increase in  $V_{DS}$  and exhibits a sharp decrease at some critical  $V_{DS}$ . This is because, the increase in  $V_{DS}$  leads to P-switching via domain wall motion near the drain side (Fig. **18(b)**) due to negative  $V_{GD}$ . This results in an increase in drain-to-channel barrier which reduces  $I_D$ . In the reverse  $V_{DS}$  sweep, the reduced  $I_D$  state is retained due to P-retention and follows a different path from the forward sweep. The critical  $V_{DS}$  (which triggers P-switching) increases with the increasing  $T_{FE}$  (Fig. **19**) and increasing  $V_{GS}$  (Fig. **19**).

Note that, due to the 2D nature of our simulation, the MD state formation occurs only along the channel length. However, due to the polycrystalline nature of HZO, MD state can form along the width direction also. Consideration of MD state along the width can potentially lead to a gradual  $V_T$  shift, as well as a gradual  $V_{DS}$  induced decrease in  $I_D$  as shown in experiments [2,11].

# V. NON-HYSTERETIC OPERATION WITH ENHANCED EFFECTIVE PERMITTIVITY $(\epsilon_r)$ OF GATE STACK

Let us now turn our attention to  $T_{FE}$  scaling below 5nm. The measured capacitance-voltage (C-V) characteristics of MFIS stack (Fig. **20(a)**) with 2.5nm HZO/0.8nm SiO<sub>2</sub>/p-Si as is shown in Fig. **20(b)** suggesting a higher capacitance compared to a physically equivalent MOS capacitor (2.5nm HfO2/0.8nm SiO<sub>2</sub>/p-Si). This

suggests enhanced- $\epsilon_r$  behavior of HZO. The simulated C-V characteristics show good agreement with the experimental results (Fig. **20(b)**). The *P* and *E*-field profile of FE-DE layer of the MFIS stack (Fig. 21) show that the FE layer is in the MD state with a dense domain pattern and no P-switching occurs within the operational voltage range (0-1V). Instead, the charge response is due to the background permittivity of FE and change in polarization magnitude (|P|) and not due to the change in Pdirection through domain nucleation or DW motion. Enhanced- $\epsilon_r$ stems from the fact that some stray E-field lines between the domains with opposite P at OV (Fig. 21(a)) transforms into out-ofplane component at -0.5V(Fig. 21(b)). To understand this phenomenon, a two dipole (with opposite dipole moment) model has been depicted in Fig. 22, where the dipoles are analogous to the domains with opposite P direction in FE. First, note that at 0V, there are stray E-fields lines between the dipoles (Fig. 22). Now, with the increase in voltage, |P| in +P domain increase and -P domain decreases. That leads to a decrease in the stray E-field between the dipoles. Instead, an out-of-plane E-field by compensating more charges at the FE-DE interface. The additional charges on the FE interface from this transformation of stray Efield to out-of-plane E-field leads to an enhanced effective  $\epsilon_r$ behavior of the FE layer. The simulated FEFET characteristics with 2nm HZO is shown in Fig. 23 showing an improved ID and subthreshold swing (SS) compared to a FET with 2nm HfO2. Note that the stray E-fields exist near the DW. As the domain and DW density increases with the decrease in  $T_{FE}$ , the stray E-field in the FE layer increases. Thus, stray E-field can be transformed into outof-plane E-field to a larger extent leading to an increase in effective- $\epsilon_r$  of FE with the decrease in  $T_{FE}$  (Fig. 24). Therefore, the SS improvement in FEFET is more significant compared to HfO2-FET with the decrease in oxide thickness (Fig. 25).

#### VI. CONCLUSION

In summary, we analyze the FEFET characteristics with  $T_{FE}$ =5nm-10nm for memory operation where the number of memory state can be increased in scaled FEFET by reducing the  $T_{FE}$  so that the DW motion becomes preferable P-switching mechanism over domain nucleation. In addition, we investigate the origin of enhanced- $\epsilon_r$  behavior of thin (1.5nm-3nm) FE layer as an outcome of electrostatic multi-domain interaction in dense MD state and analyze its influence in FEFET for logic operation. The trends obtained from our model for different  $T_{FE}$  in terms memory and improved-SS logic operation show good agreement with existing experimental FEFET results (Fig. 26).

### ACKNOWLEDGMENT

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# **FEFET Applications and Modeling Approaches** (b) High FE thickness (c) Low FE thickness (a) FEFET 1 1 1 1 Channel Ferroelectric: FF: Dielectric: DF Gate: G; Source: S; Drain: D

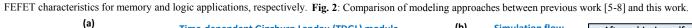
Fig. 1: (a) FeFET structure; (a) Hysteretic and (b) non-hysteretic

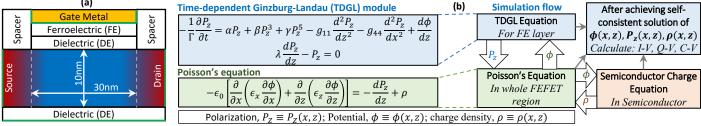
## Previous work -

- [6] Uses 1D Landau-Khalatnikov equation. Assumes Homogenous P along FE thickness.
- [5] Area dependent number of domains. Number of domain does not depend on FE thickness (T<sub>FF</sub>).
- [7,8] Captures multi-domain polarization as a macroscopic/average effect.

#### This work -

- Uses 2D Ginzburg-Landau equation Captures P gradient along FE thickness.
- Electrostatic driven domain formation. Number of domain depends on  $T_{FE}$ .
- Captures Electrostatic driven domain nucleation and domain-wall motion
- Captures microscopic interaction of multi-domain polarization.





(c)  $\alpha$ = -1.63 × 10 $^{9}Vm/C$ ;  $\beta$ =-6.36 × 10 $^{10}Vm^{5}/C^{3}$ ;  $\gamma$ =-7 × 10 $^{11}Vm^{9}/C^{5}$ ;  $g_{11}$ =0.5 × 10 $^{-10}m^{3}V/C$ ;  $g_{44}$ =2 × 10 $^{-10}m^{3}V/C$ ;  $\lambda$ =3nm;  $\epsilon_{x,FE}$ =25;  $\epsilon_{z,FE}$ =18

Fig. 3: (a) Simulated FeFET structure, (b) self-consistent simulation flow among TDGL, Poisson's and semiconductor (silicon) charge equations (c) FE parameters used in FEFET simulation. For the simulation of metal-FE-insulator-metal (MFIM) stack only Poisson's and TDGL equations are used

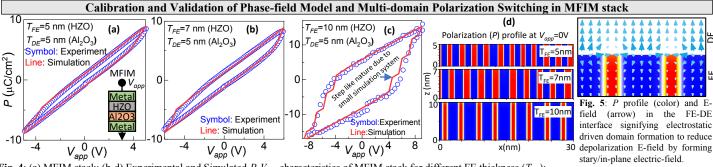


Fig. 4: (a) MFIM stack; (b-d) Experimental and Simulated P- $V_{app}$  characteristics of MFIM stack for different FE thickness ( $T_{FE}$ ); (e) P profile of FE in MFIM stack for  $T_{FE}$  at negative remanent state (- $P_R$ ). Red (blue) color corresponds to +(-)P domains

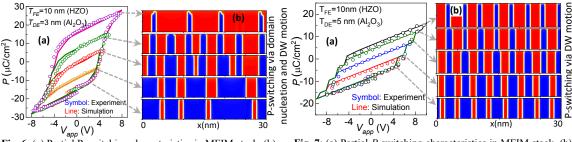


Fig. 6: (a) Partial P-switching characteristics in MFIM stack. (b) Corresponding P profile signifying domain nucleation and domainwall (DW) motion based P-switching.  $T_{FF}$ =10nm,  $T_{DF}$ =3nm.

Fig. 7: (a) Partial P-switching characteristics in MFIM stack. (b) Corresponding P profile at different  $V_{app}$  signifying DW motionbased P-switching.  $T_{FE}$ =10nm,  $T_{DE}$ =5nm.

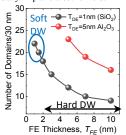
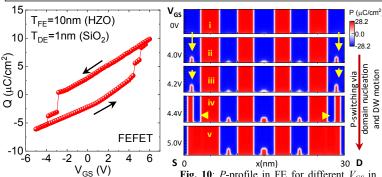
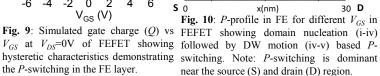
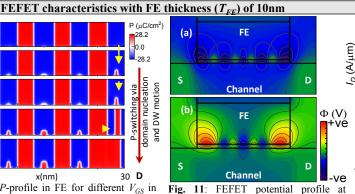


Fig. 8: Number of domains as a function of  $T_{FE}$  for different  $T_{DE}$  showing that the domain pattern becomes denser with decreasing  $T_{FE}$ 



the P-switching in the FE layer.





 $V_{\textit{GS}}$ =0V after (a)  $V_{\textit{RESET}}$ = $V_{\textit{GS}}$ = -5V and (b)  $V_{SET} = V_{GS} = 6V$  showing P-switching induced change in channel potential as well as non-uniform potential in channel.

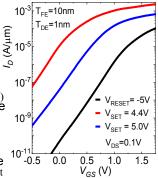
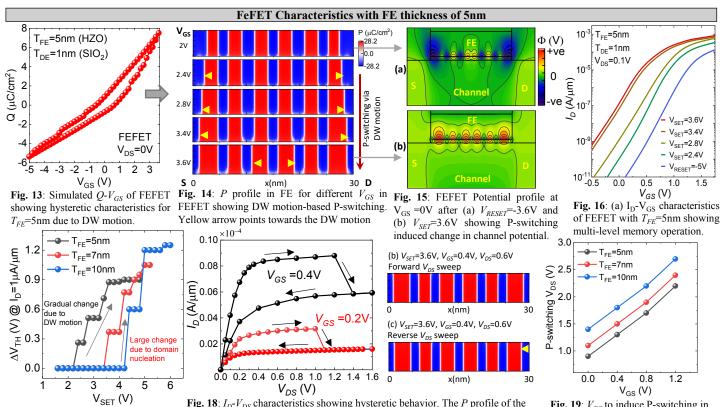


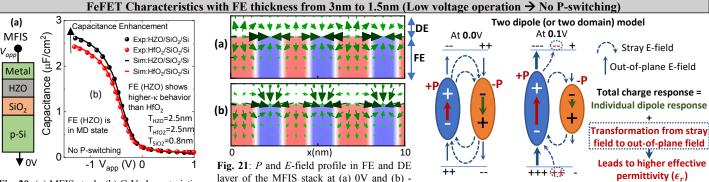
Fig. 12: (a)  $I_D$ - $V_{GS}$  characteristics of FEFET for different  $V_{SET}$ showing the P-switching induced  $V_T$  shift and memory behavior.



**Fig. 17**: (a) Memory window ( $\Delta V_{TH}$ ) for different  $T_{FE}$  and  $V_{SET}$  showing that MW increases with the increase in  $T_{FE}$ .

Fig. 18:  $I_D$ - $V_{DS}$  characteristics showing hysteretic behavior. The P profile of the FE layer is shown for  $V_{DS}$ =0.6V considering (b) forward sweep and (c) reverse  $V_{DS}$  sweep. Due to the P-switching near drain side the drain barrier height increases leading to a lower current in reverse sweep compared to forward sweep.

Fig. 19:  $V_{DS}$  to induce P-switching in drain-side for different  $V_{GS}$  showing P-switching occur at higher  $V_{DS}$  for increasing  $T_{FE}$ ..



**Fig. 20**: (a) MFIS stack; (b) C-V characteristics of MFIS stack compared to MOS capacitor showing high-k behavior of the FE layer.

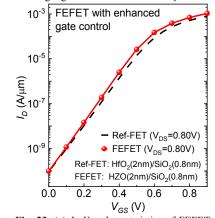
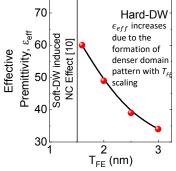
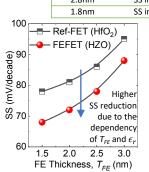


Fig. 23: (a)  $I_D$ - $V_{GS}$  characteristics of FEFET showing higher-k behavior with enhanced gate control compared to Ref-FET (with  $HfO_2/SiO_2$  gate stack).

Fig. 21: P and E-field profile in FE and DE layer of the MFIS stack at (a) 0V and (b) - 0.5V showing the transformation of stary/in-plane E-field to out-of-plane E-field due to the magnitude change of P. This phenomena leads to extra charge in FE-DE interface leading to enhanced effective  $\epsilon_r$  in FE layer. (a)



**Fig. 24**: Effective permittivity of the FE layer for different  $T_{FE}$  showing that permittivity increase with the decrease in FE thickness.



**Fig. 25**: Subthreshold swing (SS) for different  $T_{FE}=T_{HfO2}$  showing more SS reduction in FEFET.

induced change in dipole interaction associated electric field.

T<sub>FE</sub> Experimental Characteristics Model Prediction

10nm Hysteretic [Ref-2] Hysteretic

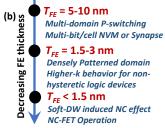
5.5nm Hysteretic [Ref-1] Hysteretic

2.8nm SS improvement [Ref-4] SS improvement

1.8nm SS improvement [Ref-3] SS improvement

Fig. 22: Two dipole (or domain) scenario describing the origin of

enhanced- $\epsilon_r$  of hard multi-domain thin FE film as a voltage



**Fig. 26**: (a)  $T_{FE}$  dependent experimental characteristics of FEFET in different earlier works. (b) Co-relation between  $T_{FE}$  and FEFET characteristics and its application.