

# MODELING A THIN-FILM TRANSISTOR (TFT)

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## MODELING A THIN FILM TRANSISTOR (TFT)

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## ABSTRACT

#### MODELING A THIN FILM TRANSISTOR (TFT)

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In this thesis, we studied the working principles and modeling the static current – voltage characteristics of a Thin Film Transistor (TFT). TFTs are extensively used in flat panel displays for switching the pixels. We first analyzed the characteristics of Amorphous Indium Gallium Zinc Oxide (a-IGZO) TFTs, and reviewed the analytic models to estimate the static electrical behavior of the TFTs in general and the a-IGZO TFT, in particular. We developed a new analytical model for predicting the static electrical behavior of the a-IGZO TFTs and for analyzing and simulating the new electrical circuits when these transistors are used to form new circuits. Finally, we compared the predictions of this model with the measured data and other models.

Keywords: Semiconductor, TFT Models, IGZO, TFT Output Characteristics

ÖZ

## **İNCE FİLM TRANSİSTOR MODELLEME**

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Bu tez çalışmasında, ince film transistorların (TFT) çalışma prensipleri ve statik akım – gerilim karakteristiklerini çalıştık. TFT'lar yoğunlukla düz panel göstergelerinde görüntü gözelerini (piksel) sürmekte kullanılmaktadır. Öncelikle Amorf İndiyum Galyum Çinko Oksit (a-IGZO) TFT'lerin özelliklerini analiz ettik. Daha sonra genel olarak TFT'lerin özel olarak da a-IGZO TFT'lerin statik akım – gerilim davranışlarını kestirebilmek için geliştirilen analitik modelleri inceledik. a-IGZO TFT'lerin statik elektriksel davranışlarını kestirmek ve bunlar yeni devrelerde kullanıldığında, bu devrenin davranışını analiz ve simüle edebilmek amacıyla yeni bir analitik model geliştirdik. Bu modelin tahminlerini ölçülen değerler ve diğer modeller ile kıyasladık.

Anahtar Kelimeler: Yarıiletkenler, İnce Film Transistor Modelleri, IGZO, TFT Çıktı Özellikleri

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# LIST OF ABBREVIATIONS

AMLCD	Active Matrix Liquid Crystal Display
AMOLED	Active Matrix Organic Light Emitting Diode
a-Si:H	Hydrogenated Amorphous Silicon
a-IGZO	Amorphous Indium Gallium Zinc Oxide
CdSe	Cadmium Selenide
DOS	Density of States
FET	Field Effect Transistor
E <sub>IGZO</sub>	Transverse Electric Field of IGZO TFT
$kT_{EFF}$	Characteristic Energy of Activation
LCD	Liquid Crystal Display
LTPS	Low Temperature Polycrystalline Silicon
MOSFET	Metal Oxide Semiconductor Field Effect Transistor
$N_{\scriptscriptstyle EFF}$	Effective Density of States (DOS)
n <sub>EFF</sub>	Effective Carrier Density
n <sub>FREE</sub>	Free Carrier Density
n <sub>LOC</sub>	Localized Charge Density
TFT	Thin Film Transistor
ZnO	Zinc Oxide
$\Phi_s$	Surface Potential
$\Phi(x)$	Transverse Electrical Potential
$\Phi_{_{C\!H}}$	Fermi Potential Lowering by the Drain Bias

$\Phi_{_{FO}}$	Fermi Potential Under Thermal Equilibrium
$\mathcal{E}_{IGZO}$	Permittivity of $\alpha$ -IGZO Active Layer
$\mu_{\scriptscriptstyle BAND}$	Conduction Band Mobility
$\mu_{_{CH}}$	Channel (or Effective) Mobility

### **CHAPTER 1**

#### 1. INTRODUCTION

The image on a large LCD TV screen, or a small LCD mobile phone screen is produced by turning on some pixels and by turning off the others in accordance with the pattern determined by the image. The pixels are turned on by applying electrical current to the liquid crystal properly. The selected pixel generates the shade and not selected pixel passes the light with proper intensity and color. There are thousands of pixels on a display and all of them should be turned on, if needed, by applying electric currents to the selected pixels. This means every pixel on the display could be turned on or off by using a dedicated switch to that pixel. This switch is mostly a transistor operated non-linearly by driven either to ON state (driven into saturation), or to the OFF state (driven into cut-off) from a signal generated in accordance with the image to be displayed on the screen.

On a flat panel the transistors should be fabricated in a matrix array to switch the pixels. The most suitable type of transistor that would be for such an application is a metal-oxide semiconductor field-effect transistor (MOSFET), as it can be easily manufactured cost-effectively on a substrate planarly. Normally MOSFETs are produced on crystalline semiconductor substrate by either a p or n-type epitaxial growth of a crystalline semiconductor layer, and then two terminals at the two ends of a channel are produced by first doping these regions with proper atoms to make

them either p or n-type (opposite doping of the epitaxial region), and subsequently depositing metal regions on top of these doped regions. These metal terminals are called the drain (D) and the source (S). In between the drain and the source an oxide layer is formed and on top of this oxide layer a gate terminal is formed by using a proper conducting material. The channel formed between the source and the drain is turned on by applying a proper voltage to the gate and turned off by applying a voltage of the opposite polarity. This transistor, MOSFET, is the mostly used transistor type today in electronic devices, and the crystalline silicon (Si) is the most mature material to fabricate the very small size MOSFETs, making today the very large scale integrated circuits (VLSI) possible, where nearly 2 billion transistors can be manufactured on a semiconductor die of area of for example of 1mm<sup>2</sup> [1].

However, this type of MOSFETs are not proper to be used as a switch to turn on and off the pixels on a relatively large flat panel display, since firstly on flat displays we do not need the high density of transistors like we do in VLSI circuits, secondly these transistors are not to be produced on a semiconductor die of a very small size (such as 1mmx1mm) but on a much larger flat glass displays (such as from 2cm x 2cm to 2m x 2m). In a display we have generally a large glass panel and whatever to be deposited should be deposited on this panel of glass.

Therefore, on one hand to switch the pixels on the large displays it is sufficient to use relatively larger MOSFETS than those of on a crystalline silicon substrate, where one has to squeeze all the transistors in a very small area to make a very large scale integrated circuit chip (VLSIC). On the other hand, these bigger MOSFETs used to drive the pixels on the displays should also be manufactured with a cost effective way, on relatively large glass type materials. The most suitable technology to manufacture such planar devices on large screens is the thin film deposition technology. A large glass screen can be put into a vacuum system, and a semiconductor, or metal or an organic material can be coated on it by using the thin film technology. The thin film is a film of thickness up to 1-2 micrometer. The vacuum system is made such that the thickness and the purity of the thin film can be controlled to a certain level of precision. When a semiconductor sheet is deposited on an amorphous material such as glass at relatively low temperatures ( $\leq 400^{\circ}C$ ) the semiconductor thin film also becomes amorphous. However, after annealing the amorphous thin silicon films above  $400^{\circ}C$ , the material becomes polycrystalline. The annealing temperature should not exceed  $500^{\circ}C$  when a glass substrate is used because glass cannot withstand these temperatures. Quartz can be used as a substrate to anneal the amorphous thin film semiconductor at higher temperatures to convert to crystalline semiconductor but quartz is not economical to be used on large displays.

As seen in Fig. 1, a thin film transistor (TFT) can be manufactured such that, an active layer that forms the channel is deposited first on a glass substrate. Two metal terminals for the source and the drain can be coated on the two ends of the channel by again metal deposition technique. Then an oxide layer is coated on the channel by chemical techniques. On top of the oxide between the two terminals another metal region can be formed as the gate [2].

In the TFT, as stated above, due to the glass substrates the channel layer formed cannot be made crystalline but it can be made mostly in amorphous or sometimes polycrystalline forms. To make p or n-type material from an amorphous or a polycrystalline material is very difficult but using some techniques such as sputtering some amorphous semiconductors can be doped. Hydrogenated amorphous silicon (a-Si:H) is such a material that can be made slightly p or n-type. This is due to the passivation of the dangling bonds in the amorphous structure by the hydrogen atoms [3].

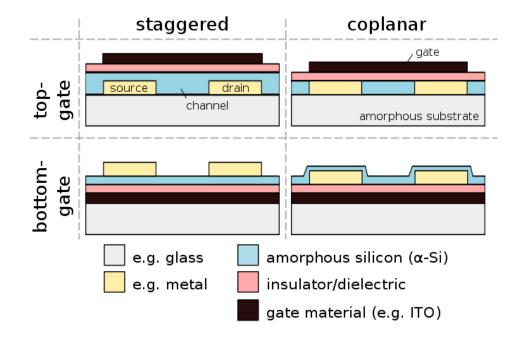


Figure 1 Some different types of basic TFT constructions [4]

In Fig. 1 we see some types of TFT configurations [4]. These are differing from each other in manufacturing order or physical properties. Thin film depositing technique is used to manufacture the devices in all these configurations.

a-Si:H TFTs are the first transistors that were used to switch the pixels on the flat displays. This type of TFT has hydrogenated amorphous silicon as channel material. Since a-Si:H is easily deposited on a substrate material such as glass by using mostly the sputtering technique under vacuum, it has been the first and mostly used material to form a TFT [3]. Although they were widely used, a-Si:H TFTs have some disadvantages. Some of the disadvantages are low field - effect mobility, sensitivity to visible light, threshold voltage shift in off state, etc. These drawbacks are increasing the length between pixels and reducing some driving abilities of TFTs. a-Si:H TFTs were frequently used for AMLCDs backplanes, because high resolution is the most important parameter for AMLCDs, and an a-Si:H TFT can provide this. Drawbacks of a-Si:H TFTs show themselves in active-matrix organic light emitting diode (AMOLED) displays, hence they are not a good choice to drive AMOLEDs. AMOLED displays can be used not only for large flat panel application but also for flexible panel applications. This makes producing flexible electronic devices such as electronic papers, wearable technologies, etc. possible.

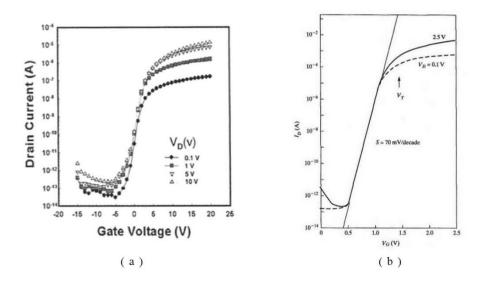
It is not easy and cost effective to run AMOLEDs with a-Si:H TFT driver circuit that can be used in AMLCDs [6-8]. So, researches have been investigating to find different materials to be used in the TFTs. Some examples for the alternative active channel material for the TFTs are low-temperature polycrystalline Si (LTPS) [9], pentacene-based organic materials [10], and metal-oxide semiconductors, which are mostly based on polycrystalline zinc oxide (*ZnO*) [11, 12] and amorphous metal oxide semiconductors [13].

The TFTs using pentacene-based organic materials in active layers were proven to be not reliable and stable. LTPS TFTs are used in some devices and they have relatively good performances. Especially, they provide good device performances with high field effect mobility values. However, they suffer from low level of uniformity when they were used on backplane of large panel areas [9,10]. Metal-oxide semiconductors show better performances than the above cited materials in these applications. They are able to be uniformly deposited on large backplane areas at low temperatures. They have higher mobility values, and are also not affected from visible light. These materials, when being mostly polycrystalline, have large grain boundaries at room temperature. This causes low performance at this temperature and decreases their stability on large backplane areas.

The recently introduced amorphous Indium Gallium Zinc Oxide Thin Film Transistors (a-IGZO TFTs) show much higher field effect mobilities than a-Si:H TFTs and do not suffer from grain boundaries as the polycrystalline materials do[14]. Backplanes that are produced by using that type of TFTs are used in both AMLCDs and AMOLEDs [15, 16].

#### 1.1. Why a-IGZO ?

Although the TFTs are very suitable to be manufactured on large flat displays, regarding their switching performance they are rather much inferior to the MOSFETs made on crystalline semiconductors. The TFTs form the active switches in an active-matrix liquid crystal display (AMLCD), where the ON current determines the rate of pixel charging and the OFF current is associated with the leakage of the pixel voltage [3]. The switching characteristics of a MOSFET of a crystalline semiconductor and a TFT are shown in Fig. 2-a and -b, where the drain current is the current that flows between the drain and the source. As seen a MOSFET has much higher ON currents than the TFT, since in a MOSFET the mobility of carriers is much higher than a TFT.



**Figure 2** (a) Switching characteristics of an a-Si:H TFT (W=39µm, L=5µm)[3] (b) of a MOSFET[5]

In improving TFTs, efforts were increasingly focused on the issues of stability, ON/OFF Ratio, OFF current and leakage. The TFT, when switched, the ON/OFF currents ratio for proper LCD addressing applications should not be less than  $10^6$ . The OFF current should be less than 1nA. As described by Lee et al. [20] stability was always an important issue, the threshold voltage could change in years for example in the TFTs of *CdSe* active layers. TFTs have been notorious for their poor DC stability, which manifest itself as a slow decay of drain current when the device is operated with a steady gate voltage [3].

In 2010 the state-of-art Gen-8 facility is able to process glass substrate approximately 2160x246 mm in size, which produced up to six 52-inch flat panel TV screens from one single substrate [21]. The typical plasma enhanced chemical vapor deposition (PECVD) a-Si:H TFT has field effect mobility ( $\mu_{EFF}$ ) of 0.6–0.8 cm<sup>2</sup>/Vs, OFF-state drain current ( $I_{D_off}$ ) below 10<sup>-13</sup>A, and an ON-OFF ratio about 10<sup>7</sup> [22].

These properties are suitable for LCD switching but the  $\mu_{EFF}$  would be insufficient for the new requirements of next generation displays, especially when the number of pixels are increasing (higher resolution) and more than 100 Hz of frame rate is needed on the TV screens to prevent the blurring effect [21].

The other issue in the active-matrix flat panel display (AM-FPD) is the emissive display such as active-matrix organic light emitting diode (OLED), where AMOLED is directly integrated with the TFT pixel electrode circuit. AMOLED does not need backlight as the AMLCD does, can have extremely high contrast ratio and delivers much better picture quality than AMLCD. However, AMOLED has more stringent requirements on the TFT backplane. Unlike LCD which only requires external electric field to change its phase, OLED takes significant amount of current to produce light by itself. The  $\mu_{EFF}$  of a-Si:H TFT is not high enough to drive a large-area AMOLED. The TFT may constantly operate under high bias; a long term electrical instability is another concern and can also make the pixel electrode circuit design more complicated [21].

The application of a-Si:H TFT backplane is not limited to AMFPD. By integrating the photodiode or radiation sensitive layer with the backplane, a-Si:H TFT can also be used for sensor readout [21, 23]. One important example is the flat panel X-ray imager used for medical imaging [21, 24, 25, 26]. In order to achieve desirable signal gain on the active pixel sensor for signal readout TFTs used for both amplifying and readout TFTs should have a large width-over-length ratio (W/L) of about  $150\mu m/25\mu m$  when using a-Si:H as the channel semiconductor material for the TFT. These transistors take the significant amount of the pixel area (about 50% of

250  $\mu m^2$  pixel size) resulting in a smaller resolution ( $\leq 200 dpi$ ) [21, 27, 28]. In a TFT if the mobility is lower than required, like a-Si:H TFT, one has to increase the W/L to increase the channel current, which causes a larger TFT area and reduces the resolution. Therefore, for a definite gate length (L) using a TFT having a larger mobility is one possible solution to achieve a smaller pixel area without sacrificing the gain performance. It is clear that a new high mobility semiconductor material yet with a uniform amorphous phase over a large area is highly desired [21].

Since 2004, there has been great interest in adapting TFT made of ionic amorphous oxide semiconductors [29]. Specifically, the ternary oxide system that consists of  $In_2O_3$ ,  $Ga_2O_3$  and ZnO have shown promising electrical performance for TFT active layer with a high  $\mu_{EFF}$  (3–12  $cm^2/Vs$ ), low  $I_{D_off}$  ( $\leq 10^{-12}A$ ) and good uniformity compatible with the state-of-the-art Gen-8 substrate size [21, 30, 31].

Table 1 compares amorphous In-Ga-Zn-O (a-IGZO) with other TFT Technologies [21]. It is seen that a-IGZO is the only technology available today that can achieve a desired balance between high mobility and large area uniformity. Although polycrystalline (poly-Si) TFT is currently has the largest mobility it required additional re-crystallization steps such as excimer laser annealing [21, 32], metal seeding [33], or solid phase crystallization [34]. These add more complexity and cost to the process. Also the substrate area used by poly-Si TFT technology (Gen-4, 730*x*920 *mm*) is about 4 generations behind what a-Si:H TFT can achieve today (Gen-8) [21,35].

Technology	Mobility (cm²/Vs)	Visible Light Transparency	Large Area Uniformity	Comments
a-Si:H	<1	Poor	Good (Gen-8)	Low mobility, limited current driving capability
Poly-Si	~100	Poor	Poor (Gen-4)	Additional crystallization process required
ZnO	20~50	Good	Poor	Strong tendency to form poly phase
a-IGZO	3~12	Good	Good (Gen-8)	Balance between mobility and uniformity

 Table 1 Comparison of different TFT technologies [21]

Mono-oxide semiconductor, ZnO, has been used as the active layer in TFT channel. ZnO layer can be deposited by pulse laser deposition (PLD), RF magnetron sputtering or atomic layer deposition (ALD) and the ZnO TFT's  $\mu_{EFF}$  is around  $20-50 \ cm^2/Vs$ . Despite this high mobility, due to its strong poly-crystalline nature even when deposited at room temperature the grain boundary of such oxide semiconductor affects device electrical properties, uniformity and stability over a large area [21,36,37,38]. For a TFT to have a low  $I_{D_off}$  and a high ON-OFF ratio, it is important to control the semiconductor carrier concentration to a very low level. In this sense a-IGZO is much better than ZnO or CdSe [21, 39].

Even though a-Si:H is widely used in TFT backplane, it absorbs visible photons and has a low visible light transmittance ( $\leq 30\%$ ). This has been a major drawback for utilizing such material in optoelectronics and can even raise concern of light-induced instability. On the other hand, similar to *ZnO*, a-IGZO has a wide bandgap ( $\approx 3eV$ ) and is highly transparent in visible light. This property opens up to new path of application such as transparent electronics or see-through display. In addition, light may directly transmit through a-IGZO TFT [21]. In conclusion, a-IGZO TFTs shows unique physical properties and better electrical performance over traditional a-Si:H TFTs, poly-Si TFTs and *ZnO* or *CdSe* TFTs.

A very thorough analysis of the amorphous oxide semiconductors (AZO) in general, IGZO TFTs in particular, is given by Kamiya et.al.[40]. They asserted that the a-IGZO TFTs would meet all the requirements for organic light emitting diode (OLED) displays, large and fast LCDs, 3-D displays, which cannot be satisfied by conventional silicon and organic TFTs. The reasons why the a-IGZO TFT will dominate this area in the future were listed by them [40], some of which are that most device issues, such as uniformity, long-term stability against bias stress and TFT performance are solved for a-IGZO TFTs, and 8G sputtering apparatus was developed and mature enough to manufacture these TFTs on large substrates.

We believed that the future switching device to drive the pixels on large displays would be a-IGZO TFT, and therefore, we studied modeling this transistor in this thesis.

In this thesis, we analyzed the IGZO TFT characteristics and analytic models developed for this type of TFTs. Then, we proposed and developed a new analytic model for an a-IGZO TFT.

## **CHAPTER 2**

## 2. ANALYTICAL TFT MODELS

A transistor is a three-terminal semiconductor device. There are generally two classes of transistors: a bipolar junction transistor (BJT), and a field effect transistor (FET). A TFT is an FET type transistor. In an FET transistor a channel that carries the current is formed between the source and drain terminals. The channel's conductivity, and therefore the drain current  $(I_D)$  that flows through the channel, is controlled by the voltage applied at the gate terminal  $(V_{GS})$ . The static (direct current: DC) behavior of a transistor is sufficiently described by the input characteristics and the output characteristics. In the input characteristics of an FET,  $I_D$  versus  $V_{GS}$  is given for a fixed  $V_{DS}$ , where  $V_{DS}$  is the voltage applied between drain and the source. The output characteristics are between  $I_D$  and  $V_{DS}$  for constant values of  $V_{GS}$  as parameter. Fig. 3-a and -b are the examples for the input and the output characteristics, of an a-Si:H TFT, respectively.

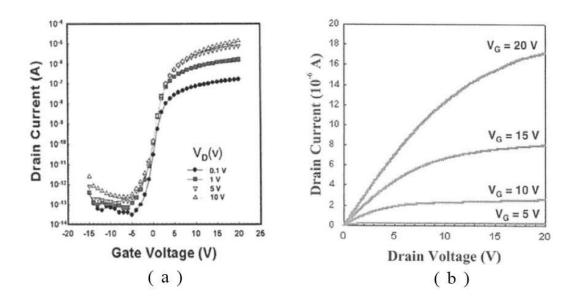


Figure 3 (a) input and (b) output characteristics of a-Si:H TFT [3]

Modeling a transistor is to obtain mathematical relations between  $I_D$ ,  $V_{GS}$  and  $V_{DS}$ , by using the transistor's geometric properties and material characteristics to understand and improve the device electrical performance. The higher the conformance of the model predictions with the measured characteristics, the better the model. Models can be analytical and numerical. Analytical models are given by closed-form mathematical equations that can be solved analytically and easier to use generally. Analytical models can be easily used in computer-aided design (CAD) tools to simulate the circuit behavior as well. However, they are limited in deriving 2- and 3-dimensional device behavior and sometimes are not accurate enough. Numerical models can better represent the device physics [3] especially in 2-and 3dimensions. These are mostly given by some partial differential equations, or integrals that can only be solved by numerical techniques for certain boundary and initial conditions. Numerical models could be more accurate than analytical models but generally they are more difficult and impractical to deal with. They mostly require relatively bigger computer resources and they take much longer times to run. The model is used to predict the behavior of the device before manufacturing it, and at the same time they could be used to derive certain parameters using measured values subsequent to production of the device. The model is mostly used for designing the device to the desired performance. It serves also as a tool for analyzing the behavior of the device produced using the measured data.

In this thesis, we developed an analytical model for predicting and representing the DC behavior of an a-IGZO TFT. Before presenting our model, in this chapter we provide the simplest model first. This is the constant mobility model where a gradual channel formation between the drain and the source is assumed. Then, we analyzed in detail a variable channel mobility model, where the mobility is given by the carrier concentrations for an a-IGZO TFT [17].

#### **2.1. The Analytical FET Models**

#### 2.1.1. The Gradual Channel Approximation FET Model

Fig. 4 shows an FET, where x denotes the direction from the gate terminal to the substrate, that is, the transverse direction, which is perpendicular to the channel; and y denotes the direction from the source terminal to the drain terminal, which is parallel to the channel.

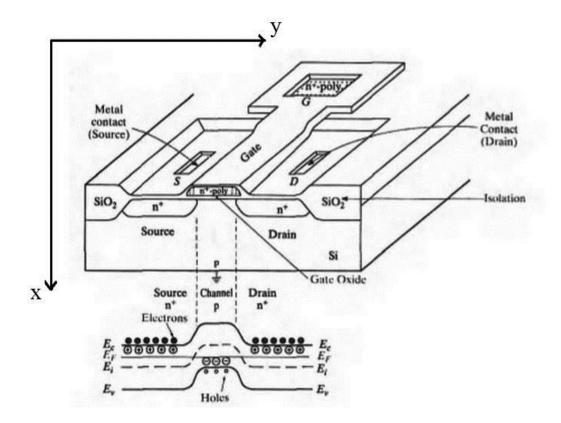


Figure 4 Cross sectional view of an n-channel FET [5]

A voltage applied to the gate changes the free charge concentration in the channel. Changing the number of free carriers per cross sectional area in the channel modulates the conductivity and hence the drain current.

If we assume a gradual channel is formed between drain and source under the influence of both  $V_{GS}$  and  $V_{DS}$  voltages, and a constant carrier mobility in the channel, when the gate potential is higher than the threshold voltage,  $V_{TH}$ , the mobile charge  $Q_I$  in the channel is related to the gate potential  $V_{GS}$  via [3]

$$Q_{I} = -C_{ox} \left[ V_{GS} - V_{TH} - V(y) \right]$$

$$(2.1)$$

where *Cox* is the capacitance per unit area of the  $SiN_x$  gate oxide, and V(y) is the channel potential at a certain y value. The current induced by the majority carriers, in this case electrons for an n-channel FET, can be given as

$$I_D = W \mu_n Q_I E_v \quad , \tag{2.2}$$

where W is the channel width,  $\mu_n$  is the electron mobility, and  $E_y$  is the electric field at y. By substituting  $E_y = \frac{-dV(y)}{dy}$  into Eq. 2.2, we get;

$$I_D dy = W \mu_n C_{ox} \left[ V_{GS} - V_{TH} - V(y) \right] dV(y)$$
(2.3)

By integrating the current along the channel (from y = 0 to y = L, the channel length), using V(0) = 0 at the source and  $V(L) = V_{DS}$  at the drain, we obtain the drain current for the linear region;

$$I_{D} = \frac{C_{ox}\mu_{n}W}{L} \left[ \left( V_{GS} - V_{TH} \right) V_{DS} - 1/2 V_{DS}^{2} \right]$$
(2.4)

In this region where  $V_{DS} \ll V_{GS}$ ;

$$I_D \cong \frac{C_{ox} \mu_n W}{L} \left( V_{GS} - V_{TH} \right) V_{DS}$$
(2.5)

As seen in Fig. 3-b, in the linear region where  $(V_{DS} \ll V_{GS})$  holds, for a constant  $V_{GS}$ , the current changes linearly with  $V_{DS}$ .

The saturation region, where the current saturates and becomes nearly constant even the  $V_{DS}$  increases, is the point where the field-induced carrier density at the drain side disappears as the drain potential increases. Eventually, when

$$V_{DS} = V_{GS} - V_{TH} \tag{2.6}$$

the channel with induced electrons becomes completely depleted of the electrons where this condition is realized. This is called the pinch-off of the channel, where the drain current saturates. For  $V_{DG} \ge V_{GS} - V_{TH}$  Eq. 2.4 and 2.5 are no longer valid. The saturation drain current can be obtained by substituting Eq. 2.6 into 2.5, giving [3]

$$I_{D} = \frac{C_{ox} \mu_{n} W}{2L} \left( V_{GS} - V_{TH} \right)^{2}.$$
 (2.7)

The model gives a fairly good fit to the curves in the linear region and also predicts the saturation point (pinch-off); however, it predicts that after this point the drain current is constant even  $V_{DS}$  continues to increase.

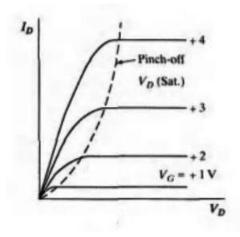


Figure 5 Gradual Channel Approximation of FET [5]

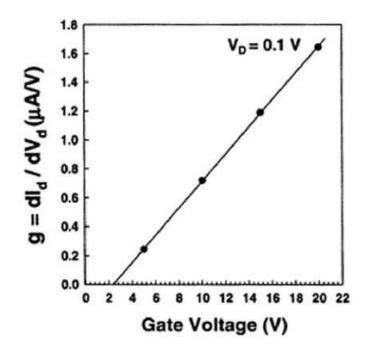


Figure 6 Field effect mobility extracted from channel conductance in the linear

#### region [3]

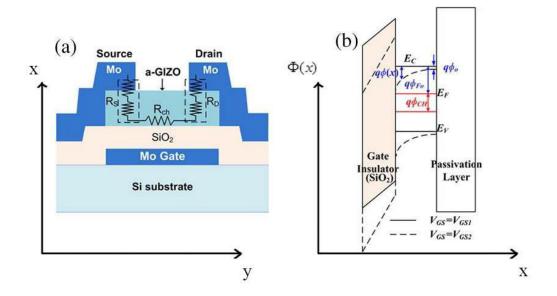
One can determine the field-effect mobility, or the channel mobility  $(\mu_n)$ , from Fig. 6 by using Eq. 2.7. Deviations from the gradual channel approximation model are often related to contact resistance effects and gate-voltage dependent mobility [3]. The field-effect mobility of the electrons in the channel of an a-Si:H TFT is only related to the 10% of the extended-state conduction-band mobility because of multiple carrier trapping in the band tail states.

# **2.1.2.** The model based on a mobility that changes with carrier concentrations in the channel for an a-IGZO TFT [17]

In this model, we provide a more accurate analytical model for an a-IGZO TFT. The geometry of the a-IGZO TFT is given in Fig. 3-a. In this model we do not assume a gradual channel. At any point in the channel the carrier concentration is dependent

on the potential  $\Phi(x)$ . The surface potential  $(\Phi_s)$  is changed with  $V_{GS}$  and  $V_{DS}$ . Amorphous - Indium Gallium Zinc Oxide (a-IGZO) is the active channel material in this transistor. We derived all the equations step by step clearly. Effective electron density  $(n_{EFF})$  is an important parameter in analytic DC model. In an n-type a-IGZO channel material, there are conduction band electrons  $(n_{FREE})$  and the carriers that hop under effect of the potential in the localized deep states within the conduction band  $(n_{LOC})$ . Because  $n_{EFF}$  is the effective concentration of electrons that contributes the channel conductivity. It is composed of  $n_{FREE}$  and  $n_{LOC}$ .

$$n_{EFF}(x) = n_{FREE}(x) + n_{LOC}(x)$$
(2.8)



**Figure 7** (a) A cross-sectional schematic of a-IGZO TFT. (b) Energy-band diagram with definitions of  $\Phi(x)$ ,  $\Phi_{CH}$  and  $\Phi_{FO}$  for the a-IGZO TFT [17]

In Fig. 3-a, the cross-sectional schematic of a-IGZO TFT with channel resistance  $(R_{CH})$  and parasitic resistances  $(R_p(V_{GS}, V_{DS}))$  is given. The energy – band diagram in Fig. 3 also shows how the potential changes in x direction  $(\Phi(x))$ , Fermi potential is lowered by the drain bias  $(\Phi_{CH})$  and Fermi potential under thermal equilibrium  $(\Phi_{FO})$ . Poisson's equation is expressed for potential  $\Phi(x)$  as

$$\frac{\partial^2 \Phi}{\partial x^2} = -\frac{\rho(x)}{\varepsilon_{IGZO}} = \frac{q n_{EFF}(x)}{\varepsilon_{IGZO}}$$
(2.9)

Eq. 2.9 defines the relation between charge density and potential.  $\rho(x)$  is charge density and  $\varepsilon_{IGZO}$  is permittivity of the a-IGZO active layer. When Eq. 2.8 and Eq.2.9 are combined, we can easily understand there is a direct relationship between charge density and free carriers and localized charges. Eq. 2.10 shows that relation clearly.

$$\frac{\rho(x)}{q} = -(n_{FREE}(x) + n_{LOC}(x)) = -n_{EFF}(x)$$
(2.10)

$$E_{IGZO} = -\frac{d\Phi}{dx}, \qquad (2.11)$$

where  $E_{IGZO}$  is the electric field generated by the applied voltage at x. Eq. 2.11 also gives important information about the relation of electric field with charge density. Eq. 2.9 and Eq. 2.11 are combined by using the algebraic rule given by

$$\frac{\partial \left[\frac{\partial f(x)}{\partial x}\right]}{\partial x} = 2 \left[\frac{\partial f(x)}{\partial x}\right] \left[\frac{\partial^2 f(x)}{\partial x^2}\right].$$
(2.12)

By using this in Eq. 2.12, Eq. 2.9 can be written by using Eq. 2.11 as

$$\frac{\partial \left[\frac{\partial \Phi}{\partial x}\right]^2}{\partial x} = 2 \left[\frac{\partial \Phi}{\partial x}\right] \left[\frac{\partial^2 \Phi}{\partial x^2}\right]$$
(2.13)

$$\frac{\partial}{\partial x} \left[ -\mathbf{E}_{IGZO} \right]^2 = 2 \left[ \frac{\partial \Phi}{\partial x} \right] \left( -\frac{\rho(x)}{\varepsilon_{IGZO}} \right). \tag{2.14}$$

Now, we are trying to find the electric field in terms of effective carrier density. We integrate both sides of Eq. 2.14.

$$\int \frac{\partial}{\partial x} [\mathbf{E}_{IGZO}]^2 dx = -\frac{2}{\varepsilon_{IGZO}} \int \rho(x) \frac{\partial \Phi}{\partial x} dx$$
(2.15)

$$E_{IGZO}^{2}(\Phi(x)) = \frac{2q}{\varepsilon_{IGZO}} \int n_{EFF}(x) d\Phi(x)$$
(2.16)

 $n_{EFF}(x)$  can be expressed in terms of potentials as

$$n_{EFF}(x) = N_{EFF} \exp\left[\frac{q(\Phi(x) - \Phi_{FO} - \Phi_{CH})}{kT_{EFF}}\right].$$
(2.17)

In Eq. 2.17,  $N_{EFF}$  is the effective Density of States (DOS) in cm<sup>-3</sup> and  $kT_{EFF}$  is characteristic slope in eVs. By combining Eq. 2.16 and Eq. 2.17, we can express the electric field in terms of potentials in Eq. 2.18.

$$E_{IGZO}^{2}(\Phi(x)) = \frac{2q}{\varepsilon_{IGZO}} \int N_{EFF} \exp\left[\frac{q(\Phi(x) - \Phi_{FO} - \Phi_{CH})}{kT_{EFF}}\right] d\Phi(x)$$
(2.18)

Finally, when we take the integral in Eq. 2.18, we find the electric field  $E_{IGZO}$  in terms of potentials.

$$E_{IGZO}^{2}(\Phi(x)) = \frac{N_{EFF}kT_{EFF}}{\varepsilon_{IGZO}} \exp\left[\frac{q(\Phi(x) - \Phi_{FO} - \Phi_{CH})}{kT_{EFF}}\right]$$
(2.19)

In next step, we derive the total charge by using the electric field that is given by Eq. 2.19 and total free charge. Total free charge is given in Eq. 2.20.

$$Q_{FREE}(\Phi(x)) = q \int_{x=0}^{x=T_{IGZO}} n_{FREE}(\Phi(x)) \, dx$$
 (2.20)

$$n_{FREE}(\Phi(x)) = N_{c} \exp\left[\frac{q(\Phi(x) - \Phi_{FO} - V_{CH}(y))}{kT}\right]$$
(2.21)

After we have combined Eq. 2.20 and Eq. 2.21, we are ready to find the total charge. In addition to that process, we combine Eq. 2.22 with Eq. 2.19.

$$Q_{FREE}(\Phi(x)) = qN_c \int_{x=0}^{x=T_{IGZO}} \exp\left[\frac{q(\Phi(x) - \Phi_{FO} - V_{CH}(y))}{kT}\right] dx$$
(2.22)

In Eq. 2.19 square of  $E_{IGZO}$  is given and electric field is also the negative derivative of  $\Phi(x)$  as it is written in Eq. 2.11. By using this equation, Eq. 2.23 is developed.

$$-\frac{d\Phi(x)}{dx} = \sqrt{\frac{2N_{EFF}kT_{EFF}}{\varepsilon_{IGZO}}} \exp\left[\frac{q(\Phi(x) - \Phi_{FO} - V_{CH})}{2kT_{EFF}}\right]$$
(2.23)

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Now, we obtain dx by using Eq. 2.23 to use that information in Eq. 2.22.

$$dx = -\frac{1}{\sqrt{\frac{2N_{EFF}kT_{EFF}}{\varepsilon_{IGZO}}}} \exp\left[\frac{-q(\Phi(x) - \Phi_{FO} - V_{CH})}{2kT_{EFF}}\right] d\Phi(x)$$
(2.24)

$$Q_{FREE}(\Phi(x)) = qN_c \int_{0}^{\Phi(x)} \exp\left[\frac{q(\Phi(x) - \Phi_{FO} - V_{CH}(y))}{kT}\right] \times \left(\sqrt{\frac{\varepsilon_{IGZO}}{2N_{EFF}kT_{EFF}}} \exp\left[\frac{-q(\Phi(x) - \Phi_{FO} - V_{CH})}{2kT_{EFF}}\right]\right) d\Phi(x)$$
(2.25)

When we take the integral in Eq. 2.25;

$$Q_{FREE}(\Phi(x)) = \frac{qN_c\sqrt{\varepsilon_{IGZO}}}{\sqrt{2N_{EFF}kT_{EFF}}\left[\frac{1}{kT} - \frac{1}{kT_{EFF}}\right]q}$$

$$\times \left(\exp\left[\left(\frac{1}{kT} - \frac{1}{kT_{EFF}}\right)\left(q\left(\Phi(x) - V_{CH} - \Phi_{FO}\right)\right)\right]\right)$$
(2.26)

We use the parameters in Eq. 2.27 to simplify the equation.

$$A^* = \frac{N_c \sqrt{\varepsilon_{IGZO}}}{B^* \sqrt{2N_{EFF} k T_{EFF}}} \qquad B^* = \left(\frac{1}{kT} - \frac{1}{2kT_{EFF}}\right) \tag{2.27}$$

$$Q_{FREE}(\Phi(x)) = A^* \exp[q B^*(\Phi(x) - V_{CH}(y) - \Phi_{FO})]$$
(2.28)

The short representation of free charge is given in Eq. 2.28. Total charge is the sum of free and localized charges. At the same time, it equals the product of permittivity and electric field as shown in Eq. 2.29 and Eq. 2.30.

$$Q_{TOT}(\Phi(x)) = Q_{FREE}(\Phi(x)) + Q_{LOC}(\Phi(x)) = \varepsilon_{IGZO} E_{IGZO}(\Phi(x))$$
(2.29)

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$$Q_{TOT}(\Phi(x)) = \sqrt{2\varepsilon_{IGZO}N_{EFF}kT_{EFF}} \exp\left[\frac{q(\Phi(x) - V_{CH}(y) - \Phi_{FO})}{2kT_{EFF}}\right]$$
(2.30)

By using Eq. 2.28 and Eq. 2.30 in Eq. 2.31,  $V_{GS}$  dependent channel mobility ( $\mu_{CH}$ ) can be found.

$$\mu_{CH}(\Phi(x)) = \mu_{BAND} \frac{Q_{FREE}(\Phi(x))}{Q_{FREE}(\Phi(x)) + Q_{LOC}(\Phi(x))}$$
(2.31)

 $\mu_{BAND}$  is the conduction band mobility in Eq. 2.31. We insert the values given in Eq. 2.28 and Eq. 2.30 into Eq. 2.31

$$\mu_{CH}(\Phi(x)) = \mu_{BAND} \frac{A^* \exp[qB^*(\Phi(x) - V_{CH}(y) - \Phi_{FO})]}{\sqrt{2\varepsilon_{IGZO}N_{EFF}kT_{EFF}}} \exp\left[\frac{q(\Phi(x) - V_{CH}(y) - \Phi_{FO})}{2kT_{EFF}}\right]$$
(2.32)

$$\mu_{CH}(\Phi(x)) = \mu_{BAND} \frac{N_c}{\left(\frac{1}{kT} - \frac{1}{2kT_{EFF}}\right) (2N_{EFF}kT_{EFF})}$$

$$\times \exp\left[\left(\frac{1}{kT} - \frac{1}{kT_{EFF}}\right) q \left(\Phi(x) - V_{CH}(y) - \Phi_{FO}\right)\right]$$
(2.33)

Now, we describe drain to source current  $(I_{DS})$  in terms of  $\mu_{CH}(\Phi(x))$  and other terms in  $\mu_{CH}(\Phi(x))$  equation. The drain to source current equation is

$$I_{DS} = W \frac{dV_{CH}}{dy} \int_{\Phi(x=T_{IGZO})}^{\Phi(x=0)} q\mu_{CH} \frac{n_{FREE}(\Phi(x))}{E_{IGZO}(\Phi(x))} d\Phi(x) .$$
(2.34)

We take the integral in Eq. 2.34.  $\mu_{CH}$  is given above in Eq. 2.33.  $n_{FREE}(\Phi(x))$  can be written by using Eq. 2.21, and  $E_{IGZO}(\Phi(x))$  can be found by using Eq. 2.19.

$$I_{DS} = W \frac{dV_{CH}}{dy} \int_{\Phi(x=T_{IGZO})}^{\Phi(x=0)} q\mu_{BAND} \frac{N_c}{(2N_{EFF}kT_{EFF})B^*} \frac{N_c \sqrt{\varepsilon_{IGZO}}}{\sqrt{2N_{EFF}kT_{EFF}}}$$
(2.35)  
  $\times \exp\left[\left(\frac{1}{kT} + \frac{1}{kT} - \frac{1}{kT_{EFF}} - \frac{1}{2kT_{EFF}}\right)q(\Phi(x) - V_{CH}(y) - \Phi_{FO})\right]d\Phi(x)$ 

To simplify, we use Eq. 2.27 and Eq. 2.36.

$$C^* = \left(\frac{2}{kT} - \frac{3}{2kT_{EFF}}\right) \tag{2.36}$$

$$I_{DS} = W \frac{dV_{CH}}{dy} q \mu_{BAND} \left( \frac{A^* N_c}{2N_{EFF} k T_{EFF}} \right)$$

$$\times \int_{\Phi(x=T_{IGZO})}^{\Phi(x=0)} \exp\left[q C^* (\Phi(x) - V_{CH} - \Phi_{FO})\right] d\Phi(x)$$
(2.37)

If we take the integral we get

$$I_{DS} = W \mu_{BAND} \frac{dV_{CH}}{dy} \left( \frac{A^* N_c}{2N_{EFF} k T_{EFF} C^*} \right) \exp \left[ q C^* (\Phi(x) - V_{CH} - \Phi_{FO}) \right].$$
(2.38)

In Eq. 2.38, we can clearly see that  $\Phi(x = T_{IGZO}) = V_{CH} + \Phi_{FO}$  and  $\Phi(x = 0) = \Phi_s$ . We have found a near analytic DC I-V model. But, the surface potential  $(\Phi_s)$  also depends on the Fermi level lowering effect along the channel [18]. So, we should find a relation between surface potential and the channel voltage  $(V_{CH})$ . At first, we derive the relation between  $I_{DS}$  and  $V_{GS}$  and  $V_{DS}$ . We will apply Gauss Law to the boundary between a-IGZO active layer and the gate oxide. This operation gives us a nonlinear relation between  $V_{GS}$  and  $\Phi_S$ 

$$V_{GS} = V_{FB} + \Phi_S + \frac{\varepsilon_{IGZO} E_{IGZO} (\Phi_S, V_{CH})}{C_{OX}}.$$
(2.39)

Flat band voltage  $(V_{FB})$  is of constant value, and surface potential  $(\Phi_s)$  is variable. Electric field can be found by using Eq. 2.19.

$$\frac{\varepsilon_{IGZO}E_{IGZO}(\Phi_{S}, V_{CH})}{C_{OX}} = \frac{\sqrt{2N_{EFF}kT_{EFF}}\varepsilon_{IGZO}}{C_{OX}}\exp\left(\frac{q(\Phi_{S}-V_{CH}-\Phi_{FO})}{2kT_{EFF}}\right)$$
(2.40)

By using Eq. 2.40, Eq. 2.39 can be expressed as

$$\frac{\left(V_{GS} - V_{FB} - \Phi_{S}\right)C_{OX}}{\sqrt{2\varepsilon_{IGZO}N_{EFF}kT_{EFF}}} = \exp\left(\frac{q(\Phi_{S} - V_{CH} - \Phi_{FO})}{2kT_{EFF}}\right)$$
(2.41)

Now, we express  $V_{CH}$  in terms of surface potential( $\Phi_s$ ). The relation between  $V_{CH}$  and  $\Phi_s$  is nonlinear

$$\Phi_{S} - V_{CH} - \Phi_{FO} = \frac{2kT_{EFF}}{q} \ln\left[\frac{C_{OX}(V_{GS} - V_{FB} - \Phi_{S})}{\sqrt{2\varepsilon_{IGZO}N_{EFF}kT_{EFF}}}\right].$$
(2.42)

Now, we take the derivative of channel voltage and use that equation in Eq. 2.38.

$$\frac{dV_{CH}(y)}{d\Phi_{s}(y)} = 1 + \frac{2kT_{EFF}}{q(V_{GS} - V_{FB} - \Phi_{s})}$$
(2.43)

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$$\frac{dV_{CH}(y)}{dy} = \frac{dV_{CH}(y)}{d\Phi_s(y)} \frac{d\Phi_s(y)}{dy} = \left[1 + \frac{2kT_{EFF}}{q(V_{GS} - V_{FB} - \Phi_s)}\right] \frac{d\Phi_s(y)}{dy}$$
(2.44)

When we put Eq. 2.42 and Eq. 2.44 into Eq. 2.38, we find a better expression for drain to source current. Eq. 2.45 is giving that equation.

$$I_{DS} = W \mu_{BAND} \left[ 1 + \frac{2kT_{EFF}}{q(V_{GS} - V_{FB} - \Phi_s)} \right] \frac{A^* N_c}{2N_{EFF} k T_{EFF} C^*} \times \left[ \frac{C_{OX} \left( V_{GS} - V_{FB} - \Phi_s \right)}{\sqrt{2\varepsilon_{IGZO} N_{EFF} k T_{EFF}}} \right]^{2kT_{EFF} C^*} \frac{d\Phi_s(y)}{dy}$$

$$(2.45)$$

Finally, when we take the integral in y direction, the drain current is derived. Integrating in y direction from 0 to L,  $\Phi_{SS} = \Phi_S(y = 0, V_{CH} = 0)$  and  $\Phi_{SD} = \Phi_S(y = L, V_{CH} = V_{DS}) I_{DS}$  is obtained.

$$\int_{y=0}^{y=L} I_{DS} dy = \int_{\Phi_{SS}}^{\Phi_{SD}} W \mu_{BAND} \left[ 1 + \frac{2kT_{EFF}}{q(V_{GS} - V_{FB} - \Phi_{S})} \right]$$

$$\times \left( \frac{A^{*}N_{c}}{2N_{EFF}kT_{EFF}C^{*}} \left[ \frac{C_{OX}\left(V_{GS} - V_{FB} - \Phi_{S}\right)}{\sqrt{2\varepsilon_{IGZO}N_{EFF}kT_{EFF}}} \right]^{2kT_{EFF}C^{*}} \right] d\Phi_{S}(y)$$
(2.46)

$$I_{DS}(N_{EFF}, kT_{EFF}) = \frac{W}{L} \mu_{BAND} \frac{A^* N_c}{2N_{EFF} kT_{EFF} C^*} \left( \frac{C_{OX}}{\sqrt{2\varepsilon_{IGZO} N_{EFF} kT_{EFF}}} \right)^{2kT_{EFF} C^*} \times \left\{ \left( 2kT_{EFF} C^* + 1 \right)^{-1} \times \left( \left( V_{GS} - V_{FB} - \Phi_{SS} \right)^{2kT_{EFF} C^* + 1} - \left( V_{GS} - V_{FB} - \Phi_{SS} \right)^{2kT_{EFF} C^* + 1} \right) - \left( qC^* \right)^{-1} \times \left( \left( V_{GS} - V_{FB} - \Phi_{SS} \right)^{2kT_{EFF} C^*} - \left( V_{GS} - V_{FB} - \Phi_{SD} \right)^{2kT_{EFF} C^*} \right) \right\}$$

$$(2.47)$$

Final version of the equation gives the relation among drain current, gate to source voltage and drain to source voltage.  $\mu_{BAND}$  is an important parameter on which the drain current is directly proportional.

If there are two different parameters for sub and above of threshold region, the process is shown above should be done for these two regions. To find the total  $I_{DS}$  value, we will use

$$I_{DS\_TOT} = \frac{1}{I_{DS\_sub}(N_{EFF1}, kT_{EFF1})} + \frac{1}{I_{DS\_above}(N_{EFF2}, kT_{EFF2})}$$
(2.48)

We have calculated all values of  $I_{DS_{-TOT}}$  by using the formula given in Eq. 2.47. There are many parameters in those equations. Those parameter values are given in [18].

Parameter	Value	Parameter	Value	
W / L [μm]	225/30	$N_{\rm EFF1} \ [{ m cm}^{-3}]$	$8.7 \times 10^{17}$	
$L_{\rm OV}$ [µm]	5			
$T_{\rm OX} / T_{\rm IGZO} $ [nm]	100/50	$N_{\rm EFF2}~[{\rm cm}^{-3}]$	$5.15 \times 10^{18}$	
$\mu_{\rm BAND}  [{\rm cm}^2 /  {\rm V.s}]$	19.7	$kT_{\rm EFF2}$ [eV]	0.0263	
$N_{c}  [{\rm cm}^{-3}]$	$4.8 \times 10^{18}$	$C_{\rm ox} [\rm nF/cm^2]$	42.37	
V <sub>FB</sub> [V]	0.3	$\phi_{\rm F0}  [{ m V}]$	0.3	

 Table 2 Geometrical and extracted model parameters [18]

By using those values given in Table 2.1, we obtained 3 similar figures. All these 3 figures given below show the relation between  $I_{DS}$ ,  $V_{GS}$ , and  $V_{DS}$ . During that process, we have used Eq. 2.39 to calculate  $\Phi_{SS}$  and  $\Phi_{SD}$  values. The steps to obtain the individual values for  $I_{DS}$  by changing  $V_{GS}$  and  $V_{DS}$  values are given in the flow-chart below.

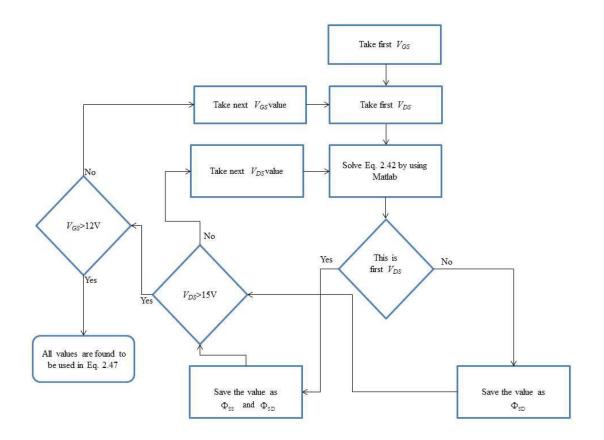


Figure 8  $\Phi_{_{SS}}$  and  $\Phi_{_{SD}}$  calculation process

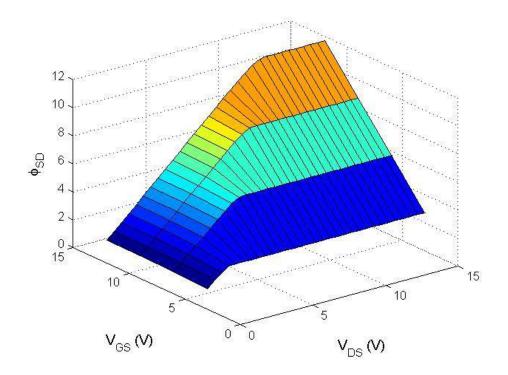
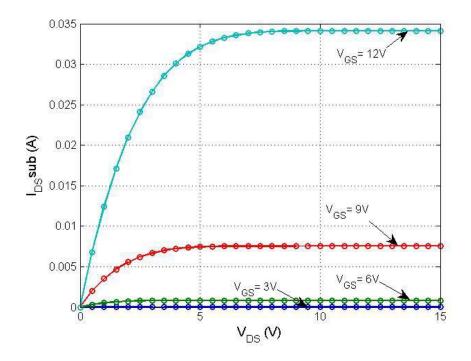


Figure 9  $\Phi_{\rm SD}$ ,  $V_{\rm GS}$  and  $V_{\rm DS}$  relation

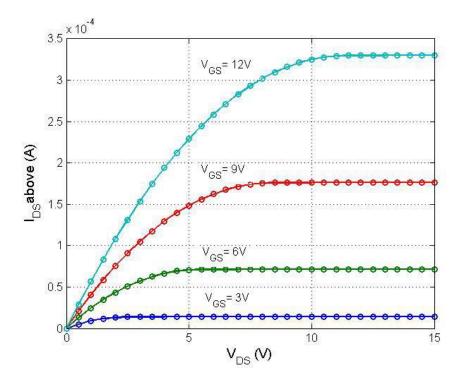
The dependence of  $\Phi_{SD}$  on  $V_{GS}$  and  $V_{DS}$  is shown in Fig. 9. As can be seen in the figure,  $V_{DS}$  is between from 0 and 15V. However,  $\Phi_{SS}$  is obtained for  $V_{DS} = 0$ .

To verify if we derived everything correctly and properly we calculated the  $I_{DS}$  versus  $V_{DS}$  values with  $V_{GS}$  as parameters by using the values given in Table 2.1. and compared all these results with Minkyung Bae et al. [18]. The Output Characteristics that we obtain are shown in Fig. 10 to 12. As seen our values fit very well with their results. Therefore we confirm that all the derivations we have made are correct and also our flowchart for solving the static characteristics for the a-IGZO TFT is working.

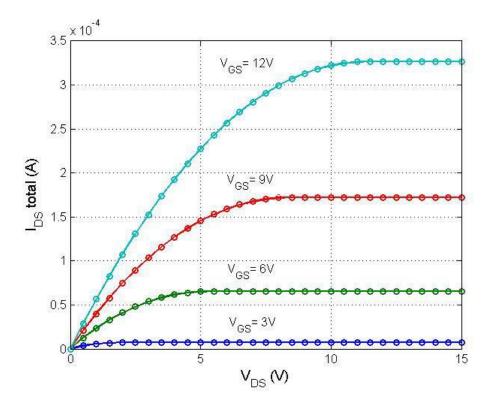
As seen in Fig. 12 this model [18] works better than the gradual channel approximation because in his model one equation simulates the behavior of the transistor in all the regions (sub-threshold, and above threshold regions (linear and saturation regions)). However, in this model, solving  $\Phi_s$  in terms of  $V_{GS}$  and  $V_{DS}$  in the nonlinear equation of 2-42, needs a numerical root-finding technique. In that sense this model of [18] can be more precisely a semi-analytic model.



**Figure 10** Drain to source current curves in sub-threshold region for the TFT having the parameters listed in Table 2.1 (lines are used for guiding eyes only)



**Figure 11** Drain to source current curves in above-threshold region for the TFT having the parameters listed in Table 2.1 (lines are used for guiding eyes only)



**Figure 12** Total drain to source current curves (both sub and above threshold regions) for the TFT having the parameters listed in Table 2.1 (lines are used for guiding eyes only)

Fig. 12 is the same with the graphics that is given in [18]. We have produced these graphs by applying the flow chart above. If noticed, Fig. 11 and Fig. 12 look very similar. Since the current values that are found in the sub-threshold region for the same  $V_{GS}$  and  $V_{DS}$  values are higher than those obtained in the above-threshold region, considering Eq. 2.48, smaller  $I_{DS}$ 's are dominant in determining the overall  $I_{DS}$ 's.

When we verified that our equations, flowchart and the method of calculation are correct and accurate in calculating the static characteristic curves of the a-IGZO TFT we decided to check the effects of some parameters in Table 2.1 on the calculated values. Since the values like  $N_{EFF1}$ ,  $kT_{EFF1}$ ,  $N_{EFF2}$ ,  $kT_{EFF2}$  are estimated and used in the model as inputs we decided to analyze the sensitivity of the model on these parameters. We changed their values only 5% and generated the same curves as Fig. 12 for these values. When we changed  $N_{EFF1}$  and  $kT_{EFF1}$ , the density of the tail states and the activation energy of these states, respectively, 5% form their values in Table 2.1 the output characteristics did not change at all. However, a 5% change in  $N_{\rm EFF2}$ and  $kT_{EFF2}$ , the density of deep states and the activation energy of these states, respectively, from their values given in Table 2.1 the output characteristics change considerably. Fig. 13 and 14 display the changes in the characteristics for a 5% change in  $N_{\rm EFF2}$  and 5% change in  $kT_{\rm EFF2}$ . Form the figures one can easily see that the drain current values increase nearly 20% for a 5% decrease in  $N_{\rm EFF2}$  (when we take  $N_{EFF2}$  4.89x10<sup>18</sup> cm<sup>-3</sup> instead of 5.15 x10<sup>18</sup> cm<sup>-3</sup>); whereas, the drain current values decrease nearly 1100% for a 5% decrease in  $kT_{\rm EFF2}$  (when we take  $kT_{\rm EFF2}$  $0.0250 \ eV$  instead of  $0.0263 \ eV$ ). Therefore, we believe that such level of error in estimating these values could be made very easily, and the sensitivity of the model to small changes in these values is the weakness of the model.

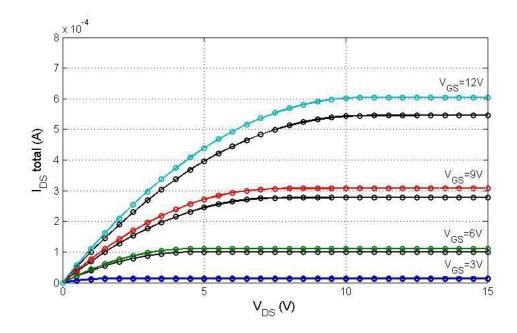


Figure 13 Changes in the characteristics for a 5% change in  $N_{EFF2}$  (curves in black color are the original curves, and the other curves for various colors are for decreased

 $N_{EFF2}$  value)

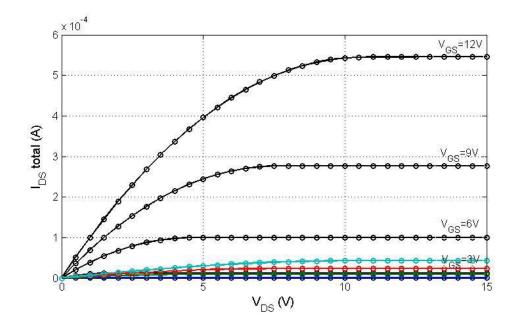


Figure 14 Changes in the characteristics for a 5% change in  $kT_{EFF2}$  (curves in black color are the original curves, and the other curves for various colors are for decreased

$$kT_{EFF2}$$
 value)

### **CHAPTER 3**

# 3. NEW SURFACE POTENTIAL-BASED ANALYTIC IGZO TFT MODEL

The Gradual Channel Approximation that we reviewed in Chapter 2 in modeling the FET, assuming constant channel mobility, is very practical because it is very simple, analytical, and quite successfully fits the empirical FET I-V behavior. However, it has different expressions for the different regions of the FET operation (one formula for the linear region, and another formula for the saturation region). On the other hand, it does not include any expression for the sub-threshold region at all. The same model has been used for a-Si:H TFTs by many authors. However, one analytical model to estimate all regions of the FET operation has always been needed.

Actually, the first analytical model for an a-Si:H TFT was introduced by Hack and Shur [41]. They obtained the channel conductance by a Taylor Series expansion, where they again developed different formulas for different operating regions of the transistor. They did not have one formula valid for all the regions of the TFT.

A unique formulation for the sub-threshold, above-threshold (linear and saturation) regions of an TFT, which accounts for a realistic distribution of tail and deep density of states (DOS) in the energy gap was developed by Colalongo [42]. This model is based on the charge-sheet analytical model of Brews for a crystalline metal-oxide-

semiconductor FET [43]. Colalongo adapted and generalized this model for a-Si:H TFTs to include the effect of trap states, namely tail and deep states, within the bandgap of hydrogenated amorphous silicon material. His model includes the relationship between the surface and the quasi-Fermi potential. One of the advantages of the model is that explicit definitions of the threshold and saturation voltages as input parameters, which cannot be accurately defined for these devices, are not required. In his model, he used constant channel mobility in calculating the drain current. However, in fitting his equations with the experimental data he took a mobility value of  $10 \text{ cm}^2 / Vs$  which is very high and seems not attainable for an a-Si:H TFT [42]. If one considers that a-Si:H mobilities are less than  $1 \text{ cm}^2 / Vs$ , that is an order less than what he used, one concludes that his model underestimates the drain current values.

At this point we believe that we should first talk about the mobility of the charge carriers in the FET's channel. Mobility is the ease with which the charge carriers in the material move under the effect of a force. In the FETs we have two kinds of mobility: the field effect mobility and the band mobility. The field effect mobility  $(\mu_{EFF})$  in the induced channel, sometimes called as the channel mobility  $(\mu_{CH})$ , of an FET of a crystalline semiconductor is equal to the conduction band mobility  $(\mu_{EAND})$ , which is generally constant, because in a crystalline semiconductor the charge carriers are the free carriers in the conduction band  $(n_{FREE})$ . These carriers are induced in the channel with the applied gate to source voltage  $(V_{GS})$ . However, in the TFTs the channel materials are generally non-crystalline. These non-crystalline materials are either amorphous semiconductors (a-Si:H) or polycrystalline

semiconductor (poly-Si), or metal oxides (*ZnO*, *CdSe* or a-IGZO), which are amorphous too. In non-crystalline semiconductors the charge carriers, for example electrons, are not only those from the conduction  $band(n_{FREE})$ , but those from the localized states within the bandgap  $(n_{LOC})$  as well. The field effect mobility in noncrystalline semiconductors is, thus, generally a function of the channel potential  $V_{CH}(y)$ , and given in one approach [19,44,45,46] as we used in Chapter 2 in deriving their formulation in detail:

$$\mu_{EFF} = \mu_{BAND} \frac{Q_{FREE}(\mathbf{x}, \mathbf{V}_{CH}(\mathbf{y}))}{Q_{FREE}(\mathbf{x}, \mathbf{V}_{CH}(\mathbf{y})) + Q_{LOC}(\mathbf{x}, \mathbf{V}_{CH}(\mathbf{y}))}$$
(2.31)

where  $Q_{FREE}$  is the total free charge density, and  $Q_{LOC}$  is the total localized charge density in the channel, x being the transverse distance from the gate oxide surface,  $V_{CH}(y)$  is the channel potential at y, y being the distance from source to drain.  $V_{CH}(y)$  is determined by  $V_{GS}$  and  $V_{DS}$ . In a-Si:H material since  $Q_{LOC} \gg Q_{FREE}$  the field effect mobility is much less than the band mobility, and therefore the mobility in a-Si:H TFTs channel,  $\mu_{EFF}$ , is nearly constant or a weak function of  $V_{CH}(y)$  [44]. However, in poly-Si or in a-IGZO the total charge of free carriers are comparable with the total charge of the localized carriers or even larger, and the field effect mobility can get a values that are somewhat closer to the band mobility. Moreover, in these TFTs the field effect mobility is a strong function of  $V_{CH}(y)$ .

Eq. 2.31 is an empirical approximation introduced to explain the difference of the field effect mobility from the band mobility in a TFT. This formula has been used by

some [18, 19] in their model for the a-IGZO TFTs for a quite wide range of I-V characteristics. Actually, in the model we analyzed in detail in Chapter 2, Minkyung Bae et al. [18] used the same approach to incorporate the tail and deep states in the band gap with Colalongo [42] in deriving the drain current for sub-threshold and above-threshold regions; however, since a-IGZO has free carriers contributing to the channel conductivity considerably, they used  $n_{FREE}$  as well in addition to the localized states and also they, like Colalongo, introduced a channel potential dependence on the channel mobility.

In some other works, the mobility's dependence on channel potential  $V_{CH}(y)$  is modeled by another formula [45,47-50]

$$\mu_{EFF} = \mu_{BAND} \left( \frac{V_{GS} - V_T}{V_{AA}} \right)^{\gamma}$$
(3.1)

where  $V_T$  is the threshold voltage,  $V_{AA}$  is a fitting voltage and  $\gamma$  is another fitting parameter. It was used with a success to some extent for a-Si:H TFTs, since their mobilities are weakly dependent on  $V_{CH}(y)$ . This formula was not preferred for the a-IGZO TFTs because in those TFTs mobilities are somewhat stronger function of  $V_{CH}(y)$ .

In his work for TFTs Tze-Ching Fung [21] used the gradual channel approximation but with a gate-to-source voltage dependent mobility to model the a-IGZO TFT. He introduced the channel mobility as

$$\mu_{EFF}(V_{GS}) = \mu_0 \left[ \frac{V_{GS} - V_T}{V_C} \right]^{\alpha}$$
(3.2)

where  $\mu_0$  is the intrinsic band mobility,  $V_C$  is the material-dependent critical voltage, and  $\alpha$  is a coefficient which describes the dependence of mobility  $\mu_{EFF}$  on  $(V_{GS} - V_T)$ . Then, in the simplified form he expressed the mobility as

$$\mu_{EFF}\left(V_{GS}\right) = K \left(V_{GS} - V_T\right)^{\gamma - 1} \tag{3.3}$$

where K is a material dependent parameter as  $K = \mu_0 (1/V_c)^{\alpha}$ , and  $\gamma = \alpha + 1$ .

The drain current in the Gradual Channel Approximation, using a voltage-dependent mobility in the linear region can be expressed as

$$I_D = K Cox(W/L) (V_{GS} - V_T)^{\gamma} V_{DS}$$
(3.4)

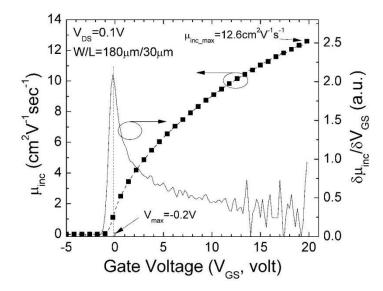
He also defined the incremental mobility

$$\mu_{inc} = \left(\frac{\delta I_D}{\delta V_{GS}}\right) \left[\frac{L}{WCoxV_{DS}}\right]$$
(3.5)

where L is the channel length and W is the channel width of the TFT. And, from Eq. 3.4 and Eq. 3.5 he obtained the incremental mobility as

$$\mu_{inc} = K \ \gamma \ \left( V_{GS} - V_T \right)^{\gamma - 1} \tag{3.6}$$

Fig. 15 shows the incremental mobility dependence on the gate-to-source voltage extracted from the measured  $I_D$  versus  $V_{GS}$  curve from an a-IGZO TFT [21].



**Figure 15** Incremental field effect mobility  $(\mu_{inc})$  of an a-IGZO TFT [21]

The extracted  $\mu_{inc}$  is a strong function of  $V_{GS}$  and reaches a maximum value of 12.6  $cm^2/Vs$ , and  $\mu_{inc}$  does not saturate. Such behavior was also observed in other oxide semiconductors [51,52]. The  $\gamma$  value here extracted, by linear fitting, from the experimental curves is equal to 1.52.

As we redeveloped all the equations for the model in [18] in Chapter 2 again and obtained the values of  $I_{DS}$  versus  $V_{GS}$  and  $V_{DS}$  by using their parameters given in their manuscript, we showed that the model using Eq. 2.31 is very sensitive to parameters  $N_{EFF2}$ 's and  $kT_{EFF2}$ , that are either extracted from the experimental characteristics or estimated. We noticed that even very small change on these parameters cause very big changes in the I-V curves. For example a change in  $N_{EFF2}$  from  $5.15 \times 10^{18}$  to  $4.89 \times 10^{18}$ , causes a 20% increase in  $I_{DS}$ , whereas,  $I_{DS}$  values

decrease nearly 1100% for a 5% decrease in  $kT_{EFF2}$  (when we take  $kT_{EFF2}$  0.0250 eV instead of 0.0263 eV). Since it is nearly impossible to experimentally determine or estimate these parameters about this level of accuracy, and the model's success is solely dependent on selecting these parameters that accurate in the beginning, we thought that we could introduce another model for mobility. By considering the results and findings of [21] about the dependence of the mobility on the  $V_{GS}$ , and by considering Eq. 2.31 for the mobility dependence on the channel voltage  $V_{CH}(y)$  in their model [18], and by comparing their results with the measured values given, we think that a new formula for the mobility dependence on the channel voltage, and an average of all the carriers contributing to the conduction process in the channel (carriers from free and localized states) would produce a better model for an a-IGZO TFT.

Towards this goal, we first checked the empirical formula Eq. 3.1 to estimate the I-V characteristics for an a-IGZO TFT. We tried a range of constant values for either  $V_{AA}$  and for  $\gamma$  and saw that they were not effective to fit the experimental I-V values of the TFT. Therefore, we introduced a new formula inspired from this equation as:

$$\mu_{EFF} = \mu_{\text{BAND}} \left[ \frac{\left( V_{GS} - V_T \right)}{V_{AA}} \right]^{\gamma(V_{GS}, V_{DS})}$$
(3.7)

Here  $\gamma$  is not constant as in Eq. 3.3 but a function of  $V_{GS}$  and  $V_{DS}$ , and  $V_{AA}$  is a fix voltage to provide unit consistency. We used the same approach as given in [5,18,42] developing our model to estimate the I-V characteristics of an a-IGZO TFT, not by using Eq. 2.31 as they did in [18] but by using our formula Eq. 3.7 for the effective

mobility. This  $\gamma$ , to be determined in mobility equation, Eq. 3.7, is not a constant as in Eq. 3.1 but a function of both  $V_{GS}$  and  $V_{DS}$ , that is  $V_{CH}(y)$ .

Jun Hyung Park et al. introduced another model [17], where they also used Eq. 2.31 for mobility but instead of differentiating tail states and deep states for the localized states and used  $N_{EFF}$ 's ( $N_{EFF1}$ ,  $N_{EFF2}$ ) and characteristic slopes (activation energies) for each of them ( $kT_{EFF1}$ ,  $kT_{EFF2}$ ) as Minkyung Bae et al., they took only one localized carriers of  $N_{EFF}$  and one characteristic slope  $kT_{EFF}$ . They even went one step further and represented all carriers in the channel with effective carriers ( $n_{EFF}$ ), which is the sum of the free carriers ( $n_{FREE}$ ) from the conduction band and localized carriers ( $n_{LOC}$ ) from the states in the bandgap. Although we did not use their approach for mobility, we used the same approach of representing all the carriers contributing conduction in the channel as effective carriers ( $n_{EFF}$ ).

Since the potential in the TFT's channel is not only determined by  $V_{GS}$  but  $V_{DS}$  as well, the mobility should also dependent on  $V_{GS}$  and  $V_{DS}$ . We believe that our formula, Eq. 3.7, for mobility in this sense is more realistic in representing the channel mobility than the other models cited above.

#### 3.1. Developing Our Model for an a-IGZO TFT

As we stated above we start with a new formula for the channel mobility to derive our model

$$\mu_{CH} = \mu_{BAND} \left( \frac{V_{GS} - V_T}{V_{AA}} \right)^{\gamma}$$
(3.8)

We assume that there is no dependence of  $\mu_{CH}$  on geometry. In other words, if there is a specific difference of TFT in geometrical features, this will be reflected into given parameters.

Eq. 2.34 gives the relation between effective channel mobility and  $I_{DS}$ . So, we combine Eq. 2.34 and Eq. 3.8

$$I_{DS} = W \frac{dV_{CH}}{dy} \int_{\Phi(x=T_{IGZO})}^{\Phi(x=0)} q \mu_{BAND} \left(\frac{V_{GS} - V_T}{V_{AA}}\right)^{\gamma} \frac{n_{FREE}(\Phi(x))}{E_{IGZO}(\Phi(x))} d\Phi(x)$$
(3.9)

By using the information given in Eq. 2.19, we can derive

$$E_{IGZO}(\Phi(x)) = \sqrt{\frac{2N_{EFF}kT_{EFF}}{\varepsilon_{IGZO}}} \exp\left[\frac{q(\Phi(x) - \Phi_{FO} - \Phi_{CH})}{2kT_{EFF}}\right].$$
(3.10)

Eq. 2.21 provides information about  $n_{FREE}(\Phi(x))$ . Using eq.2.21, and Eq.3.10 in Eq.3.9 we can obtain

$$I_{DS} = Wq\mu_{BAND} \left(\frac{V_{GS} - V_T}{V_{AA}}\right)^{\gamma} \frac{dV_{CH}}{dy} \times \int_{\Phi(x=T_{IGZO})}^{\Phi(x=0)} \frac{N_c \exp\left[\frac{q(\Phi(x) - \Phi_{FO} - V_{CH}(y))}{kT}\right]}{\sqrt{\frac{2N_{EFF}kT_{EFF}}{\varepsilon_{IGZO}}} \exp\left[\frac{q(\Phi(x) - \Phi_{FO} - \Phi_{CH})}{2kT_{EFF}}\right]} d\Phi(x)$$
(3.11)

While we are recombining Eq. 3.5 to make it easier to take an integral, we will use Eq. 2.27 to simplify the equation. When we make these changes

$$I_{DS} = \frac{Wq\mu_{BAND}N_c}{\sqrt{\frac{2N_{EFF}kT_{EFF}}{\varepsilon_{IGZO}}}} \left(\frac{V_{GS} - V_T}{V_{AA}}\right)^{\gamma} \frac{dV_{CH}}{dy} , \quad (3.12)$$
$$\times \int_{\Phi(x=T_{IGZO})}^{\Phi(x=0)} \exp[qB^*(\Phi(x) - \Phi_{FO} - V_{CH}(y))] d\Phi(x)$$

where 
$$B^* = \left(\frac{1}{kT} - \frac{1}{2kT_{EFF}}\right)$$
.

The surface potential  $\Phi(x=0) = \Phi_s$ . In addition to this, we use the relation given in Eq. 2.44 to convert the integral dependence from  $\Phi_s$  to y.

$$I_{DS} = \frac{Wq\mu_{BAND}N_c}{B^*\sqrt{\frac{2N_{EFF}kT_{EFF}}{\varepsilon_{IGZO}}}} \left(\frac{V_{GS} - V_T}{V_{AA}}\right)^{\gamma} \left[1 + \frac{2kT_{EFF}}{q(V_{GS} - V_{FB} - \Phi_S)}\right]$$

$$\times \exp[qB^*(\Phi_S - \Phi_{FO} - V_{CH}(y))] \frac{d\Phi_S(y)}{dy}$$
(3.13)

By using Eq. 2.42, we can convert  $I_{DS}$  equation into Eq. 3.14.

$$I_{DS} = \frac{Wq\mu_{BAND}N_c}{B^*\sqrt{\frac{2N_{EFF}kT_{EFF}}{\varepsilon_{IGZO}}}} \left(\frac{V_{GS} - V_T}{V_{AA}}\right)^{\gamma} \left[1 + \frac{2kT_{EFF}}{q(V_{GS} - V_{FB} - \Phi_S)}\right]$$

$$\times \exp\left\{B^*2kT_{EFF}In\left[\frac{C_{OX}(V_{GS} - V_{FB} - \Phi_S)}{\sqrt{2\varepsilon_{IGZO}N_{EFF}kT_{EFF}}}\right]\right\} \frac{d\Phi_S(y)}{dy}$$
(3.14)

We got rid of  $V_{CH}(y)$  by this way. After making some simplifications in Eq. 3.14., we obtain

$$I_{DS} = \frac{Wq\mu_{BAND}N_c}{B^*\sqrt{\frac{2N_{EFF}kT_{EFF}}{\varepsilon_{IGZO}}}} \left(\frac{V_{GS} - V_T}{V_{AA}}\right)^{\gamma} \left[1 + \frac{2kT_{EFF}}{q(V_{GS} - V_{FB} - \Phi_S)}\right] \times \left[\frac{C_{OX}(V_{GS} - V_{FB} - \Phi_S)}{\sqrt{2\varepsilon_{IGZO}N_{EFF}kT_{EFF}}}\right]^{\left(\frac{2kT_{EFF}}{kT} - 1\right)} \frac{d\Phi_S(y)}{dy}$$
(3.15)

If we take the integral along the channel (from y = 0 to y = L)

$$I_{DS} \int_{0}^{L} dy = \frac{Wq\mu_{BAND}N_{c}}{B^{*}\sqrt{\frac{2N_{EFF}kT_{EFF}}{\varepsilon_{IGZO}}}} \left(\frac{V_{GS} - V_{T}}{V_{AA}}\right)^{\gamma} \times \left\{ \int_{\Phi_{S}(y=0)}^{\Phi_{S}(y=L)} \left[\frac{C_{OX}(V_{GS} - V_{FB} - \Phi_{S})}{\sqrt{2\varepsilon_{IGZO}N_{EFF}kT_{EFF}}}\right]^{\left(\frac{2kT_{EFF}}{kT} - 1\right)} d\Phi_{S}(y) \right. (3.16) + \int_{\Phi_{S}(y=0)}^{\Phi_{S}(y=L)} \left[\frac{2kT_{eff}}{q(V_{GS} - V_{FB} - \Phi_{S})}\right] \left[\frac{C_{OX}(V_{GS} - V_{FB} - \Phi_{S})}{\sqrt{2\varepsilon_{IGZO}N_{EFF}kT_{EFF}}}\right]^{\left(\frac{2kT_{EFF}}{kT} - 1\right)} d\Phi_{S}(y) \right\}$$

Here we have a sum of two integrals, which can be called as  $I_1$  and  $I_2$ .

$$I_{DS} \int_{0}^{L} dy = \frac{Wq\mu_{BAND}N_c}{B^* \sqrt{\frac{2N_{EFF}kT_{EFF}}{\varepsilon_{IGZO}}}} \left(\frac{V_{GS} - V_T}{V_{AA}}\right)^{\gamma} \times \left(I_1 + I_2\right)$$
(3.17)

When we take the first integral in Eq. 3.16, the result is

$$I_{1} = \left(\frac{C_{OX}}{\sqrt{2\varepsilon_{IGZO}N_{EFF}kT_{EFF}}}\right)^{\left(\frac{2kT_{EFF}}{kT}-1\right)} \left(\frac{kT}{2kT_{EFF}}\right) \times \left[\left(V_{GS}-V_{FB}-\Phi_{SS}\right)^{\left(\frac{2kT_{EFF}}{kT}\right)}-\left(V_{GS}-V_{FB}-\Phi_{SD}\right)^{\left(\frac{2kT_{EFF}}{kT}\right)}\right].$$
(3.18)

When we calculate the second integral, the result is

$$I_{2} = \left(\frac{2kT_{EFF}}{q}\right) \left(\frac{C_{OX}}{\sqrt{2\varepsilon_{IGZO}N_{EFF}kT_{EFF}}}\right)^{\left(\frac{2kT_{EFF}}{kT}-1\right)} \frac{1}{\left(\frac{2kT_{EFF}}{kT}-1\right)} \\ \times \left[\left(V_{GS}-V_{FB}-\Phi_{SS}\right)^{\left(\frac{2kT_{EFF}}{kT}-1\right)} - \left(V_{GS}-V_{FB}-\Phi_{SD}\right)^{\left(\frac{2kT_{EFF}}{kT}-1\right)}\right]$$
(3.19)

If we insert  $I_1$  and  $I_2$  into (3.16), the final expression is obtained as

$$\begin{split} I_{DS} &= \frac{W}{L} \frac{q\mu_{BAND}N_{c}}{B^{*}\sqrt{\frac{2N_{EFF}kT_{EFF}}{\varepsilon_{IGZO}}}} \left(\frac{V_{GS} - V_{T}}{V_{AA}}\right)^{\gamma} \left(\frac{C_{OX}}{\sqrt{2\varepsilon_{IGZO}N_{EFF}kT_{EFF}}}\right)^{\left(\frac{2kT_{EFF}}{kT}-1\right)} \\ &\times \left\{ \left(\frac{kT}{2kT_{EFF}}\right) \left[ (V_{GS} - V_{FB} - \Phi_{SS})^{\left(\frac{2kT_{EFF}}{kT}\right)} - (V_{GS} - V_{FB} - \Phi_{SD})^{\left(\frac{2kT_{EFF}}{kT}\right)} \right] \\ &+ \left(\frac{2kT_{EFF}}{q}\right) \frac{1}{\left(\frac{2kT_{EFF}}{kT} - 1\right)} \\ &\times \left[ (V_{GS} - V_{FB} - \Phi_{SS})^{\left(\frac{2kT_{EFF}}{kT} - 1\right)} - (V_{GS} - V_{FB} - \Phi_{SD})^{\left(\frac{2kT_{EFF}}{kT} - 1\right)} \right] \right\} \end{split}$$
(3.20)

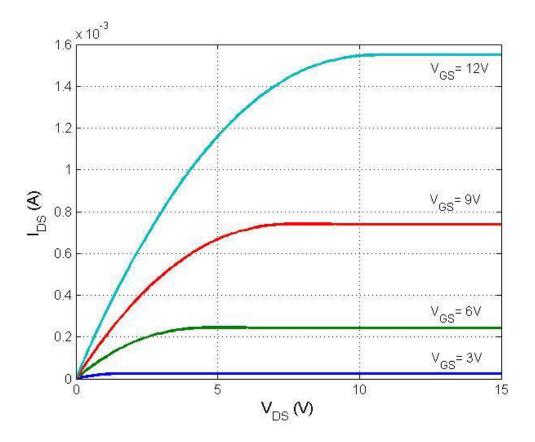
This equation is the main equation of our model that gives the relation between  $I_{DS}$ ,  $V_{GS}$  and  $V_{DS}$ . We compared the calculations from this equation firstly with the model that is more similar to our model then the model in [18] in the sense that it also uses

the  $n_{EFF}$  for the number of carriers taking part in the channel conduction process for an a-IGZO TFT [17], however, differing from our model in modeling the mobility. In [17] Park et al. used exactly the same formulations as we analyzed in Chapter 2 in detail, however, they used only one  $N_{EFF}$  and  $kT_{EFF}$  for an effective carrier concentration  $n_{EFF}$ . Therefore, to make the comparison we took the same parameter values as used by Park et al. [17], shown in Table 3.1.

Parameter	Value	Parameter	Value	
W / L [μm]	200/50	$N_{\rm EFF}~[{ m m}^{-3}]$	$3.12 \times 10^{18}$	
$\mu_{\rm BAND}  [{\rm cm}^2 /  {\rm V.s}]$	17.4	$kT_{\rm EFF}$ [eV]	0.028	
$N_{c}  [{\rm cm}^{-3}]$	$3 \times 10^{18}$	$C_{\rm ox} [\rm nF/cm^2]$	42.37	
$V_{_{ m FB}}$ [V]	0.2	$\phi_{_{ m F0}}$ [V]	0.5	
<i>V</i> <sub>AA</sub> [V]	1	<i>V</i> <sub>T</sub> [V]	0.6	

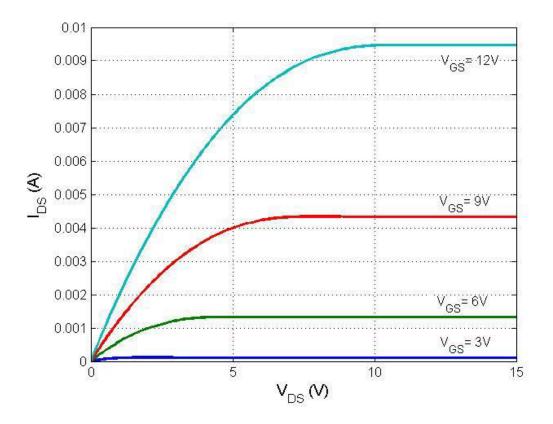
 Table 3 Geometrical and extracted model parameters [17]

Using these values, we calculated the output characteristics of the TFT and generated Fig. 16 with Eq. 3.20, by using  $\gamma = 0.06$  as used by [19].



**Figure 16**  $I_{DS}$  curves by using Eq. 2.47

To compare the results from our formula with the model by [17], we calculated the  $I_{DS}$  values, by changing  $V_{GS}$  and  $V_{DS}$  using Eq. 2.47 (Park et al.'s formulation) with the same parameter values given in Table 3.1. We obtained Fig. 17.



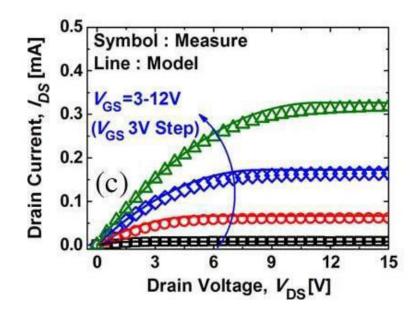
**Figure 17**  $I_{DS}$  curves by using Eq. 3.20

When we compare Fig. 16 and Fig. 17, we can see that although the general behaviors of the curves are the same, the numerical values are at least an order of different. This is due to the reason that we took a constant  $\gamma$  value. However, as we explained above, our  $\gamma$  is a function of  $V_{GS}$  and  $V_{DS}$ . We need to determine the best function for  $\gamma$  to successfully fit the TFT characteristics.

# **3.2. Fitting Theoretical Parts with Experimental Values**

In finding a proper function to represent  $\gamma$ , we used the measured data of the TFTs available in the literature. We found the measured data in the figures from the manuscripts in the literature but we needed to accurately transcribe the correct numerical values from the figures. We used some image processing to obtain the

numerical data values from the figures. Fig 3.4 shows the output characteristics of an a-IGZO TFT from measured data given by Minkyung Bae et al. [18]. We obtained the numerical values from Fig. 18 by applying the image processing methods to form our measured data set.



**Figure 18**  $I_{DS}$  -  $V_{DS}$  graphics (output characteristics) of an a-IGZO TFT from [18] used to obtain the measured data values for our model verification

In the graphic given in Fig. 18, symbols show the measured data values. It is not easy to get numerical values by eye inspection, and moreover, we need to have accurate data values to be compared properly with our model values. So, we used image processing methods and got numerical values. The derived numerical values are given in Table 3.2.

V <sub>GS</sub>	=12V	$V_{ m GS}$	<sub>s</sub> =9V	$V_{ m G}$	s=6V	$V_{ m GS}$	=3V
$V_{\rm DS}$ (v)	$I_{\rm DS}~({\rm mA})$	$V_{\rm DS}$ (v)	$I_{\rm DS}~({\rm mA})$	$V_{\rm DS}$ (v)	$I_{\rm DS}~({\rm mA})$	$V_{\rm DS}$ (v)	$I_{\rm DS}~({\rm mA})$
0,465	0,0103	0,468	0,00948	0,624	0,00558	0,938	0,00727
0,95	0,0382	0,796	0,0231	0,874	0,0123	1,38	0,0117
1,43	0,0658	1,08	0,0348	1,14	0,0205	1,91	0,015
1,91	0,093	1,45	0,0489	1,46	0,0279	2,39	0,0161
2,39	0,116	1,91	0,0644	1,9	0,0359	2,86	0,0171
2,89	0,14	2,42	0,0806	2,38	0,0424	3,36	0,0181
3,36	0,161	2,9	0,0943	2,87	0,0484	3,85	0,0182
3,85	0,18	3,39	0,106	3,36	0,0538	4,33	0,0182
4,34	0,199	3,86	0,117	3,85	0,058	4,82	0,0182
4,83	0,215	4,34	0,127	4,33	0,0618	5,3	0,0182
5,31	0,23	4,83	0,135	4,82	0,0638	5,76	0,0182
5,79	0,244	5,8	0,148	5,29	0,0662	6,26	0,0185
6,27	0,255	6,29	0,153	5,79	0,0673	6,76	0,0185
6,75	0,267	6,77	0,157	6,27	0,0686	7,25	0,0185
7,24	0,277	7,25	0,161	6,75	0,0688	7,71	0,0185
7,72	0,285	7,74	0,163	7,24	0,0688	8,19	0,0185
8,21	0,292	8,22	0,165	7,71	0,0692	8,67	0,0188
8,7	0,298	8,7	0,167	8,21	0,0694	9,16	0,0188
9,18	0,303	9,2	0,168	8,69	0,0696	9,65	0,0188
9,66	0,306	10,6	0,17	9,16	0,0701	10,1	0,0188
10,2	0,311	11,1	0,171	9,64	0,0701	10,6	0,0188
10,6	0,313	12,1	0,172	10,1	0,0701	11,1	0,0188
11,1	0,315	12,6	0,172	10,6	0,0701	11,6	0,0188
11,6	0,317	13,1	0,172	11,1	0,0703	12,1	0,0188
12,1	0,318	13,6	0,172	11,6	0,0703	12,6	0,0191
12,6	0,32	14	0,172	12,1	0,0703	13,1	0,0191
13,1	0,32	14,5	0,172	12,6	0,0703	13,5	0,0191
13,5	0,322	14,9	0,172	13,1	0,0705	14	0,0191
14	0,322			13,5	0,0705	14,5	0,0191
14,5	0,322			14	0,0705	14,9	0,0191
14,9	0,322			14,5	0,0705		
				14,9	0,0705		

**Table 4** The measured data values obtained from Fig. 18 by an image processing

 method (all data values are for 4 significant digits)

We then tried to best fit our model calculations to the measured data by changing  $\gamma$ . To do this, first we analyzed  $\gamma$  dependency of our model. In order to determine the  $\gamma$  dependency on  $V_{GS}$  and  $V_{DS}$  we decided to fit the calculated  $I_{DS}$  values to the experimental values given in Fig. 18. In doing this, we firstly used the image processing method and formed a data table (Table 3.2) from measured  $I_{DS}$  values for the applied  $V_{DS}$  and  $V_{GS}$  values given in Fig. 18. Secondly, we calculated  $I_{DS}$  values by using Eq. 3.20. Then, we compared it with measured value for the same  $V_{GS}$  and  $V_{DS}$ . If it is different, we changed  $\gamma$  value. We calculated  $I_{DS}$  with new  $\gamma$  and compared it with measured data again. We repeated this process until we caught the measured value within a tolerance we decided. Thus, we fitted all the measured data points as given in Table 3.2 with the values obtained by using Eq. 3.20 as shown in Fig. 20, by calculating a  $\gamma$  value that provides the best fit for each data point.

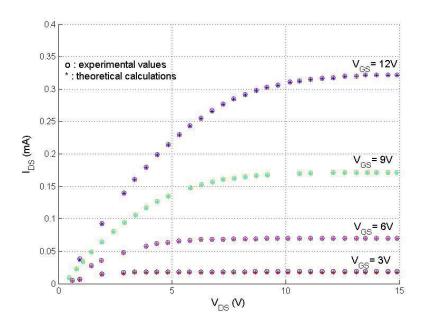


Figure 19 Experimental data and our model fitting by changing  $\gamma$ 

In Fig 19, 'o' symbols show the measured data from Table 3.2. '\*' symbols denote the  $I_{DS}$  values that are calculated by best fitting  $I_{DS}$  values using Eq. 3.20 via changing  $\gamma$ . These  $\gamma$  values that make  $I_{DS}$  values best fit to the experimental data are sketched in Fig. 20. We noticed from this figure that the biggest changes for  $\gamma$  occurred in the linear region of the  $I_{DS}$  versus  $V_{DS}$  curves in Fig. 18. For every measured data point in the linear region to fit, we needed a different  $\gamma$  value, as seen in Fig. 20.

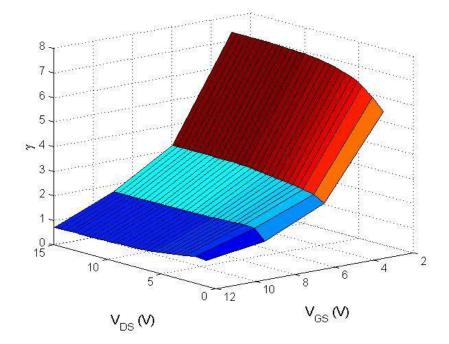


Figure 20 Drain and Gate voltage and best fitting  $\gamma$  relation

We developed a 2-D fitting equation for  $\gamma$  by using the Matlab's curve fitting tool for surfaces [53].

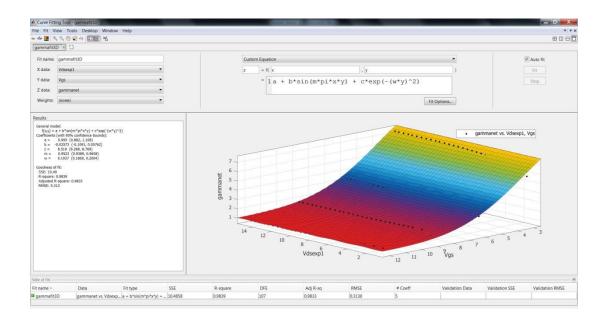


Figure 21 Snapshot of the custom equation surface fitting process screen

In fitting the 2-D equation to our 2-D surface in Fig. 20, we determined that a "custom equation" method in Matlab tools is the best way among others available. The fitted 2-D surface is given in Fig. 21. By using that method, we were able to find a function that expresses  $\gamma$  successfully as a function of  $V_{GS}$  and  $V_{DS}$ , given in the window of Fig. 21 as

$$\gamma = 1 - 0.025 \sin\left(\frac{\pi V_{\rm GS} V_{\rm DS}}{V_{\rm AA}^2}\right) + 8.5 e^{-\left(\frac{0.19 V_{\rm GS}}{V_{\rm AA}}\right)^2}$$
(3.21)

When we combined Eq. 3.20 and Eq. 3.21, we found the final form of the  $I_{DS}$  equation of our model.

$$I_{\rm DS} = \frac{W}{L} \frac{q\mu_{\rm BAND} \, N_{\rm c}}{B^* \sqrt{\frac{2N_{\rm EFF} kT_{\rm EFF}}{\varepsilon_{\rm IGZO}}}} (V_{\rm GS} - V_{\rm T})^{(1-0.025 \, \rm sin \left(\frac{\pi V_{\rm GS} V_{\rm DS}}{V_{\rm AA}^2}\right) + 8.5e^{-\left(\frac{0.19V_{\rm GS}}{V_{\rm AA}}\right)^2})} \times \left\{ \left(\frac{C_{\rm OX}}{\sqrt{2\varepsilon_{\rm IGZO}} N_{\rm EFF} kT_{\rm EFF}}\right)^{\left(\frac{2kT_{\rm EFF}}{kT} - 1\right)} \times \left\{ \left(\frac{kT}{2kT_{\rm EFF}}\right) \left[ (V_{\rm GS} - V_{\rm FB} - \Phi_{\rm SS})^{\left(\frac{2kT_{\rm EFF}}{kT}\right)} - (V_{\rm GS} - V_{\rm FB} - \Phi_{\rm SD})^{\left(\frac{2kT_{\rm EFF}}{kT}\right)} \right] \right\}$$
(3.22)
$$+ \left(\frac{2kT_{\rm EFF}}{q}\right) \frac{1}{\left(\frac{2kT_{\rm EFF}}{kT} - 1\right)} \times \left[ (V_{\rm GS} - V_{\rm FB} - \Phi_{\rm SS})^{\left(\frac{2kT_{\rm EFF}}{kT} - 1\right)} - (V_{\rm GS} - V_{\rm FB} - \Phi_{\rm SD})^{\left(\frac{2kT_{\rm EFF}}{kT}\right)} \right] \right\}$$

This equation best fits our experimental values as shown in Fig. 19. The Matlab code to derive these current values is given in Appendix. Hence we derived a new formula for an a-IGZO TFT to estimate its static electrical characteristics. Our formula comprises an empirical model for the channel mobility ( $\mu_{CH}$ ) and based on the channel carrier concentration ( $n_{EFF}$ ) dependence on the electrostatic potential, which is a function of  $V_{CS}$  and  $V_{DS}$  voltages. We used the approach taken by L. Colalongo [42], M. Bae et al. [18] and J. H. Park et al. [17], to solve the nonlinear equation to determine the electrostatic potential in the channel as a function of applied voltages. This model calculates the  $I_{DS}$  versus  $V_{DS}$  characteristics with one semi-analytic formula and does not need a different formula for each region. Our model differs from L. Colalongo and M. Bae et al. models in determining the mobility and the  $n_{EFF}$  in the channel. In their calculations they took the band tail states and midgap states that contribute to the carrier concentration as localized carriers ( $n_{LOC}$ ) together

with the free carriers  $(n_{FREE})$  in the conduction band of a-IGZO channel. However, in our model, like J. H. Park et al., we used only one type of carriers  $(n_{EFF})$  with one characteristic slope  $(kT_{EFF})$  to express the effective carrier concentrations in the channel together with the free carriers in the conduction band.

M. Bae et al. and J. H. Park et al. introduced a mobility model that is dependent of the free and localized carriers, however in our model we introduced a different empirical model for the channel mobility to enable us to 100% fit the model calculations of  $I_{DS}$  as a function of  $V_{GS}$  and  $V_{DS}$  if we have the experimental values available.

Our model is very powerful and flexible in the sense that if we have the experimental values of the  $I_{DS}$  as a function of  $V_{GS}$  and  $V_{DS}$  we can adjust our mobility formula to fit the  $I_{DS}$  values completely at every point of characteristics to the experimental values. Therefore, once one has the experimental characteristics of an a-IGZO TFT, using those data points in the measured data set we can adjust the best fitting mobility formula in our model for the specific TFT that we can 100% fit (our error range is 0.1%) to the DC  $I_{DS}$  values as a function of  $V_{GS}$  and  $V_{DS}$ . By using this formula afterwards, one can determine the circuit characteristics that are configured using the same type of TFTs as required. Yong et al. [54] reported that their static current model for an a-IGZO TFT could fit the voltage transfer characteristics (VTC) of an IGZO inverter, where two a-IGZO TFTs are used, within error ratio of 6%. We expect that a lower ratio of error than theirs could most probably be obtained in such an application if our model were used in the simulation. However, it should be

reminded that our model does not yet include the effect of the back potential between the channel and the substrate, as their model does.

# **CHAPTER 4**

### 4. CONCLUSION

# 4.1. Conclusion

In this thesis we first analyzed the thin film transistors (TFTs) that are generally used as switching elements on large panel displays to switch the pixels on Liquid Crystal Displays (LCDs), light emitting diode (LED) displays, flexible displays, and on electronic paper and so on. These transistors have mostly an amorphous material for the channel. Amorphous hydrogenated silicon (a-Si:H) TFT was the first extensively used switching device because a-Si:H material is easy to deposit on a glass substrate at relatively low temperatures by using low-cost vacuum deposition techniques. However, it was found that a-Si:H is absorbing visible light photons and loses its stability in time. Recently a new amorphous metal oxide material of indium gallium zinc oxide (a-IGZO) has replaced the a-Si:H as the channel material in the TFTs due to its higher channel mobility, and better optical and electrical reliability.

We analyzed the material properties of a-Si:H and a-IGZO materials and the properties of TFTs made by using them. We discussed the requirements for a good TFT to drive the pixels on a flat panel display of different technologies, and the channel material properties that meet these requirements mostly. We found that a-IGZO is the best material available for the TFTs today.

We then discussed the analytic models to describe the static electric characteristics of a TFT, which is a very similar device to a metal-oxide-semiconductor field effect transistor (MOSFET). We analyzed first the gradual channel approximation analytic model for a TFT. We pointed out that this model requires different expressions for each region of the transistor in the output characteristics, namely the sub-threshold, above-threshold, and saturation regions.

Searching for a better analytic model for an a-IGZO TFT, we analyzed a semianalytic model for the TFT that estimates its behavior mainly by expressing the channel conductance as a function of the surface electrostatic potential, which is in turn dependent on the gate to source and drain to source voltages,  $V_{GS}$  and  $V_{DS}$ . We then indicated the importance for expressing the mobility of the channel in this model by relating it to the localized and free carriers in the amorphous materials.

We first reviewed Colalongo's [42] approach to model an a-Si:H TFT, where he used the tail and the deep states within the band gap of the amorphous silicon material to calculate the drain current using the charge-sheet analytical model. He used a constant mobility in deriving the drain current. We observed that his model needs rather high value of channel mobility  $(10cm^2/Vs)$  than measured for a-Si:H to fit the drain current predictions to the experimental data.

We then investigated Tze-Ching Fung's model for an a-IGZO TFT [21]. Fung used the gradual channel approximation to calculate the drain current. He used a channel potential dependent mobility formula in his model. In his model he used different formulas of drain current for different regions of the TFT like gradual channel approximation, however, he introduced an empirical mobility expression in addition to better fit the predictions of the model to the experimental data.

Another analytic model was introduced by Minkyung Bae et al. [18] to represent the a-IGZO TFT static electrical behavior with one formula being valid in all the regions of operation. They used Colalongo's model of the drain current dependency of the localized states (tail states, and deep states), but they also add free carriers in their model because unlike the case for a-Si:H, in an a-IGZO free carrier concentration is also important. In addition to this, like Colalongo they used a mobility expression which is dependent on the carrier concentrations, that is, the channel potential. Their model incorporated tail states and deep states to represent the behaviors in subthreshold and above-threshold regions of the TFT. We analyzed this model by deriving all the equations in the model in detail and developed the Matlab codes to solve the nonlinear equation to calculate the surface potential for the applied voltages. We verified that with our flowchart and codes we obtained the same values of drain currents for the applied voltages with theirs. We showed that their model is good in predicting the a-IGZO TFT behavior provided that the density of states in deep states  $(N_{\rm EFF2})$  and the characteristic slope  $(kT_{\rm EFF2})$  of the tail states are estimated very accurately, which seems impossible.

Jun Hyun Park et al. [17] introduced a new model following Minkyung Bae et al.'s way of development. The only difference they have from [18] is that instead of using tail and deep states and their specific activation energies for representing the channel conductivity, they introduced an effective carrier density  $(n_{EFF})$  with one effective density of states  $(N_{EFF})$  and one effective activation energy  $(kT_{EFF})$ . However, for mobility they did the same thing as [18] and used the formula, namely they took

$$\mu_{EFF} = \mu_{BAND} \left[ \frac{n_{FREE}}{\left( n_{FREE} + n_{LOC} \right)} \right]$$
(4.1)

Although this model resolves somewhat the sensitivity of estimating the certain parameters, it is yet not predicting the behavior of the a-IGZO TFT's as accurately as [18] in all the operating regions of the TFT.

Then we developed a new analytic model for an a-IGZO TFT using the same approach as Jun Hyun Park, by taking only one effective carrier density  $(n_{EFF})$  with one effective density of states  $(N_{EFF})$ , and one activation energy  $(kT_{EFF})$  to calculate the drain current of the a-IGZO TFTs. However, unlike associating the channel mobility with the free and localized carrier densities as Minkyung Bae et al. and Jun Hyun Park et al. did, and associating the channel mobility with channel potential indirectly, we used an empirical formula that associates channel mobility with the channel potential directly, and derived the formulation to express  $I_{DS}$  with  $V_{GS}$  and  $V_{DS}$ . In this formulation, we improved the mobility expression by fitting the model's predictions with the measured data in an a-IGZO TFT. We used the best fitting mobility formula in our last equation. Our model strength comes from the fact that it is capable of fitting the experimental data nearly 100% by arranging our mobility formula. This strength of our model could be employed in the applications where the model that predicts the behavior of the TFT is used in developing new circuits involving the many TFTs of the same or similar physical and geometric properties and simulating the overall behavior of the circuit prior to manufacturing. In this direction, an example case that Yong et al. [54] reported that they developed a numerical model for the static electrical behavior for an a-IGZO TFT and they used their model to fit the voltage transfer characteristics (VTC) of an a-IGZO inverter, where two a-IGZO TFTs are used, within error ratio of 6%. We expect that a lower ratio of error than theirs could most probably be obtained in such an application if our model were used in the simulation. However, it should be reminded that our model does not yet include the effect of the back potential between the channel and the substrate, as their model does.

#### 4.2. Future Work Plan

Our last equation could be improved to include the back potential to make it suitable to use in the circuit simulations.

Our model predictions should be tested on simulating the device behavior, such as an inverter for instance, and the results should be benchmarked with the literature.

In developing and testing to improve our model, main problem we had was to find enough sample size of measured data from the TFTs of known physical and geometric parameters. We are planning to manufacture some a-IGZO TFTs in cooperation with the research groups elsewhere and make the required measurements to generate the required data set to improve our analytical model further.

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## **APPENDICES A**

#### SAMPLE MATLAB CODES

```
%This code is written to sketch the graphs that
%are given in Chapter 2.
clc;
clear all;
close all;
%%% constants
W=225*10^(-6); %m
L=30*10^(-6); %m
Lov=5*10^(-6); %m
Tox=100*10^(-9); %m
%Tigzo=50*10^(-7); %cm
Mband=19.7*10^(-4); %m^2/V.s
Nc=4.8*10^(12); %m^(-3)
Vfb=0.3; %V
Neff1=8.7*10^(11); %m^(-3)
kTeff1=0.045*(1.6*10^(-19)); %eV
Neff2=5.15*10^(11); %m^(-3)
kTeff2=0.0263*(1.6*10^(-19)); %eV
Cox=42.37*10^(-5); %F/m^2
phiF0=0.3; %V
q = 1.6*10^(-19);%e
Eo=8.85*10^(-12);%F/m
eIGZO=11.5*Eo; %F/m
kT=0.026*q;
୫୫୫୫୫୫୫୫୫୫<u></u>
syms phis;
Vgs=[3 6 9 12]; %V
Vds=0:0.5:15; %V
C=((2/kT)-(3/(2*kTeff1)));
B=((1/kT) - (1/(2*kTeff1)));
A=(Nc*sqrt(eIGZO))/(B*sqrt(2*Neff1*kTeff1));
for i=1:length(Vgs)
    for j=1:length(Vds)
        phiss(i)=solve(phis-phiF0-2*(kTeff1/q)...
            *log(Cox*(Vgs(i)-Vfb-phis)/sqrt(2*eIGZO*Neff1*kTeff1)));
```

```
phisd(i,j)=solve(phis-phiF0-Vds(j)-2*(kTeff1/q)...
            *log(Cox*(Vgs(i) - Vfb -phis)/sqrt(2*eIGZO*Neff1*kTeff1)));
        Idssub(i,j) = (W/L) * Mband*((A*Nc)/(2*Neff1*kTeff1*C))...
            *((Cox/sqrt(2*eIGZO*Neff1*kTeff1))^(2*C*kTeff1))*...
            (((2*C*kTeff1+1)^(-1))*(((Vgs(i)-Vfb-phiss(i))...
            ^(2*C*kTeff1+1))-((Vgs(i)-Vfb-phisd(i,j))...
            ^(2*C*kTeff1+1)))-((q*C)^(-1))*(((Vgs(i)-Vfb-phiss(i))...
            ^(2*C*kTeff1))-((Vgs(i)-Vfb-phisd(i,j))^(2*C*kTeff1))));
    end
end
figure(1)
plot(Vds,Idssub,'-o','LineWidth',1.5);
%set(gca, 'FontSize',12);
xlabel('V D S (V)', 'FontSize',12);
ylabel('I D Ssub (A)', 'FontSize', 12);
grid on;
C2=((2/kT) - (3/(2*kTeff2)));
B2=((1/kT) - (1/(2*kTeff2)));
A2=(Nc*sqrt(eIGZO))/(B2*sqrt(2*Neff2*kTeff2));
X2=(W/L) *Mband* ((A2*Nc)/(2*Neff2*kTeff2*C2));
Y2=((Cox/sqrt(2*eIGZO*Neff2*kTeff2))^(2*C2*kTeff2));
for i=1:length(Vgs)
    for j=1:length(Vds)
        phiss2(i)=solve(phis-phiF0-2*(kTeff2/q)...
            *log(Cox*(Vgs(i)-Vfb-phis)/sqrt(2*eIGZO*Neff2*kTeff2)));
        phisd2(i,j)=solve(phis-phiF0-Vds(j)-2*(kTeff2/q)...
            *log(Cox*(Vgs(i) - Vfb -phis)/sqrt(2*eIGZO*Neff2*kTeff2)));
        Idsabove(i,j) = (W/L)*Mband*((A2*Nc)/(2*Neff2*kTeff2*C2))...
            *((Cox/sqrt(2*eIGZO*Neff2*kTeff2))^(2*C2*kTeff2))*...
            (((2*C2*kTeff2+1)^(-1))*(((Vgs(i)-Vfb-phiss2(i))...
            ^(2*C2*kTeff2+1))-((Vgs(i)-Vfb-phisd2(i,j))...
            ^(2*C2*kTeff2+1)))-((q*C2)^(-1))...
            *(((Vgs(i)-Vfb-phiss2(i))^(2*C2*kTeff2))...
            -((Vgs(i)-Vfb-phisd2(i,j))^(2*C2*kTeff2))));
    end
end
figure(2)
plot(Vds,Idsabove,'-o','LineWidth',1.5);
%set(gca, 'FontSize', 12);
xlabel('V_D_S (V)','FontSize',12);
ylabel('I D Sabove (A)', 'FontSize', 12);
grid on;
for i=1:length(Vgs)
    for j=1:length(Vds)
        Ids(i,j) = (1/Idssub(i,j)) + (1/Idsabove(i,j));
        Ids(i,j)=1/Ids(i,j);
```

```
end
end
figure(3)
plot(Vds,Ids,'-o','LineWidth',1.5);
%set(gca,'FontSize',12);
xlabel('V_D_S (V)','FontSize',12);
ylabel('I_D_S total (A)','FontSize',12);
grid on;
```

```
figure(4)
```

```
surf(Vds,Vgs,double(phisd));
```

xlabel('V\_D\_S'); ylabel('V\_G\_S'); zlabel('phi\_S\_D');

## **APPENDICES B**

## **CURRICULUM VITAE**

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## **EDUCATION**

Degree	Institution	Year of Graduation
M.Sc	Çankaya University Electronic and Communication Engineering	January, 2015
B.Sc	Çankaya University Electronic and Communication Engineering	June, 2011
High School	19 Mayıs High School, Mersin	June, 2006

# FOREIGN LANGUAGE

Advanced English

### **PUBLICATION**

Bozkurt İ., Çil C. Z., Çalbıyık G., "Lazer ile Mesafe Ölçüm Sistemi", Çankaya Üniversitesi 5. Mühendislik ve Teknoloji Sempozyumu, Ankara, Türkiye, Apr. 2012. Çankaya Üniversitesi 5. Mühendislik ve Teknoloji Sempozyumu Kitapçığı, 55-60, 26-27 Nisan 2012.

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