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An Empirical Modeling of Gate Voltage-Dependent Behaviors of Amorphous Oxide Semiconductor Thin-Film Transistors including Consideration of Contact Resistance and Disorder Effects at Room Temperature

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Copyright: © 2021 by the author. Licensee MDPI, Basel, Switzerland. This article is an open access article distributed under the terms and conditions of the Creative Commons Attribution (CC BY) license (https:// creativecommons.org/licenses/by/ 4.0/). Department of Electronics Engineering, Pusan National University, Pusan 46241, Korea; sungsiklee@pusan.ac.kr; Tel.: +82-51-510-3123

Abstract: In this paper, we present an empirical modeling procedure to capture gate bias dependency of amorphous oxide semiconductor (AOS) thin-film transistors (TFTs) while considering contact resistance and disorder effects at room temperature. From the measured transfer characteristics of a pair of TFTs where the channel layer is an amorphous In-Ga-Zn-O (IGZO) AOS, the gate voltage-dependent contact resistance is retrieved with a respective expression derived from the current–voltage relation, which follows a power law as a function of a gate voltage. This additionally allows the accurate extraction of intrinsic channel conductance, in which a disorder effect in the IGZO channel layer is embedded. From the intrinsic channel conductance, the characteristic energy of the band tail states, which represents the degree of channel disorder, can be deduced using the proposed modeling. Finally, the obtained results are also useful for development of an accurate compact TFT model, for which a gate bias-dependent contact resistance and disorder effects are essential.

Keywords: amorphous oxide semiconductor; thin-film transistor; degree of disorder; bias-dependent contact resistance; compact transistor model; empirical modeling; transfer characteristics

1. Introduction

Amorphous semiconducting materials such as amorphous Silicon and amorphous oxide semiconductors (AOSs) have been widely used as the channel layer for thin-film transistors (TFTs) [1–3]. In particular, it is believed that the AOS TFT has become one of the most promising candidates for futuristic electronics due to its high transparency and low temperature processability [3–5]. With a low temperature process, the AOS film is more likely to be in amorphous phase; thus, the inevitable presence of localized traps (e.g., deep and tail states) is associated with structural disorder in the amorphous phase [2,6]. This additionally alludes to a poor quality of metal contacts at the source and drain [1,2,7]. It is known that these non-ideal properties have a significant influence on the electrical performances of the AOS TFTs [2,3,8]; for example, the field-effect mobility of AOS TFTs is inversely proportional to the density of the localized tail states, while poor contact can lead to a higher contact resistance, and thus a lower mobility [9]. In addition, the contact resistance is typically extracted with a transmission line method (TLM), yielding a constant value without a gate-bias dependency [10]. Therefore, it is necessary to analyze and model the correlation between the transistor characteristics and those parasitic properties which include gate bias-dependent contact resistance.

In this paper, to capture gate bias-dependent contact resistance (R_C) and Intrinsic channel conductance (G_{int}) in the AOS TFT, an empirical method is proposed based on transfer characteristics measured for the two different AOS TFTs, as the parasitic effects are reflected in the current-voltage characteristics of the transistors. For this, we examine three fabricated In-Ga-Zn-O (IGZO) TFTs. From them a pair of TFTs, chosen from three examined



TFTs (i.e., three pairs are available as three combinations made from three examined TFTs), is used to retrieve the gate voltage-dependent R_C and G_{int} while applying an analytical expression derived from the current–voltage relations of TFTs. Based on the retrieved trend of the R_C vs. V_{GS} , which is modelled with a power law, the G_{int} can also be accurately extracted and modelled, yielding the characteristic energy of band tail states as a measure of disorders (assuming the dominance of the trap-limited conduction). From these results, it is found that the gate-bias dependencies of both R_C and G_{int} are well explained with a power-law function. Finally, it is believed that the presented results could be useful for an accurate compact TFT model, where the gate-bias dependencies of the contact resistance and disorder effects are crucial.

2. Materials and Methods

In this work, we examined AOS TFTs where amorphous IGZO is incorporated as the AOS material, with disorder in the amorphous phase (see Figure 1). Following a typical bottom-gate TFT process as reported in [11,12] for the deposition of the IGZO layer on a glass substrate, an IGZO ceramic target was employed for RF sputtering with Ar plasma. During this sputtering process, the oxygen partial pressure was kept at a low level (e.g., 5%), followed by backend processes such as thermal annealing, patterning, metallization with Mo for the electrodes (i.e., source, drain, and gate), and passivation. Using this process, we prepared three IGZO TFTs with three different channel lengths (L_1 , L_2 , L_3). Here, the channel widths for those TFTs remain the same ($W = W_1 = W_2 = W_3$).



Figure 1. Schematic cross-sectional view of the examined IGZO TFT describing the contact resistance (R_C) and disorder (traps) within the IGZO channel layer.

Using the three fabricated TFTs (TFT-1, TFT-2, TFT-3), the transfer characteristics, i.e., the drain current (I_{DS}) vs. gate voltage (V_{GS}), were measured at room temperature (300 K), which formed the basis for the proposed empirical modeling. The channel geometrical details of the examined TFTs are summarized for the three possible combinations out of three TFTs in Table 1. Along with the examined TFTs, empirical modeling was performed with relevant mathematical formulations, as explained in the following sections.

Table 1. Summary of channel geometrical details of three examined IGZO TFTs.

Examined IGZO TFTs	Channel Length (L)	Channel Width (W)
TFT-1	$L_1 = 10 \ \mu m$	$W = 50 \ \mu m \ (common) (W = W_1 = W_2 = W_3)$
TFT-2	$L_2 = 20 \ \mu m$	
TFT-3	$L_3 = 30 \ \mu m$	-

3. Results

3.1. Mathematical Formulations for AOS TFTs

Considering the channel disorder and contact resistance, the transfer characteristics in the linear regime, i.e., I_{DS} vs. V_{GS} , of n-channel AOS TFTs can be represented as follows:

$$I_{DS} = K_n \frac{W}{L} (V_{GS} - V_T)^{1 + \alpha_{cm}} (V_{DS} - 2R_C I_{DS}).$$
(1)

where K_n is a pre-factor, V_T is the threshold voltage, α_{cm} is an exponent related to the conduction mechanisms (e.g., trap-limited conduction or percolation conduction), V_{DS} is the drain voltage, and R_C is the contact resistance. Note that in Equation (1), the physical meaning of α_{cm} can be changed when the dominant conduction mechanism is determined from the trap-limited conduction ($\alpha_{cm} \equiv \alpha_t$, a trap-related exponent) or the percolation conduction ($\alpha_{cm} \equiv \alpha_p$, a percolation-related exponent) [13–15]. For example, depending on the position of the Fermi level (E_F), the dominant conduction mechanism is determined [13,15]. In the present study, we presume that the E_F is well below the E_C for the given range of the V_{GS} [15]. Thus, assuming that the trap-limited conduction is dominant, α_{cm} in Equation (1) can now be related to the traps, thus $\alpha_{cm} = \alpha_t$. Based on this assumption, K_n and α_t are further defined, respectively, as follows:

$$K_n = \mu_n \frac{C_{ox}^{1+\alpha_t}}{Q_{ref}^{\alpha_t}},\tag{2}$$

$$x_t = \frac{kT_t}{kT} \text{ or } \frac{T_t}{T},\tag{3}$$

where μ_n is the electron mobility, C_{ox} is the gate insulator capacitance per area, Q_{ref} is a reference charge-density per area, kT_t is the characteristic energy of tail states, and kT is the thermal energy (Here, k is Boltzmann's constant). Note that for a given temperature (T), Equation (3) can be rewritten as $kT_t = \alpha_t kT$ or $T_t = \alpha_t T$. Here, a larger value of kT_t or T_t means a higher degree of disorder at a fixed T [15,16]. As can be seen in Equations (1)–(3), the formula for the drain current is strongly related to both the contact resistance and the disorder through α_t as a function of kT_t at a given temperature (T). In other words, Equation (1), where α_t is included, is more general compared to the textbook, where an ideal case with $R_C = \alpha_t = 0$ is only covered for a perfect crystalline semiconductor; hence, the universality of Equation (1).

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With Equation (1), the measured transfer characteristics in the linear regime (i.e., $V_{GS}-V_T >> V_{DS} = 0.1$ V) can be explained rather than using the ideally linear equation for $R_C = \alpha_t = 0$. Indeed, as can be seen in Figure 2, the curvature in the above threshold region (at $V_{GS} > V_T$) looks somewhat like a root function; thus $R_C \neq 0$ and $\alpha_t \neq 0$. Note that V_T is extracted with the second derivative method rather than a linear extrapolation [16]. This suggests the presence of contact resistance and disorder (i.e., tail states); thus, the R_C and α_t should be extracted in order to capture that behavior, as seen in Figure 2. In the following two sub-sections, the retrieval procedure for R_C and α_t is shown in detail.



Figure 2. Measured transfer characteristics for $V_{DS} = 0.1$ V of the fabricated IGZO TFT with three different *L* (see Table 1).

3.2. Contact Resistance and Transfer Characteristics of two TFTs

To extract the contact resistance (R_C), we need to derive the respective formula based on Equation (1). For the first combination with TFT-1 and TFT-2, for example, referring to Table 1, the current-voltage relations are as follows:

$$I_{DS1} = K_n \frac{W_1}{L_1} (V_{GS} - V_T)^{1+\alpha_t} (V_{DS} - 2R_C I_{DS1}),$$
(4)

$$I_{DS2} = K_n \frac{W_2}{L_2} (V_{GS} - V_T)^{1+\alpha_t} (V_{DS} - 2R_C I_{DS2}).$$
⁽⁵⁾

Here, it is notable that I_{DS} changes with varying W/L, while other parameters, including R_C , are given or constant regardless of W/L. Note that the terms $(V_{DS} - 2R_C I_{DS1})$ and $(V_{DS} - 2R_C I_{DS2})$ in Equations (4) and (5) are called effective drain voltage, which can also be extracted using R_C . From Equations (4) and (5), we can take their ratio as

$$\frac{I_{DS1}}{I_{DS2}} = \frac{W_1 L_2 (V_{DS} - 2R_C I_{DS1})}{W_2 L_1 (V_{DS} - 2R_C I_{DS2})}.$$
(6)

Equation (6) can be expressed for the R_C as follows:

$$R_C = \frac{(W_1 L_2 I_{DS2} - W_2 L_1 I_{DS1})}{2I_{DS1} I_{DS2} (W_1 L_2 - W_2 L_1)} V_{DS} [\Omega].$$
(7)

Note that Equation (7) is an analytical expression of the R_C for the first pair of TFTs (i.e., TFT-1 and TFT-2). If $W_1 = W_2$, Equation (7) is reduced as

$$R_{C} = \frac{(L_{2}I_{DS2} - L_{1}I_{DS1})}{2I_{DS1}I_{DS2}(L_{2} - L_{1})}V_{DS} \ [\Omega].$$
(8)

Similarly, for the second combination, TFT-2 and TFT-3 for $W_2 = W_3$, Equation (8) is rewritten as

$$R_{C} = \frac{(L_{3}I_{DS3} - L_{2}I_{DS2})}{2I_{DS2}I_{DS3}(L_{3} - L_{2})}V_{DS} \ [\Omega].$$
(9)

As the third pair, Equation (8) for TFT-3 and TFT-1 for $W_3 = W_1$ is reconstructed as

$$R_{C} = \frac{(L_{3}I_{DS3} - L_{1}I_{DS1})}{2I_{DS1}I_{DS3}(L_{3} - L_{1})}V_{DS} [\Omega].$$
 (10)

Note that Equations (8)–(10) can also be normalized with the given channel width (*W*), giving the normalized contact resistance (r_c) as $r_c = R_C W [\Omega \cdot cm]$.

Now, Equations (8)–(10) can be applied to obtain R_C for those three combinations, respectively, along with the measured transfer characteristics seen in Figure 2. Figure 3 shows R_C as a function of V_{GS} - V_T , along with the modeled plots. As can be seen, they show good agreement for all three difference cases. For these modelings, we employed a power-law function:

$$R_{\rm C} = A_{\rm C} (V_{\rm GS} - V_{\rm T})^{-\alpha_c} [\Omega], \qquad (11)$$

where A_C is a pre-factor and α_c is an exponent. As indicated in Figure 3, the extracted values of A_C and α_c for all three cases are approximately $1.8 \times 10^5 \Omega/V^{-\alpha_c}$ and 0.81, respectively.



Figure 3. The retrieved R_C as a function of V_{GS} - V_T for three possible combinations, which are the three cases: (**a**) TFT-1 and TFT-2; (**b**) TFT-2 and TFT-3; (**c**) TFT-3 and TFT-1. Here, the modeled results indicating the values of the model parameters (A_C and α_c) for each case are also shown.

3.3. Intrinsic Channel Conductance and Tail States

First, the ratio between the I_{DS} and $(V_{DS} - 2R_C I_{DS})$ of Equation (1) is normalized with *W/L*. From this, an empirical expression of the intrinsic channel conductance (G_{int}) is given while replacing the term R_C with Equation (11), as follows:

$$G_{int} = \frac{I_{DS}}{\left(V_{DS} - 2A_C (V_{GS} - V_T)^{-\alpha_c} I_{DS}\right) (W/L)}.$$
 (12)

The remaining term in the right-hand-side of Equation (1) is then given independently as

$$G_{int} = K_n (V_{GS} - V_T)^{1+\alpha_t}.$$
(13)

By applying Equation (12) using the results seen in Figures 2 and 3, G_{int} vs. V_{GS} - V_T can be extracted; meanwhile, Equation (13) is applied separately, yielding the values of K_n and α_t , respectively (see Figure 4). Note that in Figure 4 the G_{int} is provided with the R_C effect removed and the channel geometry (W/L) normalized, and is thus common to all examined TFTs.



Figure 4. Plot of G_{int} as a function of V_{GS} - V_T ; the retrieved values of K_n and α_t are indicated. In particular, the α_t is 0.18; thus, $kT_t \approx 4.7$ meV (i.e., $T_t \approx 54$ K).

4. Discussion

As mentioned in the Section 3.1, the measured transfer characteristics of the AOS TFTs can be explained with Equation (1) rather than the ideally-linear equation for $R_C = \alpha_t = 0$. Indeed, as can be seen in Figure 2, it was found that the curvature in the above-threshold region (at $V_{GS} > V_T$) looks slightly like a root function, implying the presence of contact resistance and disorder (i.e., $R_C \neq 0$ and $\alpha_t \neq 0$).

Regarding the effects of contact resistance, R_C was extracted, as seen in Figure 3, by applying Equations (8)–(10) to the transfer characteristics of three pairs of TFTs (see Figure 2). As can be seen in Figure 3, it is obvious that the extracted results have been

well-matched with the proposed model using a power-law. Here, it was found that it decays with increasing V_{GS} . This is because the contact resistance is reduced by narrowing of the Schottky barrier due to a higher gate bias [2,17]. In keeping with this, the results modeled using Equation (11) showed good agreement in all three cases. The retrieved values of the R_C model parameters are listed in the first row of Table 2. As can be seen in Figure 3 and Table 2, the values of the A_C and α_c for all three cases are approximately very similar to each other, with discrepancy negligible at less than 1%. This suggests that the R_C trend for different pairs of TFTs should be the same in principle, and very similar practically as long as the same initial fabrication process was applied and the same ambient conditions were maintained.

Non-Ideal Effects	Model Equations	Model Parameters
Contact Resistance	$R_{C} = \frac{(L_{2}I_{DS2} - L_{1}I_{DS1})}{2I_{DS1}I_{DS2}(L_{2} - L_{1})}V_{DS}$ (for example of TFT-1 and TFT-2)	$A_C \approx 1.8 \times 10^5 \ \Omega/V^{-lpha_c}$ for $W = 50 \ \mu m$
	$R_C = A_C (V_{GS} - V_T)^{-nc} [\Omega]$ $r_c = R_C W [\Omega \cdot \text{cm}]$	$lpha_cpprox 0.81$
Disorder (Traps)	$G_{int} \equiv \frac{I_{DS}}{(V_{DS} - 2R_C I_{DS})W/L}$ $G_{int} = K_n (V_{GS} - V_T)^{1+\alpha_t}.$	$K_n \approx 8.29 \times 10^{-8} \text{ S/V}^{1+\alpha_t}$ $\alpha_t \approx 0.18$ $kT_t \approx 4.7 \text{ meV}$ i.e., $T_t \approx 54 \text{ K}$

Table 2. Summary of model equations and parameters for the same W at T = 300 K.

Using the retrieved model parameters for $R_{C_{i}}$ the intrinsic channel conductance (G_{int}) was then accurately extracted using Equation (12). In other words, we were able to remove the effect of R_C for accurate extraction of G_{int} . As seen in Figure 4, it is clear that good agreement between the extracted and modeled results was achieved. Here, it was found that G_{int} increased without any root-function behavior. However, it exhibited slightly parabolic behavior, with an exponent > 1 using the power law. From this result, it was found that the trend of G_{int} could be modelled with a power law with the exponent 1.18 (i.e., Equation (13) with $1 + \alpha_t$ for $\alpha_t = 0.18$). While the crystalline material-based transistor has the exponent of unity, in our case, the retrieved value of the exponent, i.e., 1.18, was found to be slightly greater than unity. This implies that the channel material is non-crystalline, and thus a disorder. As seen in Table 2, kT_t is 4.7 meV < kT at 300 K, which is consistent with the literature [16,18]. This suggests that the degree of disorder is less than amorphous Silicon, with $kT_t > kT$. As for a physical interpretation, the conduction band of the AOS, e.g., IGZO, is composed of spherical orbitals (i.e., s orbitals) of metal cations; thus, the AOS is insensitive to the bonding angle disorder, whereas amorphous Silicon has strong bonding directivity due to the sp3 orbitals of its conduction band [18–20]. However, the AOS still has a bonding distance error in the amorphous phase, alluding to the existence of localized traps associated with band tail states, which can be estimated using extraction methods based on the current-voltage and capacitance-voltage characteristics [18,21]. In order to minimize these localized traps, process conditions can be optimized in terms of the AOS target compositions for the sputtering process, oxygen partial pressure, and annealing temperature [1,3,22].

Consequently, the presented empirical model for the transfer characteristics, in which the gate-bias dependencies of the contact resistance and disorder effects are considered, could be easily added into a compact TFT model thanks to its simplicity.

5. Conclusions

In this paper, we provided an empirical model to explain gate bias-dependent contact resistance and disorder effects in AOS TFTs at room temperature. As an intermediate outcome, we were also able to obtain the gate-bias dependency of the intrinsic channel conductance where the disorder effects were viewed after removal of the parasitic effect TFTs (i.e., a pair of TFTs), was derived and applied for its retrieval. Interestingly, it was found to follow a simple power law, giving the values of the pre-factor and exponent. In addition, these values were found to be approximately the same for all three pairs examined here. This allowed accurate extraction of the gate bias-dependent intrinsic channel conductance as modelled using a power law. From this analysis, assuming the domination of trap-limited conduction, it was shown that the characteristic energy and temperature of the band tail states could be estimated as a measure of the degree of disorder, which is consistent with the literature. Consequently, the presented results could be useful for the development of an accurate compact TFT model, in which the gate bias-dependent behaviors of the contact resistance and disorder effects are crucial.

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