Determination of energy barrier profiles for high-\(k\) dielectric materials utilizing bias-dependent internal photoemission

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(Received 15 April 2004; accepted 14 September 2004)

We utilize bias-dependent internal photoemission spectroscopy to determine the metal/dielectric/silicon energy barrier profiles for Au/HfO\(_2\)/Si and Au/Al\(_2\)O\(_3\)/Si structures. The results indicate that the applied voltage plays a large role in determining the effective barrier height and we attribute much of the variation in this case to image potential barrier lowering in measurements of single layers. By measuring current at both positive and negative voltages, we are able to measure the band offsets from Si and also to determine the flatband voltage and the barrier asymmetry at 0 V. Our SiO\(_2\) calibration sample yielded a conduction band offset value of 3.0±0.1 eV. Measurements on HfO\(_2\) give a conduction band offset value of 2.7±0.2 eV (at 1.0 V) and Al\(_2\)O\(_3\) gives an offset of 3.3±0.1 (at 1 V). We believe that interfacial SiO\(_2\) layers may dominate the electron transport from silicon for these films. The Au/HfO\(_2\) barrier height was found to be 3.6±0.1 eV while the Au/Al\(_2\)O\(_3\) barrier is 3.5±0.1 eV. © 2004 American Institute of Physics. [DOI: 10.1063/1.1812831]

The study of high-\(k\) dielectrics as a replacement for SiO\(_2\) in complementary metal-oxide-semiconductor devices has become a field of enormous interest.\(^1\) In this letter, we investigate the band-offset characteristics of high-\(k\) dielectrics on silicon. We utilize internal photoemission spectroscopy, a simple optical method developed in the 1960s,\(^4\)\(^5\) which has seen recent renewed interest in order to gain information about barrier heights, trap states and interface dipoles in high-\(k\) dielectrics.\(^6\) In this technique, a bias is applied across a dielectric structure, while tunable monochromatic light shines on the sample. At a threshold photon energy, electrons from the substrate (or metal gate) are excited by internal photoemission over the dielectric barrier.\(^7\) This threshold energy corresponds to the barrier height of the dielectric. Using bias-dependent internal photoemission spectroscopy, we have determined a barrier height profile as a function of voltage. By measuring the barrier height at both positive and negative voltages, band offsets with respect to silicon (and also the metal gate) can be determined in addition to the flatband voltage and barrier asymmetry at 0 V.

In our experimental system, we utilize a 1000 W Hg–Xe lamp with a monochromator as our light source. We use a voltage source/femtoammeter to apply the bias across the sample and to measure the current at each bias. The system is computer controlled by LabView so that the light can be scanned between 1 and 6 eV at any bias and photon energy step size. A multifunction optical meter is used to determine the lamp output spectrum to normalize the photoemission yield. Fused silica lenses are used to focus the light onto the top gold contact of the sample, which is held vertically.

The dielectric samples are grown on degenerately phosphorus doped \(n\)-type silicon to minimize the voltage drop across the depletion region in the silicon, and enhance the accuracy of our measurement. The dielectrics presented in this letter are HfO\(_2\) and Al\(_2\)O\(_3\) grown by atomic layer deposition.\(^8\)\(^9\) Before deposition, samples were dipped in a 5% HF solution for 30 s followed by a 3 min UV/ozone cleaning. Al\(_2\)O\(_3\) films were grown using de-ionized water (DI H\(_2\)O) and trimethylaluminum (Al\([\text{CH}_3\text{]_3}\)), while HfO\(_2\) films were grown using DI H\(_2\)O and tetrakis(diethylamido)hafnium (Hf\([\text{NEt}_2\text{]_4}\)). Nitrogen was used as the carrier gas, and the deposition temperature was 225 °C. The top electrode is a 12 nm layer of evaporated gold, which is sufficiently transparent so that the light source can photoexcite carriers in the silicon. The back contact is indium.

As a calibration for our experimental system, we analyzed a 15 nm thermally grown SiO\(_2\) film deposited on \(n\)-Si (1–10 \(\Omega\) cm doping). The photocurrent was measured as the photon energy was scanned at many voltages between −10 and 10 V. Figure 1 shows the raw photocurrent versus photon energy spectra for a variety of voltages (from −0.5 to +3.0 V). It is evident from these plots that the sign of the photocurrent depends strongly on the applied voltage, and the peaks in photocurrent correspond with the peaks in Hg–Xe lamp output (see Fig. 1 inset). For voltages from −10 V up to +0.7 V, negative currents were observed. At +0.7 V, the photocurrent switched signs and was positive for all higher positive voltages. We suggest that for positive photocurrents, collected electrons originate in the silicon, while the negative photocurrents indicate that electrons are mainly being generated in the metal gate contact. We determined

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0.7 V to be the voltage where the currents from the metal matched those coming from the semiconductor. We expect this voltage to be close to flatband for single dielectric layers.

The next analysis step is to calculate the photocurrent yield

$$Y = \frac{I \cdot \hbar \omega}{P},$$  \hspace{1cm} (1)

where $I$ is measured current in amperes, $P$ is the absorbed light power in watts, $\hbar \omega$ is the photon energy in eV, and $Y$ is the yield in electrons/photon. Each individual current versus photon energy curve is divided by the incident photon energy spectrum. The square root, cube root, or 2/5 power of the yield is then plotted versus photon energy as is shown in Fig. 2 for the 15 nm SiO$_2$ sample. The x intercept is then extracted and is reported as the band offset relative to the valence band of silicon. We assume that most electrons are emitted from the valence band since the number of filled initial states is much higher than in the conduction band. The literature is in general agreement that taking the square root is the appropriate power for intercept extraction when considering electrons emitted from a metal.$^{10,11}$ The case where electrons are emitted from the semiconductor is not well resolved. Semiclassical calculations suggest that taking a $2/5$ power of the yield is correct and a quantum mechanical correction to the theory predicts that the cube root is correct$^{13}$ when analyzing photocurrent from the valence band of a semiconductor.$^{14}$ In order to most thoroughly report the relevant results, we have computed offsets based on models assuming both the $2/5$ power and the cube root of the yield for these situations.$^{15}$

After extracting band offsets for each voltage, we obtain a barrier height profile as a function of voltage, as can be seen in Fig. 3 for 15 nm SiO$_2$. The points on the left of the vertical dotted line are for electrons emitted from the metal, the points on the right are for electrons emitted from the semiconductor. As can be seen, the band offset varies greatly with applied voltage, and this illustrates that it is of the utmost importance to report a corresponding bias voltage associated with a measured band offset. We report our band offsets as the point nearest to flat band, where electrons are coming from the semiconductor, or in the case of our SiO$_2$ film, 4.13± 0.1 eV. When we subtract the 1.1 eV SiO$_2$ band gap, we find that the Si/SiO$_2$ conduction band offset is 3.03± 0.1 eV. Ultimately, our results for SiO$_2$ fit well to what is expected for SiO$_2$ films with image potential barrier lowering.$^{11}$ The image potential barrier lowering simulations are shown by the black lines and can be represented by the equation (in energy units)

$$V_i = -1.15 \lambda x^2 / \lambda (s-x),$$  \hspace{1cm} (2)

where

$$\lambda = \epsilon^2 \ln (2/8 \pi \epsilon x),$$  \hspace{1cm} (3)

$\epsilon$ is the electron charge, and $s$ and $\epsilon$ are the thickness and dielectric constant of the insulating layer. The distance of the electron from the first (source) electrode is $x$. The method for approximating image barrier lowering is described in Ref. 16. The image force was approximated as that of an insulator between two metallic electrodes. In general, the silicon substrates used in our experiments are highly doped, so this is a reasonable approximation although the thin depletion layer in the Si will somewhat modify the result. For SiO$_2$, we assumed an optical dielectric constant of 2.5.

Similar analyses were completed for Al$_2$O$_3$ and HfO$_2$ films on n-Si (0.0001 $\Omega$ cm). The resulting barrier height
profiles and barrier height simulations are shown in Fig. 4. The barrier height profiles for these two materials are not nearly as clear as for SiO$_2$. This is particularly true for HfO$_2$—a result of leakage through the barrier and greater difficulty in extracting band offsets from the yield curves. Transmission electron microscopy analysis revealed that interfacial SiO$_2$ layers 2.2 nm thick were present between the HfO$_2$ and Al$_2$O$_3$ dielectric and semiconductor. These interfacial layers could be attributed to the UV/ozone clean during the substrate preparation or a postdeposition 600 °C anneal in Ar+$\text{O}_2$ ppm, and could account for the higher measured band offsets compared with literature values, though the electrical characteristics of these layers are unknown.

Because of the presence of interfacial layers in the Al$_2$O$_3$ and HfO$_2$ samples, the analysis for these films is slightly more complicated than for SiO$_2$. The measured barrier height (from the Si valence band) for 16.1 nm Al$_2$O$_3$ is 4.6±0.1 eV (at 1.0 V). After subtracting the Si band gap (1.1 eV), the Al$_2$O$_3$ conduction band offset is found to be 3.5±0.1 eV (at 1.0 V), but this probably corresponds most directly to the interfacial layer than to the Al$_2$O$_3$ layer itself. For this reason, another quantity of interest is the Au/Al$_2$O$_3$ barrier height (from the metal (from −0.5 to 0.9 eV) and we can determine a Au/HfO$_2$ barrier height of 3.6±0.1 eV. Using the same reasoning as for Al$_2$O$_3$, based on our data, our best approximation for the Si/HfO$_2$ barrier height (from valence band) is 3.8±0.2 eV. This corresponds to a conduction band offset with respect to Si of 2.7±0.2 eV. The experimental literature reports conduction band offsets of 2.0 and −1.2 eV for HfO$_2$, while theoretical calculations predict 1.5 eV. Further efforts of barrier lowering in HfO$_2$/Al$_2$O$_3$ heterostructures will be addressed in a paper to be published soon (see Ref. 21 for background).

The high-$\kappa$ field has been plagued by the growth of unwanted interfacial layers at the dielectric/silicon interface. Progress is being made to decrease these layers by careful cleaning and passivation of the silicon surface before growth, but eliminating these layers completely remains a great challenge. Although interfacial layers will in general modify the measurements, by careful use of our barrier profile technique, and by selecting the barrier height at specific voltages (and thus specific band alignment) we can gain a very good estimate of the band offset of any barrier. We have shown that profiling the barrier heights of dielectrics on silicon as a function of applied bias is a very viable and valuable technique for determining the effective band offset at any particular voltage.

Similar analysis can be done for HfO$_2$ (open symbols in Fig. 4), but because of the degraded data quality attributed to leakage from the substrate, it is more difficult to verify our data by simulation. The dashed curve indicates a barrier lowering simulation for the HfO$_2$ barrier (HfO$_2$ $\kappa$=22, SiO$_2$, $\kappa$=3.9). The data quality is good when the electrons originate from the metal (from −0.5 to 0.9 eV) and we can determine a Au/HfO$_2$ barrier height of 3.6±0.1 eV. Using the same reasoning as for Al$_2$O$_3$, based on our data, our best approximation for the Si/HfO$_2$ barrier height (from valence band) is 3.8±0.2 eV. This corresponds to a conduction band offset with respect to Si of 2.7±0.2 eV. The experimental literature reports conduction band offsets of 2.0 and −1.2 eV for HfO$_2$, while theoretical calculations predict 1.5 eV. Further efforts of barrier lowering in HfO$_2$/Al$_2$O$_3$ heterostructures will be addressed in a paper to be published soon (see Ref. 21 for background).