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Numerical simulation of a solar cell based on a double CdTe/FeSi₂ absorber layer and an AZO window layer with an efficiency of 26.34 %

Solar Cell simulation serves to reduce the number of work hours and optimize the resources dedicated to carrying out experiments. Simulation of solar cells based on AZO as window layer, cadmium telluride (CdTe) like first absorber layer and FeSi₂ like second absorber layer with the structure (Al/AZO/CdTe/FeSi₂/Ni) using the Cell Capacitance Simulator solar-1D (SCAPS-1D) has been carried out. The effect of the thickness and carrier concentration of the first and the second absorber layer was studied. The thickness and concentration of carriers for the CdTe layer varied from 0.4 to 1 μm and from 10^{13} to 10^{15} cm^{-3} respectively, it was found that PCE decreases 0.22% with increasing thickness. The thickness and concentration of carriers in the second absorber layer varied from 0.1 to 1 μm and from 10^{14} to 10^{18} cm^{-3} , respectively. It was observed that by increasing the thickness of the FeSi₂ layer from 0.1 to 1 μm , PCE increased by 7.66%, while by increasing the concentration of carriers from 1014 to 1018 cm^{-3} in this layer, PCE increased by 2.81%. It was also observed that by increasing the defect density at the FeSi₂/CdTe interface from 1010 to 1017 cm^{-2} , PCE decreased by 22.1%. Finally, a maximum power conversion efficiency (PCE) of 26.34% is achieved with an open circuit voltage (Voc) of 0.62 V, short circuit current density (Jsc) of 51.43 mA/cm^2 and a fill factor (FF) of 82.24 %, obtained with a thickness and concentration of carriers of 1 μm and 10^{18} cm^{-3} for the FeSi₂ layer and for the CdTe layer the thickness and concentration of carriers was 0.4 μm and 10^{15} cm^{-3} , respectively. The above results provide design guidelines for a solar cell based on two absorber layers CdTe and FeSi₂ and window layer of AZO with high efficiency.

Keywords

solar cells simulation, two absorber layer, SCAPS-1D

Reference

Performance simulation of solar cell based on AZO/CdTe heterostructure by SCAPS 1D software. José Carlos Zepeda Medina. Heliyon, 9 (marzo 2023) e14547

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