

Supplementary Materials

Optimization of the Sb₂S₃ Shell Thickness in ZnO Nanowire-Based Extremely Thin Absorber Solar Cells

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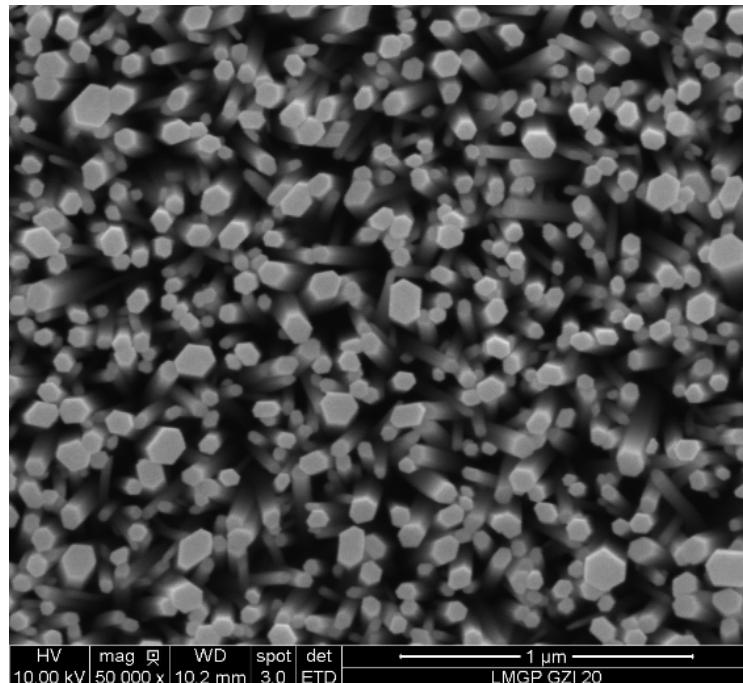


Figure S1. Typical top-view FESEM image of ZnO/TiO₂ core-shell NW heterostructures without any Sb₂S₃ shell.

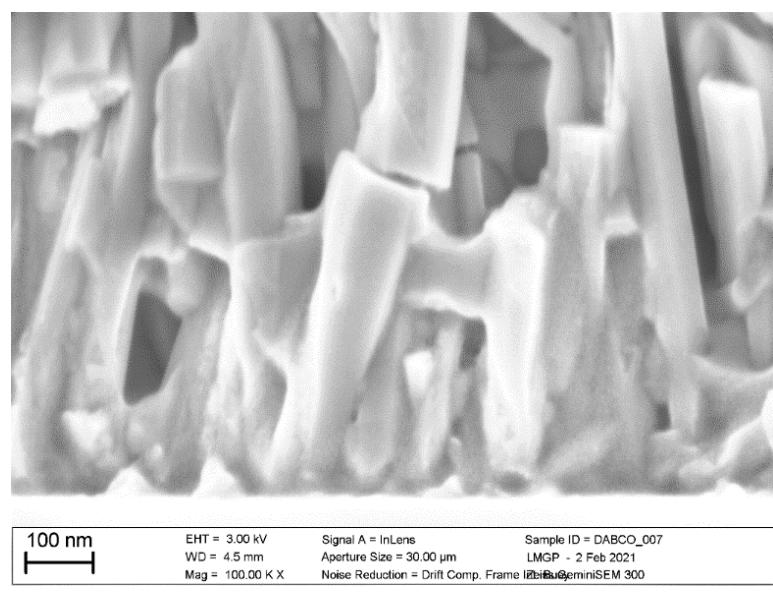


Figure S2. High magnification cross-sectional view FESEM image of ZnO/TiO₂ core-shell NW heterostructures covered by a Sb₂S₃ shell grown by CSP using 70 cycles.

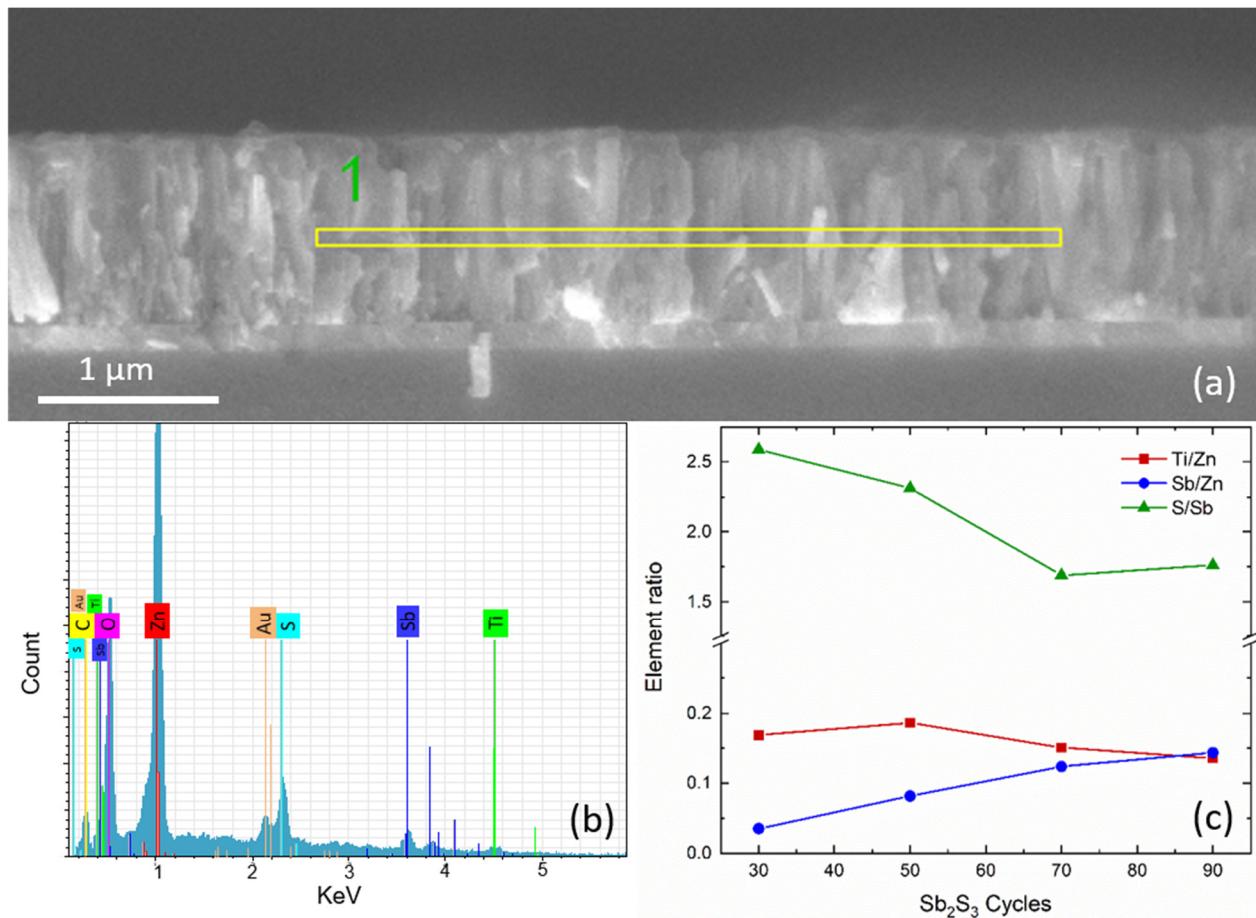


Figure S3. (a) Cross-sectional view FESEM image of the complete ETA solar cell structure of ITO/ZnO/TiO₂/Sb₂S₃/P3HT/Au when using the cycle number of 70, (b) FESEM-EDX spectra collected on the yellow rectangular area as denoted in (a), (c) Element ratio as a function of the cycle number ranging from 30 to 90.

Table S1. In-plane XRD data recapitulating the nature and position of the diffraction peaks in this work as compared to the 00-042-1393 ICDD file.

(hkl)	2-Theta (°) This Work	2-Theta (°) 00-042-1393 ICDD file
(220)	22.20	22.28
(130) – (310)	24.93	24.89 – 25.01
(111)	25.65	25.73
(121) – (211)	29.20	29.25
(301)	33.38	33.39
(131) – (311)	34.19	34.24 – 34.34
(240) – (420)	35.54	35.52 – 35.64
(231)	36.76	36.99
(340) – (430) – (141) – (411) – (510)	40.17	39.91 – 40.00 – 40.38 – 40.57 – 40.89
(250) – (520)	42.98	43.04 – 43.30
(530) – (002) / (151) – (511)	46.80	47.04 – 47.30 – 47.54
(531)	52.84	53.04
(061) – (132) / (312) – (360)	54.00	54.13 – 54.20 – 54.41 – 54.65
(630)	54.62	54.65

Table S2. Raman scattering data recapitulating the nature and position of the Raman lines in this work as compared to Ref. [1].

Wavenumber (cm ⁻¹) This Work	Wavenumber (cm ⁻¹) Ref. [1]	Vibrational modes
61	65.7 / 65.8	B _{1g} / B _{3g}
71	74.7	A _g
90	92.9	B _{2g}
100	100	A _g
115	112	B _{2g}
128	122 / 128	B _{2g} / A _g
157	159 / 159	B _{2g} / A _g
192	190	A _g
239	225 / 227	B _{1g} / B _{3g}
283	280	A _g
300	291 / 293	B _{2g} / A _g
314	294	B _{2g}

References

- Ibanez, J.; Sans, J.A.; Popescu, C.; Lopez-Vidrier, J.; Elvira-Betanzos, J.J.; Cuenca-Gotor, V.P.; Gomis, O.; Manjon, F.J.; Rodriguez-Hernandez, P.; Munoz, A. Structural, Vibrational, and Electronic Study of Sb₂S₃ at High Pressure. *J. Phys. Chem. C* **2016**, *120*, 10547–10558. <https://doi.org/10.1021/acs.jpcc.6b01276>.