

Large-area synthesis of monolayer $\text{MoS}_{2(1-x)}\text{Se}_{2x}$ with a tunable band gap and its enhanced electrochemical catalytic activity

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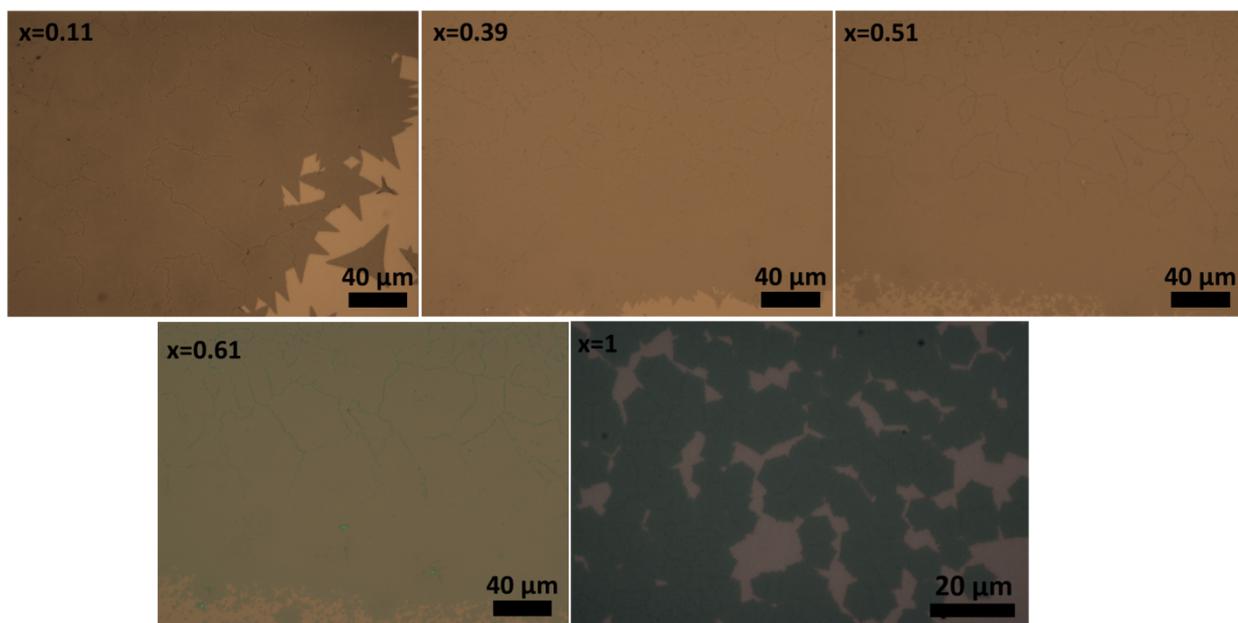


Fig. S1 Optical images of monolayer $\text{MoS}_{2(1-x)}\text{Se}_{2x}$ with different S composition.

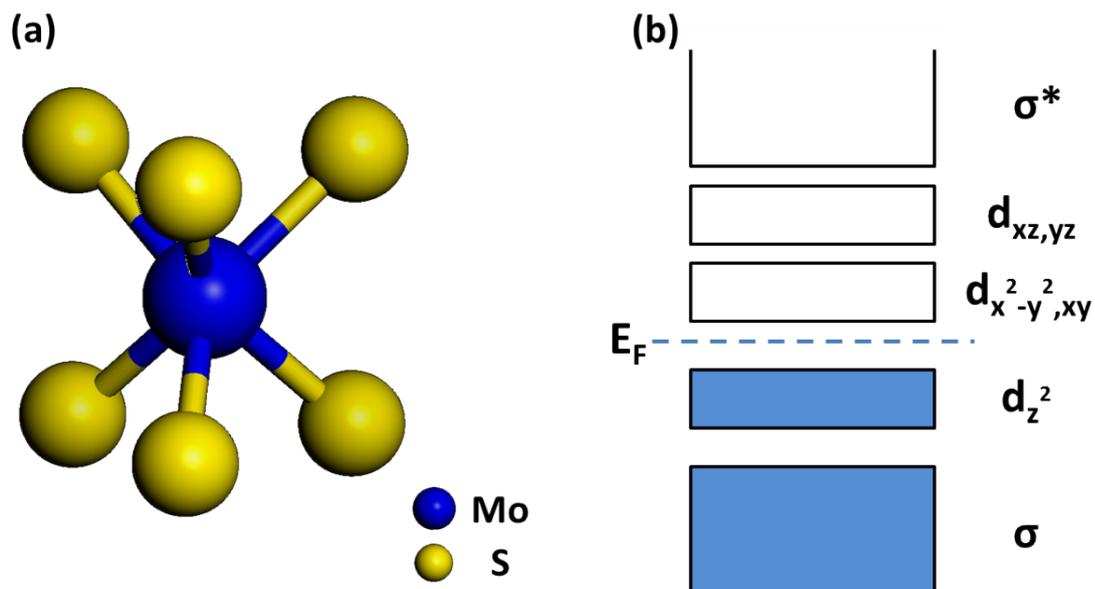


Fig. S2 (a) Schematic structure of monolayer MoS₂. The Mo atoms are bonded with six S atoms symmetrically. (b) Schematic illustration of the band structure of 1H MoX₂ (X=S, Se, Te). The non-bonding *d* bands are located within the gap of bonding (σ) and antibonding (σ^*) bands. From bottom to top, the non-bonding *d* bands are Mo- d_{z^2} , Mo- $d_{x^2-y^2, xy}$ and Mo- $d_{xz, yz}$ orbitals, respectively. The filled states are shaded with blue. The gap between the Mo- d_{z^2} and Mo- $d_{x^2-y^2, xy}$ orbitals are corresponding to the band gap of 1H MoX₂.^[S1]

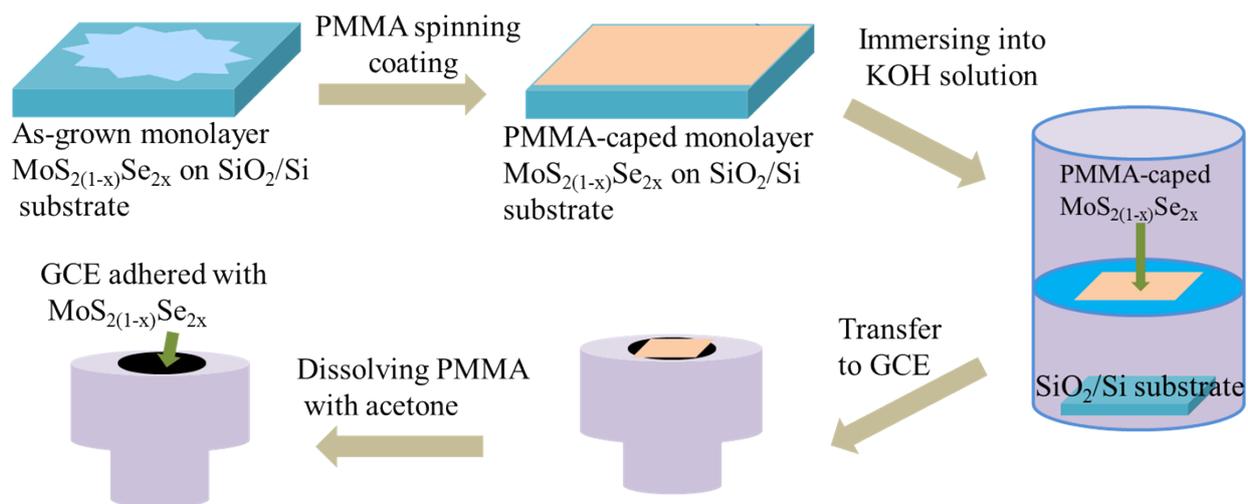


Fig. S3 Schematic illustration of transferring the as-grown monolayer $\text{MoS}_{2(1-x)}\text{Se}_{2x}$ onto GCE.

References

S1 M. Chhowalla, H. S. Shin, G. Eda, L.-J. Li, K. P. Loh and H. Zhang, *Nat. Chem.*, 2013, **5**, 263–275.