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# Small Micro

## Supporting Information

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Phosphorus Enhanced Intermolecular Interactions of  $SnO_2$  and Graphene as an Ultrastable Lithium Battery Anode

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### Supporting Information

#### Phosphorous enhanced Intermolecular Interactions of SnO<sub>2</sub> and Graphene as an Ultrastable

#### Lithium Battery Anode

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Figure S1. SEM images of SnO<sub>2</sub>@P@GO.



Figure S2. TEM images of SnO<sub>2</sub>/GO



Figure S3. SEM images of SnO<sub>2</sub>/P/GO



**Figure S4**. Nitrogen adsorption-desorption isotherms and pore size distributions (inset) of A) SnO<sub>2</sub>@P@GO and B) SnO<sub>2</sub>/P/GO.



Figure S5. XPS spectrum of SnO<sub>2</sub>/P/GO: A) O1s spectrum, B) P1p spectrum.



Figure S6. TG curves of SnO<sub>2</sub>@P@GO in air.



Figure S7. Charge-discharge curves of SnO<sub>2</sub>/P/GO at 100 mA g<sup>-1</sup>.



Figure S8. CV curves of SnO<sub>2</sub>@P@GO at different scan rate ranging from 0.1 to 1 mV s<sup>-1</sup>.



Figure S9. Kinetic analysis of  $SnO_2@P@GO$ : A) Separation of the capacitive and diffusion currents at a scan rate of 1 mV s<sup>-1</sup>. B) Contribution ratio of the capacitive and diffusion-controlled charge at various scan rates.



Figure S10. Cycling performance of bare P at 100 mA g<sup>-1</sup>.

Table S1 ICP result	
	Mass Ratio
Sn	5.4
Р	1



Figure S11. Ex-situ chemical composition analysis by XPS for A) O and B) P.