

Supplementary Information

Rice Husk-Derived Porous Silicon Dioxide Fillers for Enhancing Ionic Conductivity in a Solid-State Electrolyte of Lithium-Sulfur Batteries Under Molecular Dynamic Calculation

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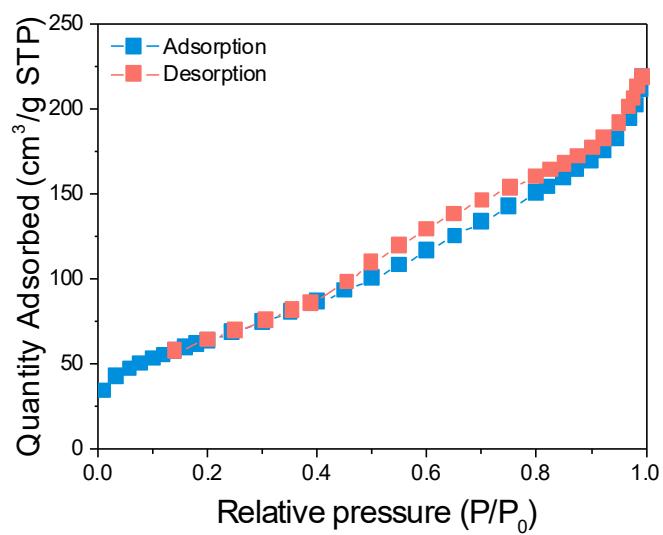


Fig. S1 The BET analysis of RH_{SiO₂}.

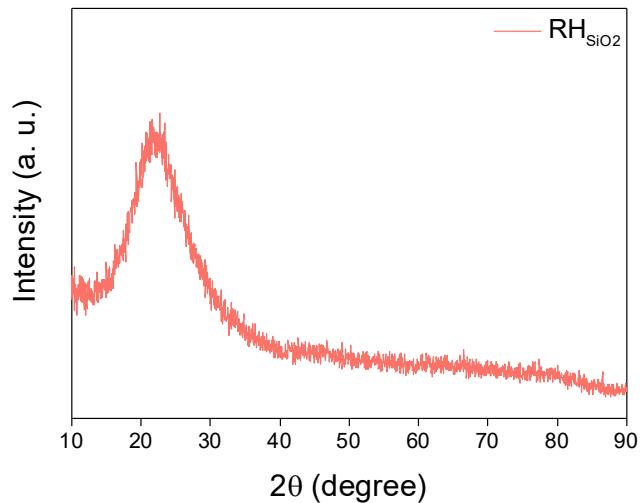
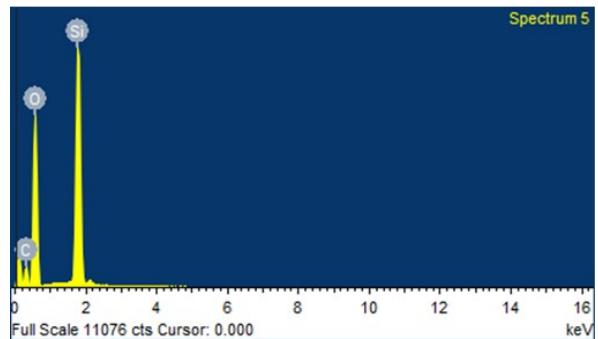


Fig. S2 The XRD analysis of RH_{SiO₂}.



Element	Weight%	Atomic%
C K	17.00	23.95
O K	57.20	60.50
Si K	25.80	15.55
Totals	100.00	

Fig. S3 The EDX analysis of RH_{SiO₂}.

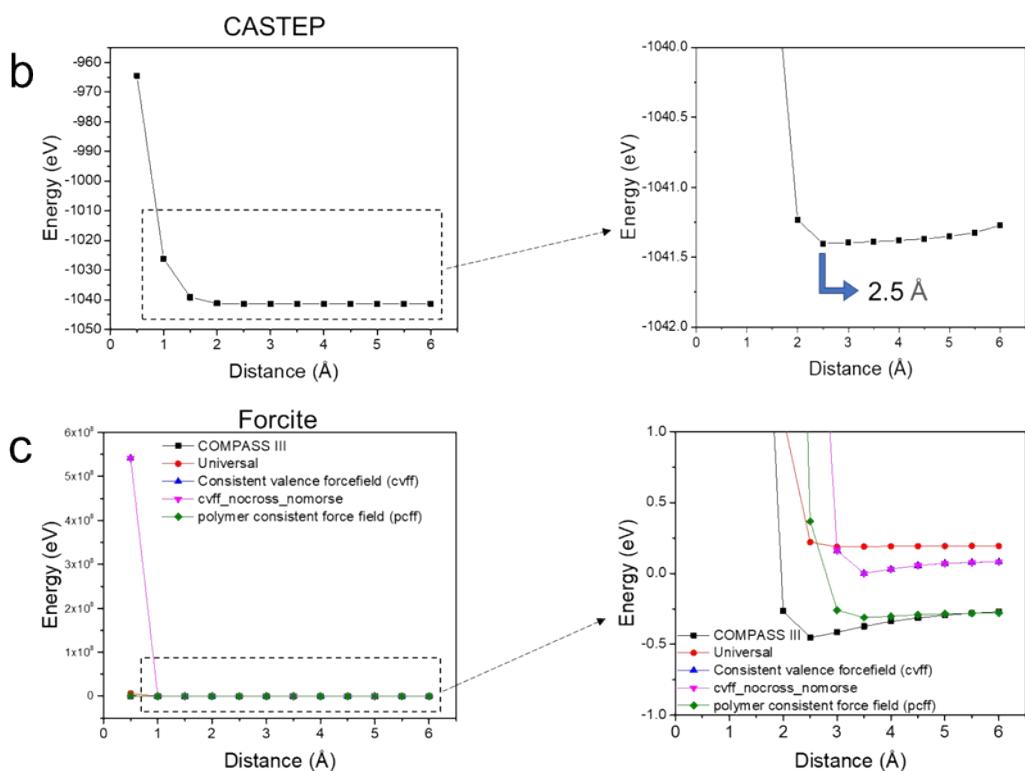
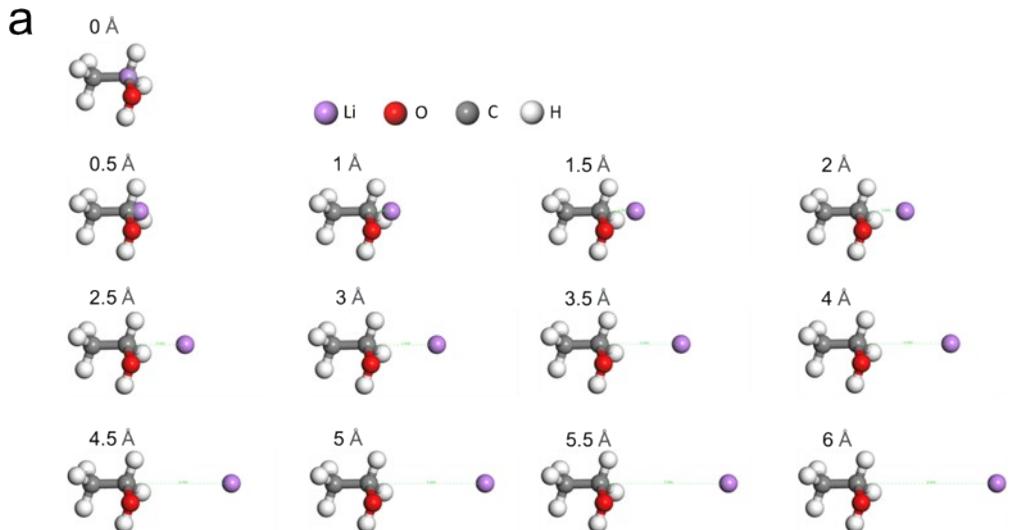


Fig. S4 (a) The calculation model of ethylene oxide monomer and Li-ion with various distances.

The energy curve as a function of the distance by (b) CASTEP calculation and (c) Forcite calculation.

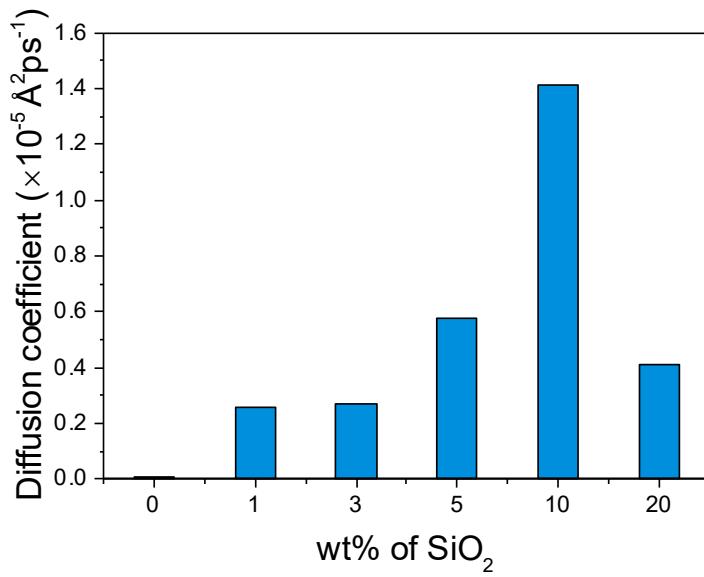
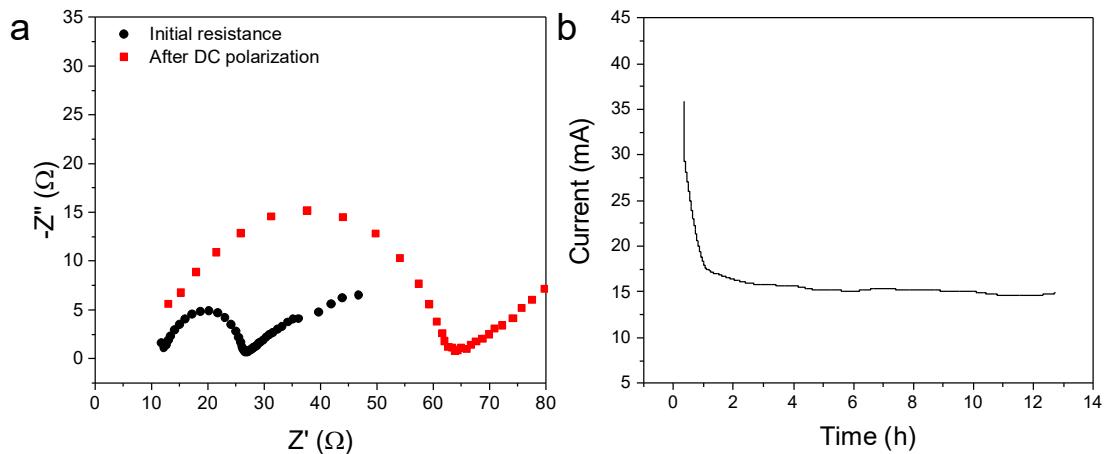
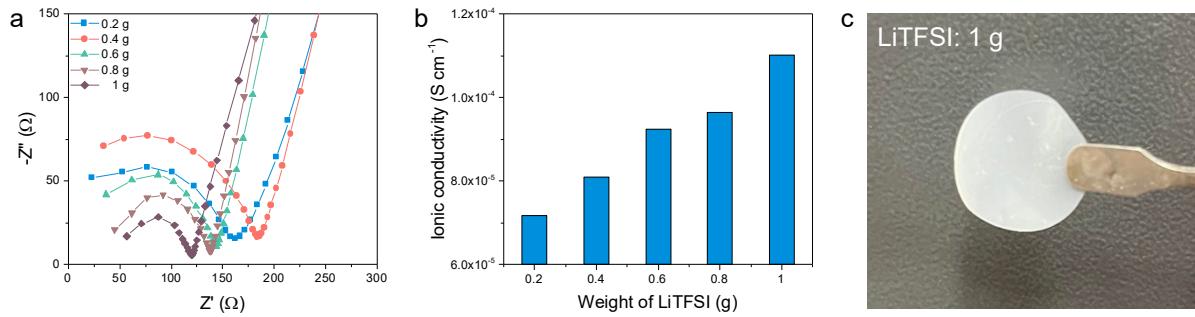


Fig. S5 The diffusion coefficients of CPEs with various content of RH_{SiO₂} from the molecular dynamics simulation.



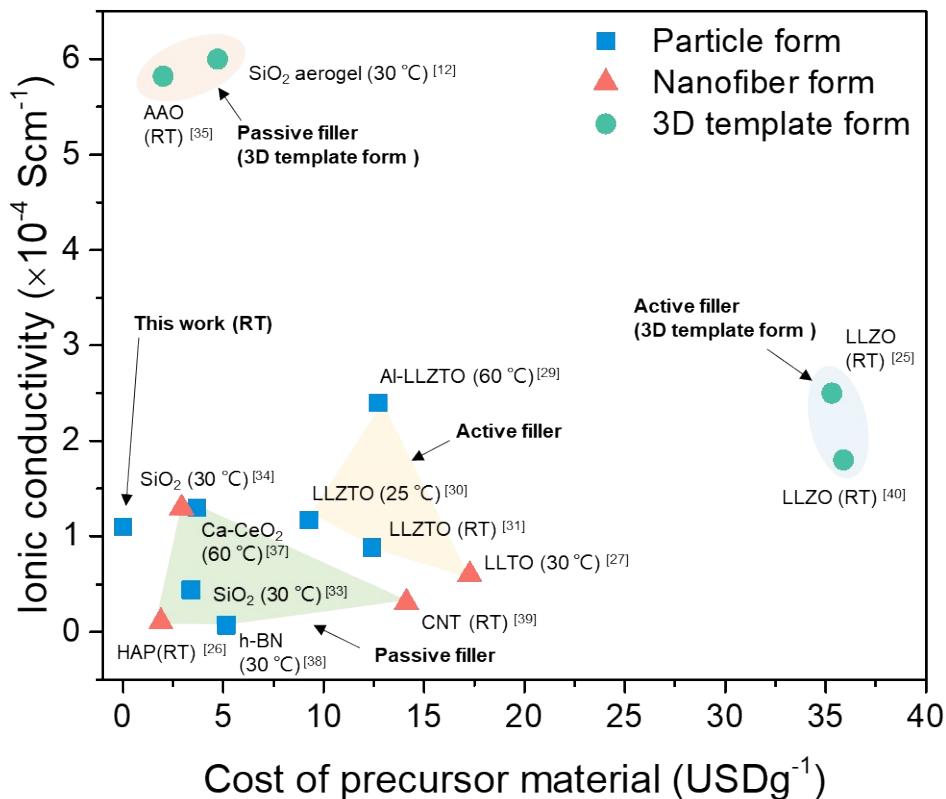


Fig. S8 The comparison of ionic conductivity versus the cost of precursor materials of our work and the recent reports in premier journals.

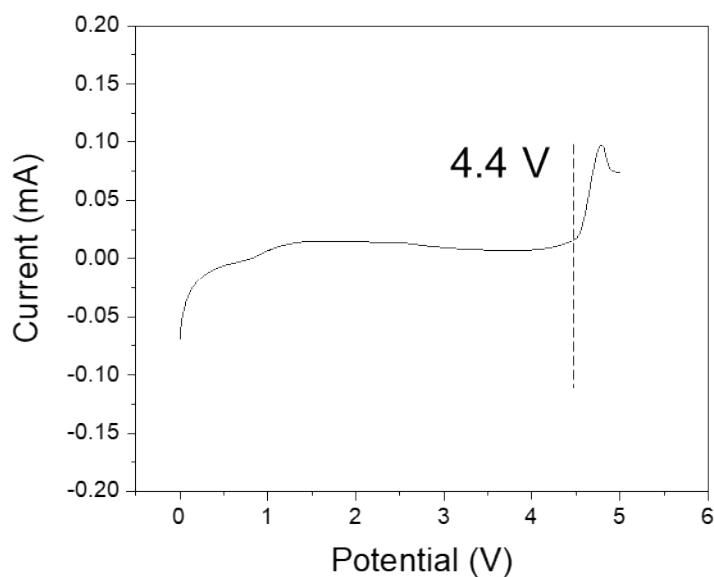


Fig. S9 The stable voltage window of RH_{SiO₂} CPE.

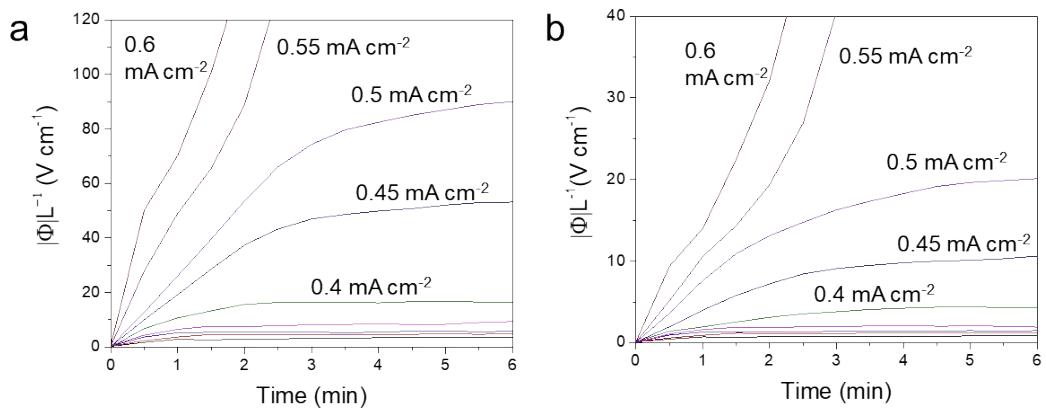


Fig. S10 The absolute value of the length-normalized potential of the (a) SPE and (b) RH_{SiO₂} CPE electrolyte for various applied current densities.

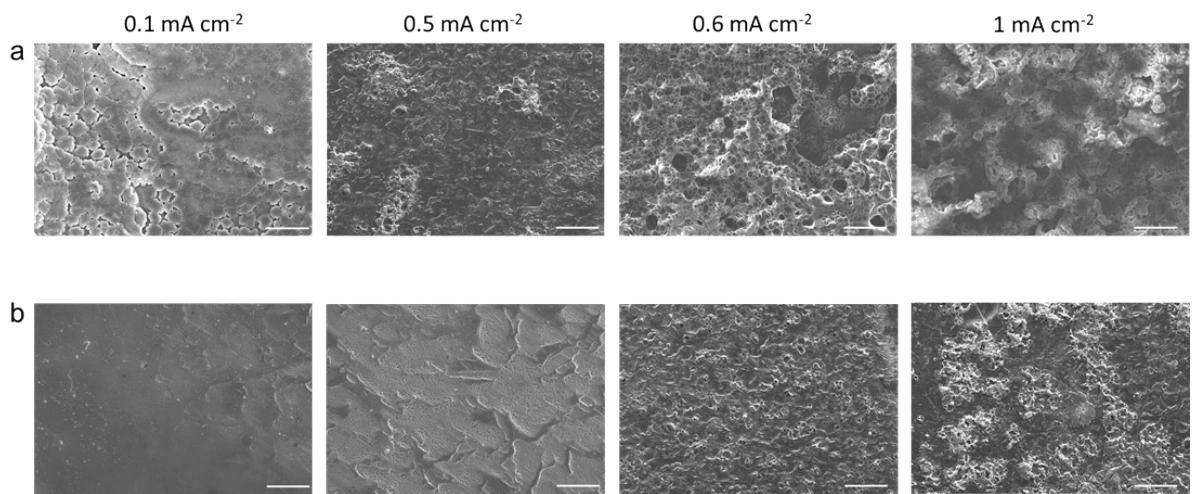


Fig. S11 The SEM images of dendrite formation with (a) SPE and (b) RH_{SiO₂} CPE at various current densities. Scale bar: 20 μ m.

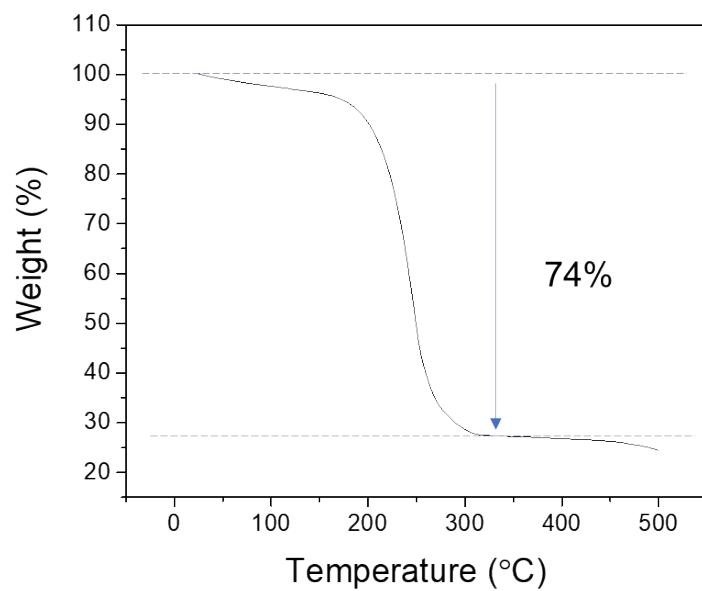


Fig. S12 The thermogravimetric analysis (TGA) of CMK-3/S material.

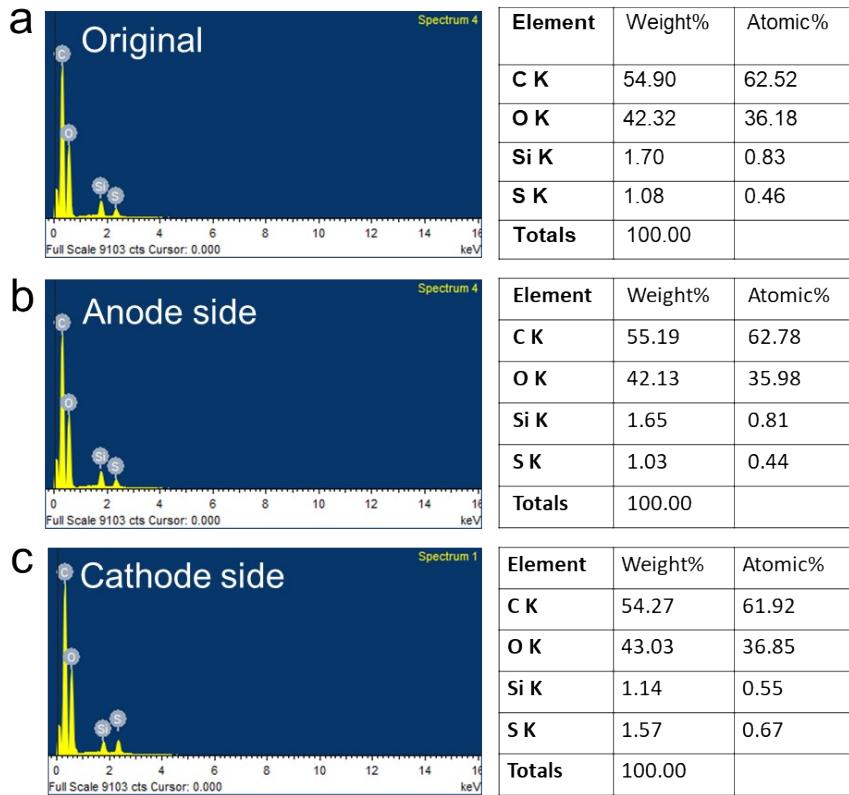


Fig. S13 The EDX analysis of the (a) original RH_{SiO_2} CPE, the (b) top side, and (c) bottom side of RH_{SiO_2} CPE after cycling.

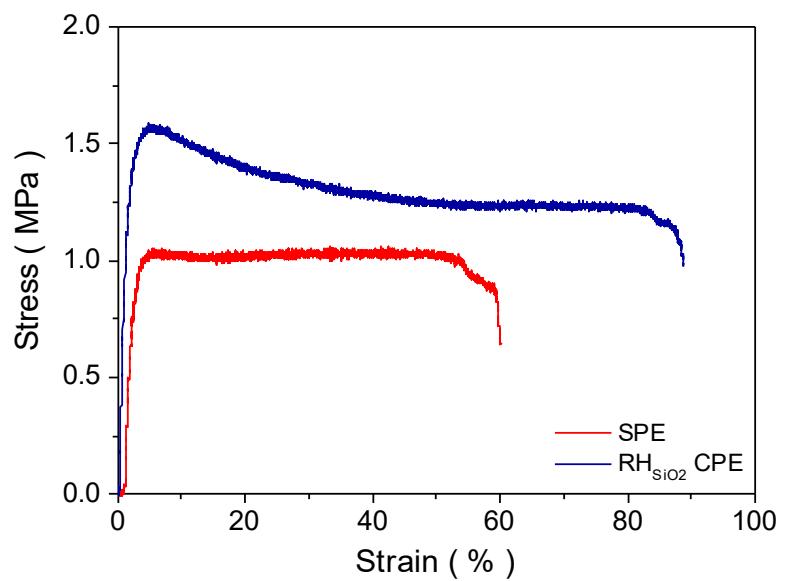


Fig. S14 The stress-strain curves of RH_{SiO2} CPE and the filler-free SPE.