

Plasma-photocatalytic conversion of CO₂ at low temperatures: Understanding the synergistic effect of plasma-catalysis

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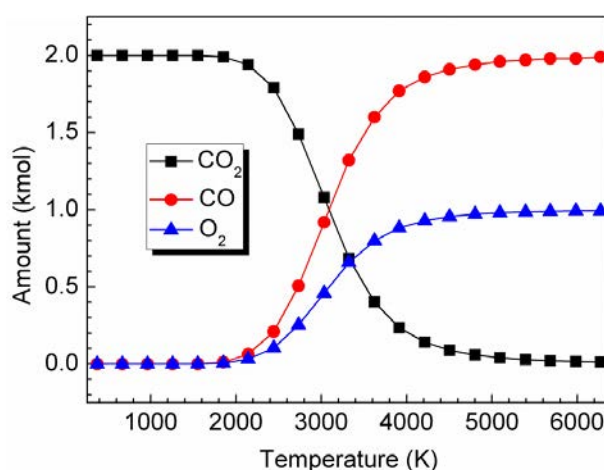
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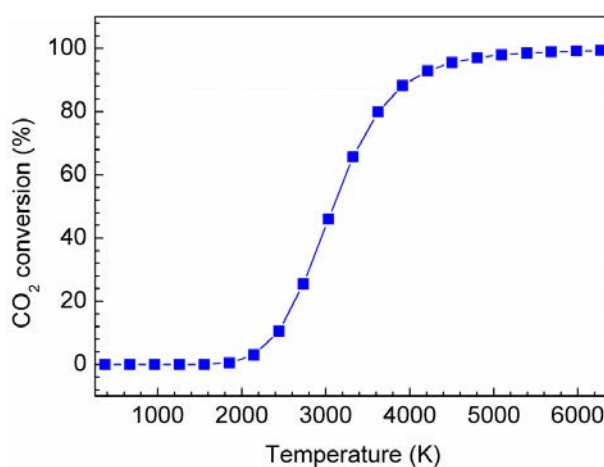
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1. Thermodynamic equilibrium calculation

The thermodynamic equilibrium calculation of CO₂ conversion was carried out using the method based on the minimization of Gibbs free energy in a closed system. The main gas products are CO and O₂. No O₃ was detected in the present experiment. We can see that CO₂ begins to decompose into CO and O₂ near 2000 K and the conversion of CO₂ is very low (< 1%). Reasonable conversion of CO₂ (~60%) can be obtained at an extraordinarily high temperature (3000-3500K), which leads to the high energy cost for thermal conversion of CO₂.



(a)



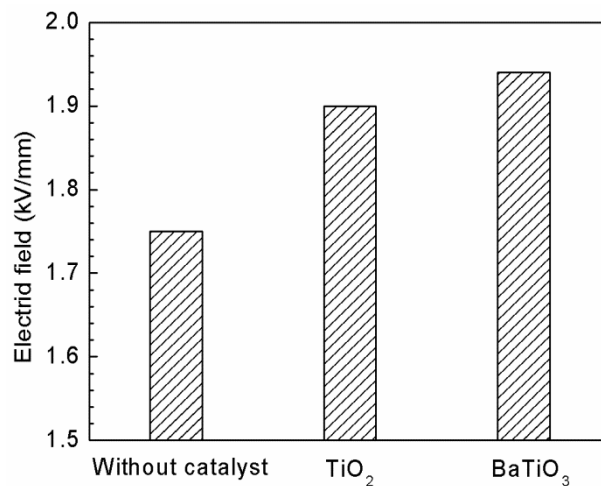
(b)

Fig. S11 Thermodynamic equilibrium calculation of CO₂ conversion as a function of

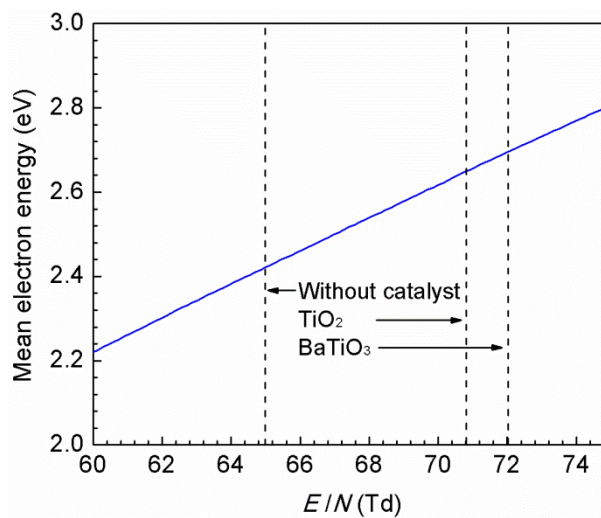
operating temperature at 1 atm (without plasma) (a) gas composition vs. temperature; (b)
CO₂ conversion vs. temperature.

2. Mean Electric field and electron energy

The average electric field of the discharge (breakdown voltage/electrode gap) and the mean electron energy for the three different experimental conditions are calculated through Lissajous figure and BOLSIG⁺ code based on electron energy distribution function (EEDF), respectively[1, 2] and the corresponding results are shown in Fig. SI2.



(a)



(b)

Fig. SI2 Average electric field (a) and mean electron energy (b) of the CO₂ DBD with and without catalyst.

3. Catalyst characterization

XRD patterns of the catalyst samples are plotted in Figure SI3. BaTiO₃ has the tetragonal phase [3], which can be approved by the peaks at $2\theta = 22.14^\circ, 31.61^\circ, 39.00^\circ, 45.37^\circ, 51.00^\circ, 56.32^\circ, 65.94^\circ$ (JCPDS 05-0626), while the fresh TiO₂ shows a crystal structure of anatase, as evidenced by X-ray reflections at $2\theta = 25.24^\circ, 36.94^\circ, 37.82^\circ, 38.56^\circ, 48.04^\circ, 53.96^\circ, 55.00^\circ, 62.69^\circ, 68.76^\circ$ (JCPDS 84-1286)[4].

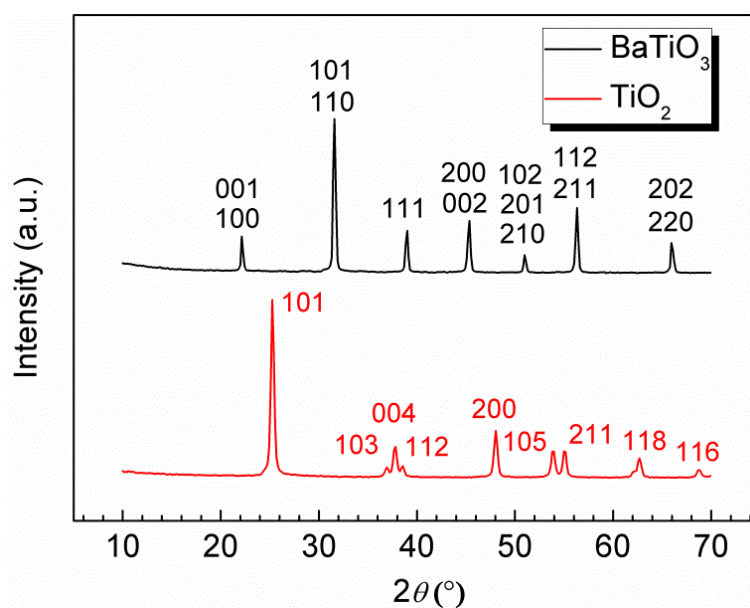


Fig. SI3 XRD patterns of the fresh catalysts.

Reference

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