

*Supporting Information for*

## **Catalytic Transfer Hydrogenation of Ethyl Levulinate to $\gamma$ -valerolactone over Supported MoS<sub>2</sub> Catalysts**

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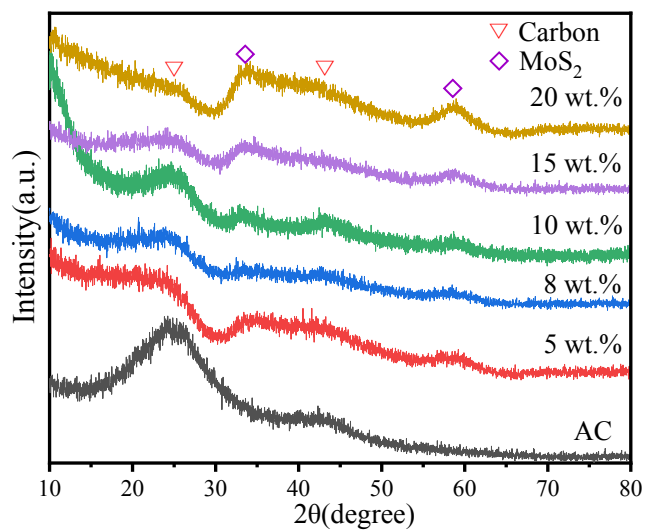
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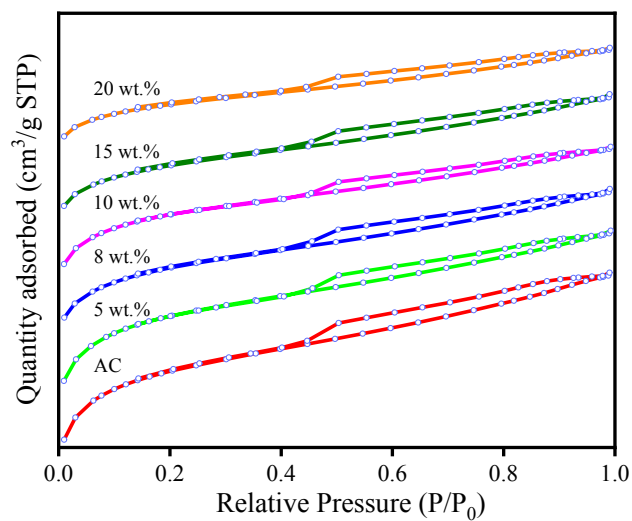
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## SECTION S1 SUPPLYMENTARY FIGURES



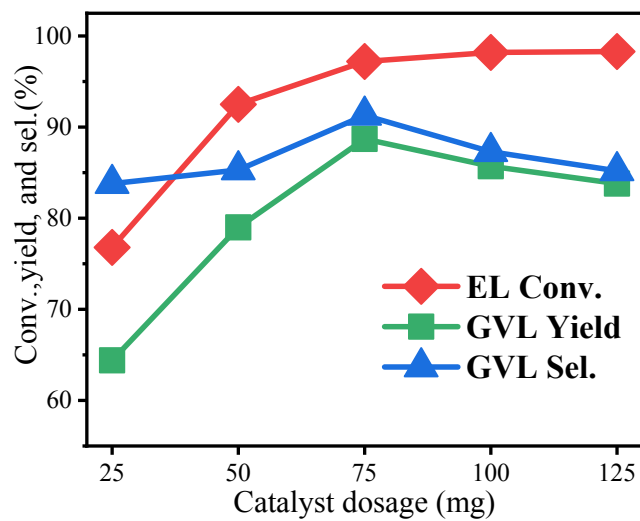
**Fig. S1** XRD patterns of the AC support and MoS<sub>2</sub>/AC catalysts with different Mo loadings.

## SECTION S1 SUPPLYMENTARY FIGURES



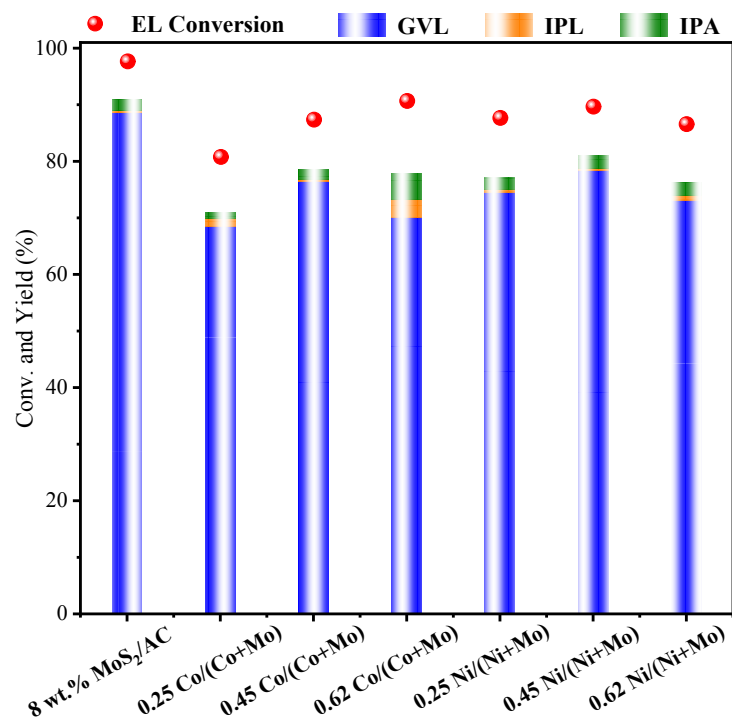
**Fig. S2** Nitrogen adsorption-desorption isotherms of the AC support and MoS<sub>2</sub>/AC catalysts with different Mo loadings.

## SECTION S1 SUPPLYMENTARY FIGURES



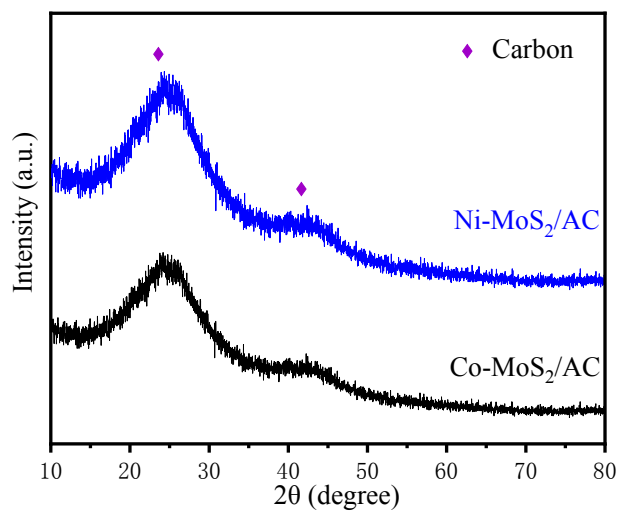
**Fig. S3** Effects of the catalyst dosage on the conversion of EL into GVL. Reaction condition: EL 2 mmol, isopropanol 20 mL, 230 °C, 1.5 h, 1 MPa H<sub>2</sub>.

## SECTION S1 SUPPLYMENTARY FIGURES



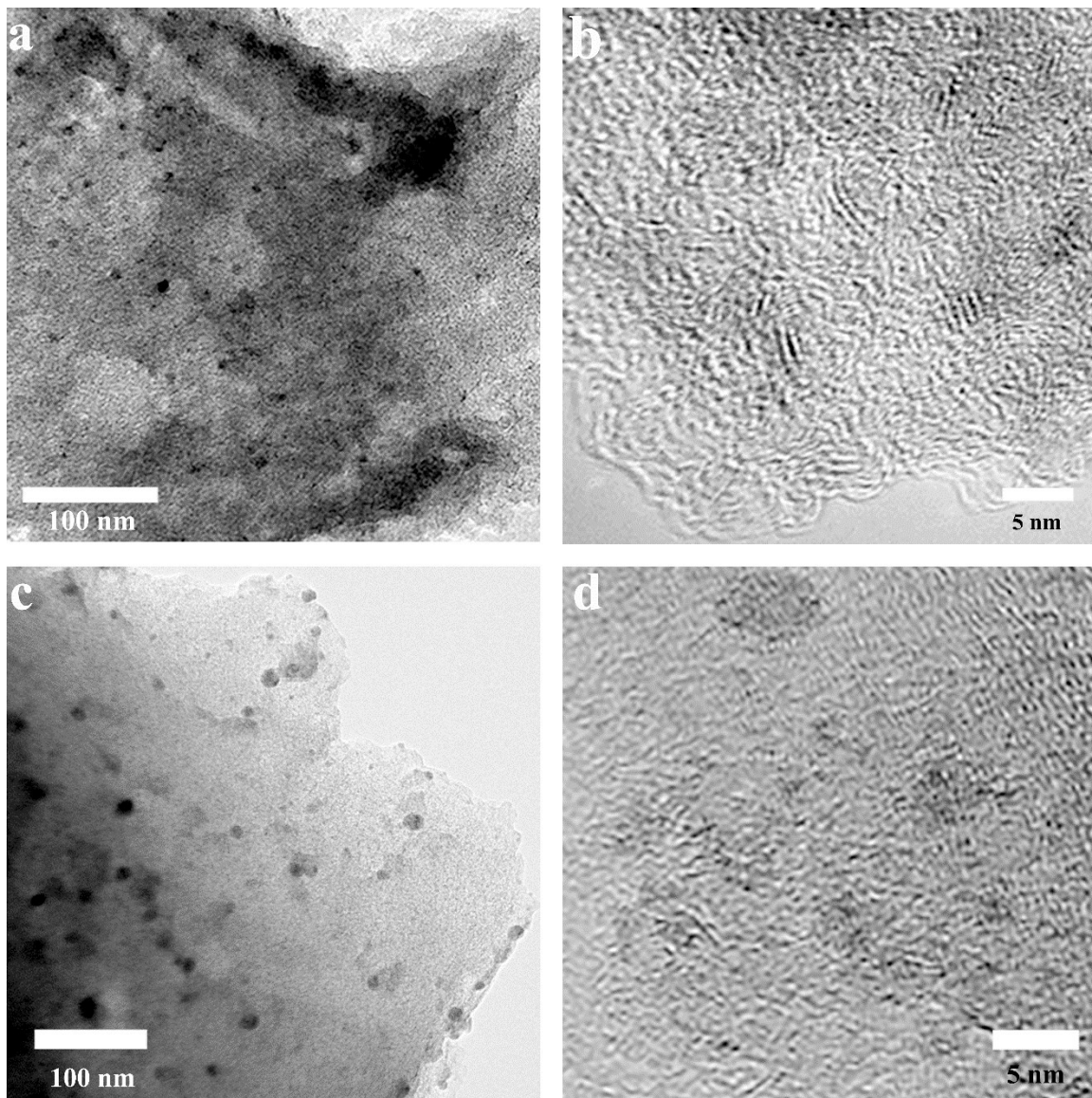
**Fig. S4** Effects of the Co/Ni doping on the conversion of EL into GVL over MoS<sub>2</sub>/AC catalyst. Reaction condition: EL 2 mmol, catalyst 75 mg, isopropanol 20 mL, 230 °C, 1.5 h, 1 MPa H<sub>2</sub>.

## SECTION S1 SUPPLYMENTARY FIGURES



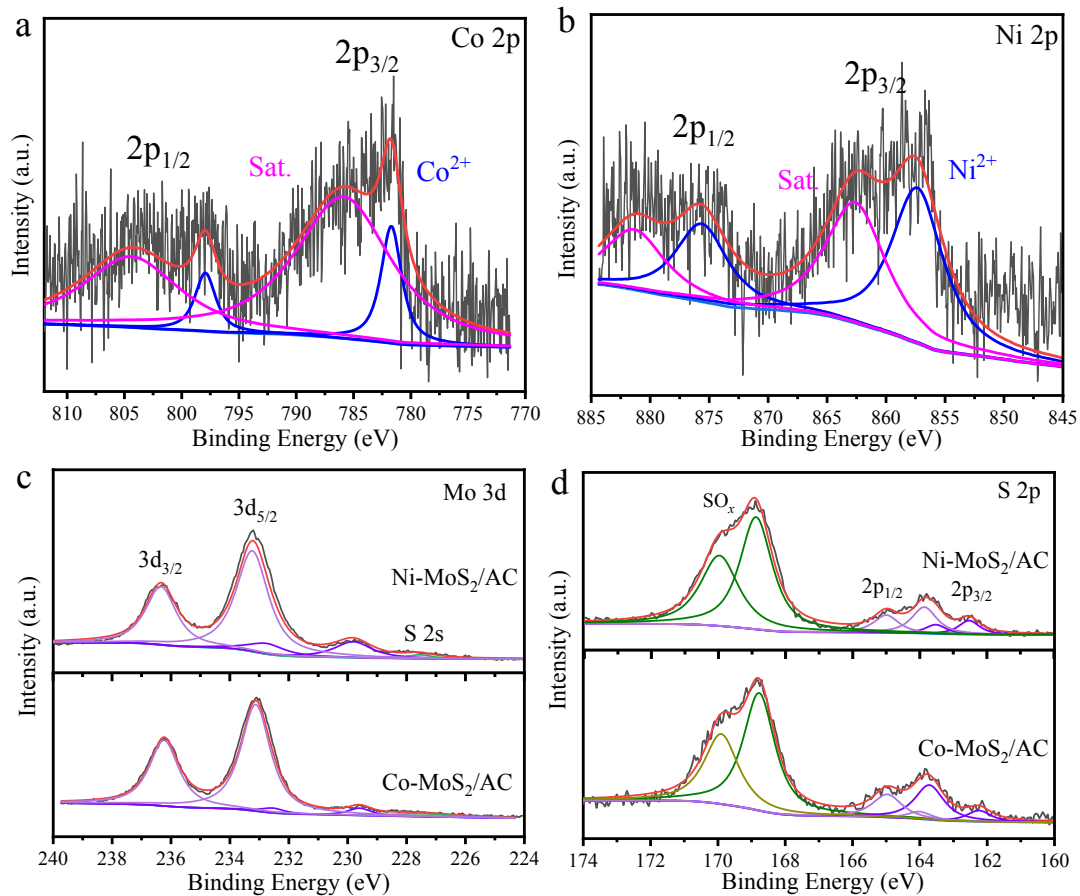
**Fig. S5** XRD patterns of Co-MoS<sub>2</sub> (Co/(Co+Mo)=0.45) and Ni-MoS<sub>2</sub> (Ni/(Ni+Mo)=0.45)

## SECTION S1 SUPPLYMENTARY FIGURES



**Fig. S6** TEM images of (a-b) Ni-MoS<sub>2</sub>/AC (Ni/(Ni+Mo) =0.45) and (c-d) Co-MoS<sub>2</sub>/AC (Co/(Co+Mo)=0.45) catalysts.

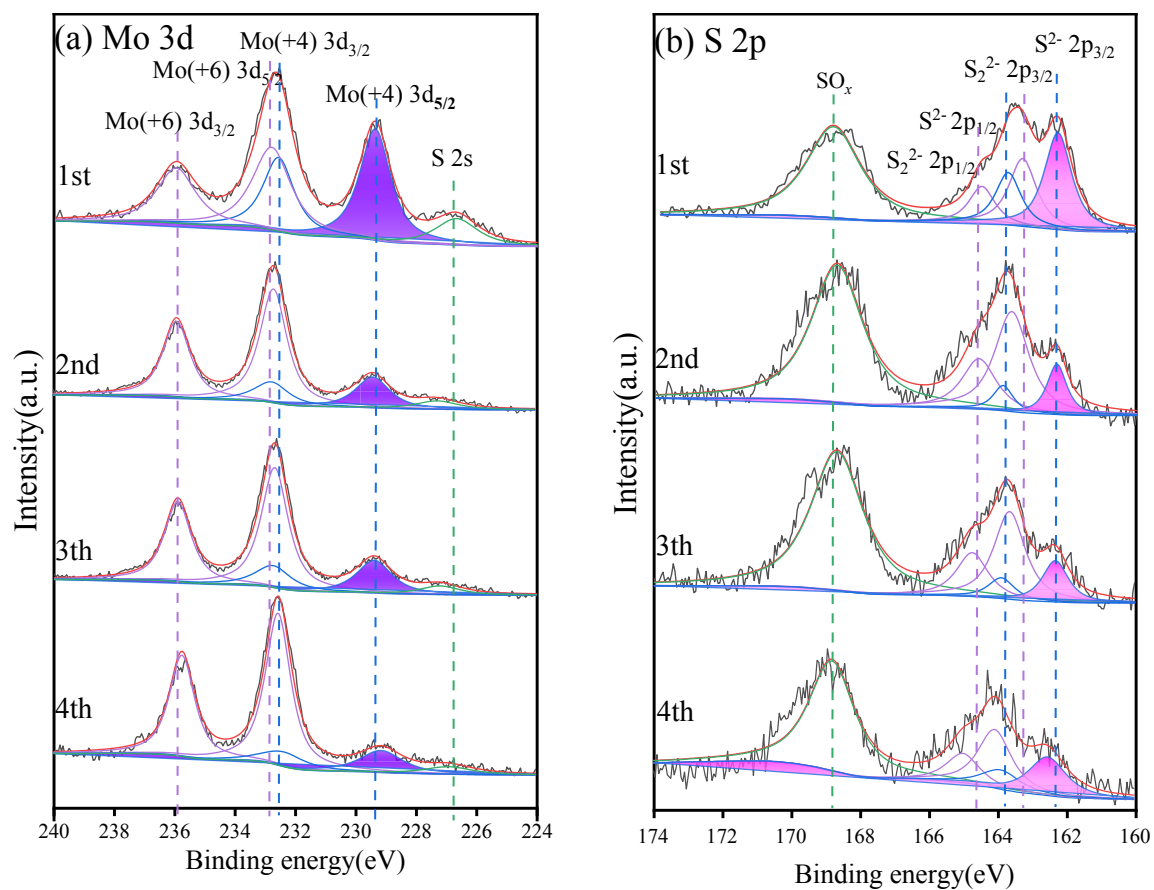
## SECTION S1 SUPPLYMENTARY FIGURES



**Fig. S7** XPS spectrum of Co-MoS<sub>2</sub>/AC (Co/(Co+Mo)=0.45) and Ni-MoS<sub>2</sub>/AC (Ni/(Ni+Mo)=0.45) catalysts.



## SECTION S1 SUPPLYMENTARY FIGURES



**Fig. S8** XPS spectrum of fresh and spent 8 wt.% MoS<sub>2</sub>/AC catalyst (a) Mo 3d and (b) S 2p.

## SECTION S2 SUPPLYMENTARY TABLES

**Table S1.** Textural properties of activated carbon and MoS<sub>2</sub>/AC catalyst

Sample	BET surface area <sup>a</sup> (m <sup>2</sup> /g)	Pore volume (cm <sup>3</sup> /g)	Pore size <sup>b</sup> (nm)
AC	1161	0.39	3.88
5 wt.% MoS <sub>2</sub> /AC	968	0.34	3.87
8 wt.% MoS <sub>2</sub> /AC	763	0.31	3.91
10 wt.% MoS <sub>2</sub> /AC	741	0.27	3.89
15 wt.% MoS <sub>2</sub> /AC	618	0.27	3.96
20 wt.% MoS <sub>2</sub> /AC	507	0.21	3.98

<sup>a</sup>BET surface area. <sup>b</sup>Average pore volume calculated by the BJH method from desorption branches

## SECTION S2 SUPPLYMENTARY TABLES

**Table S2** XPS quantitative analysis of the fresh and spent 8 wt.% MoS<sub>2</sub>/AC catalysts

Samples	Atomic (%) <sup>a</sup>			
	Mo 3d	S 2p	O 1s	C 1s
1st	1.82	0.97	11.28	85.93
2nd	1.75	0.63	11.96	85.65
3th	1.43	0.66	11.53	86.38
4th	0.8	0.36	13.62	85.66

<sup>a</sup>Calculated from XPS quantitative analysis