Supporting Information

Dual-Electrocatalysis Behavior of Star-like Zinc-Cobalt-Sulfide Decorated with Cobalt-Molybdenum-Phosphide in Hydrogen and Oxygen Evolution Reactions

Maryam Shamloofard^a, Saeed Shahrokhian^{a,b*}

^aDepartment of Chemistry, Sharif University of Technology, Tehran 11155-9516, Iran

^bInstitute for Nanoscience and Technology, Sharif University of Technology, Tehran, Iran



Figure S1. Energy-dispersive X-ray of the Zn-Co precursor.



Figure S2. Energy-dispersive X-ray of the Zn-Co-S.



Figure S3. Energy-dispersive X-ray of the Co-Mo-P (3)/Zn-Co-S.



Figure S4. Tafel plots for catalysts at the high overpotentials.



Figure S5. CV curves of catalysts in 1M KOH with different scan rates (5-100 mV s⁻¹)



Figure S6. The calculated TOF values for (A) OER, and (B) HER.



Figure S7. The multi-current step test over Co-Mo-P (3)/Zn-Co-S for OER and HER.



Figure S8. (A) The FE-SEM images, and (B) Energy-dispersive X-ray of the Co-Mo-P (3)/Zn-

Co-S after OER and HER tests.



Figure S9. The XRD patterns of the Co-Mo-P (3)/Zn-Co-S after OER and HER.



Figure S9. The XPS of the Co-Mo-P (3)/Zn-Co-S after OER test.



Figure S11. EIS Nyquist plots recorded for Co-Mo-P (3)/Zn-Co-S catalyst after (A) OER, and

(B) HER tests.

Table S1. Electrochemical parameters of the Zn-Co precursor, Zn-Co-A, Zn-Co-O, and Zn-Co-Scatalysts toward OER in 1 M KOH solution.

Catalysts	η at j = 10	η at j = 50	j at η = 0.34	
	mA cm ⁻² (mV)	mA cm ⁻² (mV)	V (mA cm ⁻²)	
Zn-Co precursor	377	-	3.95	
Zn-Co-O	394	450	1.04	
Zn-Co-A	372	413	2.9	
Zn-Co-S	321	355	22.79	

Table S2. Electrochemical parameters of the Zn-Co precursor, Zn-Co-A, Zn-Co-O, and Zn-Co-Scatalysts toward HER in 1 M KOH solution.

Catalysts	η at j = 10	η at j = 50	j at η = 0.2	
	mA cm ⁻² (mV)	mA cm ⁻² (mV)	V (mA cm ⁻²)	
Zn-Co precursor	478	550	2.29	
Zn-Co-O	414	467	1.12	
Zn-Co-A	303	383	2.56	
Zn-Co-S	176	242	17.32	

Table S3. Electrochemical parameters of the Co-P, Co-Mo-P (1), Co-Mo-P (2), Co-Mo-P (3), and Co-Mo-P (4) catalysts toward OER in 1 M KOH solution.

Catalysts	η at j = 10	η at j = 50	j at η = 0.34
	mA cm ⁻² (mV)	mA cm ⁻² (mV)	V (mA cm ⁻²)
Co-P	380	416	1.73
Со-Мо-Р (1)	378	411	1.85
Co-Mo-P (2)	357	398	5.05
Co-Mo-P (3)	348	381	6.82
Со-Мо-Р (4)	367	402	3.01

Table S4. Electrochemical parameters of the Co-P, Co-Mo-P (1), Co-Mo-P (2), Co-Mo-P (3), and Co-Mo-P (4) catalysts toward HER in 1 M KOH solution.

Catalysts	η at j = 10	η at j = 50	j at η = 0.2	
	mA cm ⁻² (mV)	mA cm ⁻² (mV)	V (mA cm ⁻²)	
Со-Р	334	425	2.57	
Со-Мо-Р (1)	309	362	2.70	
Co-Mo-P (2)	288	336	3.10	
Co-Mo-P (3)	171	268	14.85	
Co-Mo-P (4)	250	312	5.75	

Catalyst	Electrolyte	J (mA cm ⁻²)	η (mV)	Tafel slope	Ref.
				(mV dec ⁻¹)	
Co ₉ S ₈ /Zn-Co-S	1 M KOH	10	292	52	1
CoP(MoP)- CoMoO3@CN	1 M KOH	10	296	105	2
Co5MoxP NSs@NF	1 M KOH	10	270	54.4	3
CoMoPx	1 M KOH	10	305	50	4
Co-Mo-P (3)/Zn- Co-S	1 M KOH	10	273	41	This work

 Table S5A. Comparison of OER activity data with some representative non-noble catalysts.

Table S5B. Comparison of HER activity data with some representative non-noble catalysts.

Catalyst	Electrolyte	J (mA cm ⁻²)	η (mV)	Tafel slope	Ref.
				(mV dec ⁻¹)	
CoP(MoP)- CoMoO3@CN	1 M KOH	10	198	72	2
Co5MoxP NSs@NF	1 M KOH	10	173	218.6	3
СоРЗ/СоМоР	1 M KOH	10	125	61.1	5
ZnCo/NC@MoS ₂	1 M KOH	10	130	60.2	6
Co-Mo-P (3)/Zn- Co-S	1 М КОН	10	120	61.7	This work

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