Supporting information

3D Morphology of Bimodal Porous Copper with Nano-Sized and Micron-Sized Pores to Enhance Transport Properties for Functional Applications

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Further information on refinement analysis results

The synchrotron X-ray powder diffraction (XPD) refinement results of six different precursor alloys are shown in Figure S 1. The refinement results provide the information of phase composition, crystal structure, lattice parameter, phase density, weight percent and volume percent of each phase, as shown in Table S 1. Because the atomic radius of Cu and Fe are quite similar, they cannot be distinguished accurately in XPD beamline. When doing refinement, the definite crystal structures of CuFeAl phase are used as models and locally adjusting the atomic site occupancy of Cu, Fe and Al in each phase to make the total relative ratio of Cu, Fe and Al similar to that in initial precursor alloys. Therefore, the atomic site occupancy is only a reference. Moreover, although there is an unidentified phase in Al₃₀Fe₄₅Cu₂₅ alloy, the intensity of this unknown phase very low, so that it did not have

much effect on the other three phases during the refinement. Table S 1 showed that for the phases with the same crystal structure in various alloys, the atomic site occupancy is different.



Figure S 1 The synchrotron X-ray powder diffraction (XPD) refinement results of six different precursor alloys: (a) $Al_{10}Fe_{65}Cu_{25}$, (b) $Al_{20}Fe_{55}Cu_{25}$, (c) $Al_{30}Fe_{45}Cu_{25}$, (d) $Al_{40}Fe_{35}Cu_{25}$, (e) $Al_{50}Fe_{25}Cu_{25}$, (f) $Al_{60}Fe_{15}Cu_{25}$.

Alloy	Phase	Crystal structur e	Lattice parameter (Å)	Phase ratio (%)	Atomic site occupancy			_ R _{WP}
					Cu	Fe	Al	(%)
Al ₁₀ Fe ₆₅ Cu ₂₅	Cu-rich	Fm-3m	a=b=c=3.64	30.85	0.60	0.30	0.10	14.7
	Fe-rich	Im-3m	a=b=c=2.88	69.15	0.07	0.89	0.20	
Al ₂₀ Fe ₅₅ Cu ₂₅	Cu-rich	Fm-3m	a=b=c=3.67	32.88	0.50	0.30	0.20	13.4
	Fe-rich	Im-3m	a=b=c=2.90	67.12	0.13	0.67	0.20	
	Cu-rich	Fm-3m	a=b=c=3.67	4.92	0.58	0.16	0.26	
Al ₃₀ Fe ₄₅ Cu ₂₅	Fe-rich	Im-3m	a=b=c=2.91	50.69	0.15	0.75	0.10	10.0
	$Cu_{0.46}Fe_{0.82}Al_{0.72}$	Pm-3m	a=b=c=2.91	44.38	0.45	0.70	0.85	19.9
	unknown	/	/	/	/	/	/	
Al ₄₀ Fe ₃₅ Cu ₂₅	$Cu_{0.48}Fe_{0.72}Al_{0.8}$	Pm-3m	a=b=c=2.92	100	0.48	0.72	0.80	16.3
Al ₅₀ Fe ₂₅ Cu ₂₅	Cu _{0.5} Fe _{0.56} Al _{0.94}	Pm-3m	a=b=c=2.93	100	0.50	0.56	0.94	
	$Cu_{0.9}Fe_{0.2}Al_{0.9}$	Pm-3m	a=b=c=2.94	69.61	0.90	0.20	0.90	
Al ₆₀ Fe ₁₅ Cu ₂₅	CuFe ₄ Al ₁₂	C2/m	a=15.55, b=7.99, c=12.51	30.39	6	24	72	11.9

Table S 1 - Phases quantification of six different precursor alloys obtained from synchrotron X-ray

Composition of bimodal porous Cu after dealloying of Cu-Fe-Al alloys

The X-ray diffraction results of porous Cu after dealloying of six different Cu-Fe-Al precursor alloys are shown in Figure S 2. Only Cu phase diffraction peaks (with high intensity) and $Cu_{2+1}O$ phase diffraction peaks (with very low intensity) were detected in the XRD patterns of porous Cu. The results show that Fe and Al have been fully dissolved in the 5 wt% sulfuric acid (H₂SO₄) aqueous solution at 90 °C, and a small amount of Cu was oxidized during the dealloying process.



Figure S 2 The X-ray diffraction (XRD) results of porous Cu after dealloying of six different precursor alloys: Al₁₀Fe₆₅Cu₂₅, Al₂₀Fe₅₅Cu₂₅, Al₃₀Fe₄₅Cu₂₅, Al₄₀Fe₃₅Cu₂₅, Al₅₀Fe₂₅Cu₂₅, Al₆₀Fe₁₅Cu₂₅.



Average Size from the Pore Size Distribution by Gaussian Fitting

Figure S 3 – Gaussian fitting of pore size distribution

Table S 2 – Fitted results of average pore size and	d full width half maximum
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Broguesor Composition	Average Pore Size	Full Width Half Maximum		
Precuisor Composition	(Center Gravity) (nm)	(FWHM) (nm)		
Altere	1713	2629		
A1101 C65C u25	5616	3960		
	500	1269		
A130F C45C U25	2053	3230		
	74	49		
A1501 C25C U25	1005	2359		
AluFeuCura	70	2		
A1601 C15C U25	1221	2388		