Supporting Information

Particle Size Dependence of Proton Conduction in a Cationic Lanthanum Phosphonate MOF.

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Figure S1. Thermogravimetric Analysis of PCMOF21-AcO as Synthesized.



 $<38 \,\mu m$ $180 \le x < 125 \,\mu m$ $210 \,\mu m \le$

Figure S2. Post Impedance SEM Analysis of Particles from the Analyzed Distributions



Figure S3. Indexed Single Crystal of PCMOF21-AcO



Figure S4. Crystal structure of PCMOF21-AcO obtained at 100K.



Figure S5. Conductivity versus Temperature Comparison of Various Particle Sizes



Figure S6. Nyquist (Left, Second Cooling Cycle) and Arrhenius Plots (Right) for 'As Synthesized' Sample



Figure S7. Nyquist (Left, Second Cooling Cycle) and Arrhenius Plots (Right) for <38µm Particle Size



Figure S8. Nyquist (Left, Second Cooling Cycle) and Arrhenius Plots (Right) for 180>x≥125µm Particle Size



Figure S9. Nyquist (Left, Second Cooling Cycle) and Arrhenius Plots (Right) for \geq 210 μ m Particle Size



Figure S10. Extended Nyquist plots comparison of the particle sizes at various temperature ranges



Figure S11. 195K CO₂ Absorption data activated at 100°C overnight for PCMOF21-AcO



Figure S12. PXRD Pattern of PCMOF21-AcO Subject to Different Aqueous Conditions



Figure S13. Nyquist Plot Comparison of the 40 & 45 cN•m Variable Torque Experiment.

MOF	PCMOF21-AcO
Empirical formula	$C_{22}H_{46}La_2N_3O_{15}P_3$
FW (g mol ⁻¹)	963.35
Space group	P-1
λ (Å)	0.71073
Collection temp (K)	100
a (Å)	9.6056(16)
b (Å)	15.424(3)
c (Å)	16.780(3)
α (deg)	95.434(7)
β (deg)	106.266(7)
γ (deg)	91.765(9)
V (Å ³)	2371.5(7)
Crystal size (mm)	0.308 x 0.157 x 0.084
Z	2
ρ_{calc} (g cm ⁻³)	1.349
2θ range (deg)	5.174 to 56.886
Data/restraints/parameters	11739/0/410
No. reflns	32249
Ind refins $[I > 2\sigma(I)]$	11739
GOF, F ²	0.968
R factor $[I > 2\sigma(I)]$	$R_1 = 0.0494$, $wR_2 = 0.1156$
R factor (all data)	$R_1 = 0.0768$, $wR_2 = 0.1272$
CCDC ref code	

Table S1. Crystal Structure and Collection Parameters of PCMOF21-AcO