## **Supporting Information for**

## A Chiral Salen-based MOF Catalytic Material with highly thermal, aqueous and chemical stabilities

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Figure S2. IR spectrum of 1



Figure S3. UV-vis spectrum of L



Figure S4. <sup>1</sup>H NMR of L



Figure S5. <sup>13</sup>C NMR of L



Figure S6. HR-MS of L



**Figure S7.** (a) <sup>1</sup>H NMR of NiL in *d*-DMSO; (b) <sup>1</sup>H NMR of NiL in *d*-DCl and *d*-DMSO (V/V, 1:1); (c) <sup>1</sup>H NMR of **1** after dissolved in *d*-DCl and *d*-DMSO (V/V, 1:1).



 $Figure \ S8. \ A symmetric \ unit \ of \ 1$ 



Figure S9. Space-filling model of the 8-fold interpenetrated framework along *a* axis



Figure S10. TG curve of 1



<sup>1</sup>H NMR







<sup>1</sup>H NMR



<sup>13</sup>C NMR



<sup>1</sup>H NMR



<sup>13</sup>C NMR



<sup>1</sup>H NMR



Figure S11. NMR spectra of  $\beta$ -hydroxy-1,2,3-triazoles









Figure S12. GC-MS analyses of the cyclic carbonates



Figure S13. Proposed mechanism for the cycloaddition of  $\text{CO}_2$  with epoxides catalyzed by 1

## Table S1. BET plot data for 1

<b>BET Surface Area:</b>	$527.7582 \pm 3.3610 \ m^2/g$
Slope:	$0.008244 \pm 0.000052 \text{ g/cm}^3 \text{ STP}$
Y-Intercept:	$0.000003 \pm 0.000003 \text{ g/cm}^3 \text{ STP}$
C:	2873.98242
Qm:	121.252 cm <sup>3</sup> /g STP
Correlation Coefficient:	0.9999595
Molecular Cross-Sectional Area:	0.1620 nm <sup>2</sup>

<b>Relative Pressure</b>	elative Pressure Quantity Adsorbed		ative Pressure Quantity Adsorbed	
( <b>P</b> / <b>P</b> 0)	(cm <sup>3</sup> /g STP)			
0.020221646	120.6043	0.000171		
0.030745699	123.9053	0.000256		
0.055884706	128.4697	0.000461		
0.084060946	131.5677	0.000698		