

## **Electronic Supplementary Information for MS:**

### **Interpenetrating amine-functionalized metal-organic framework as an efficient and reusable catalyst for selective synthesis of tetrahydro-chromenes**

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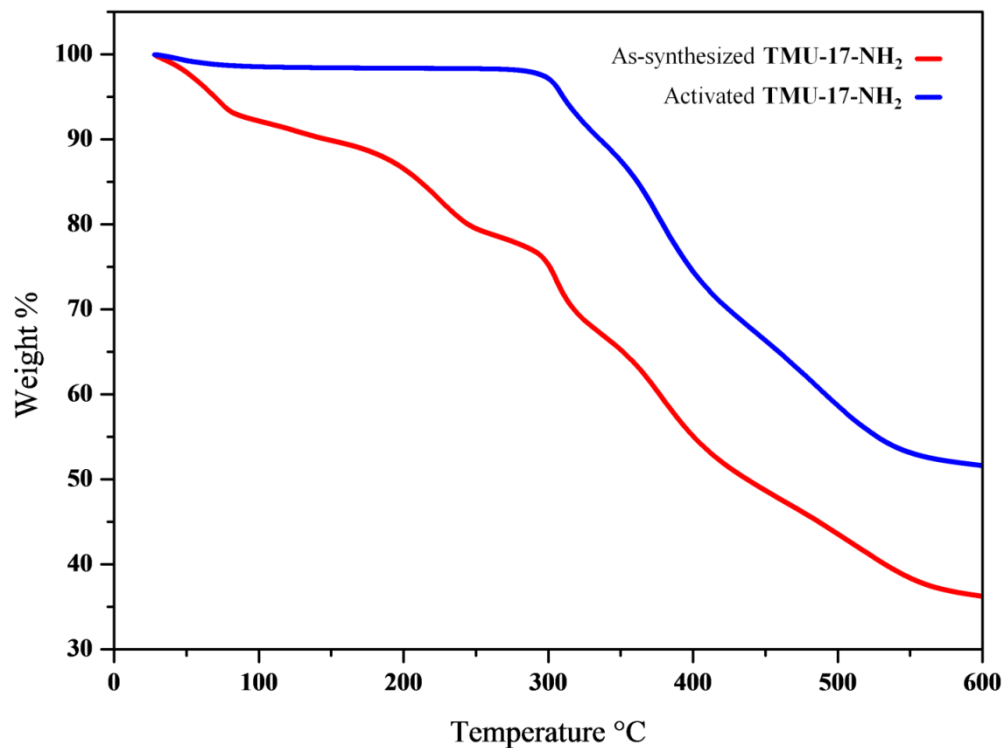


Fig. S1 TGA curves of as-synthesized (red) and activated (blue) TMU-17-NH<sub>2</sub>.

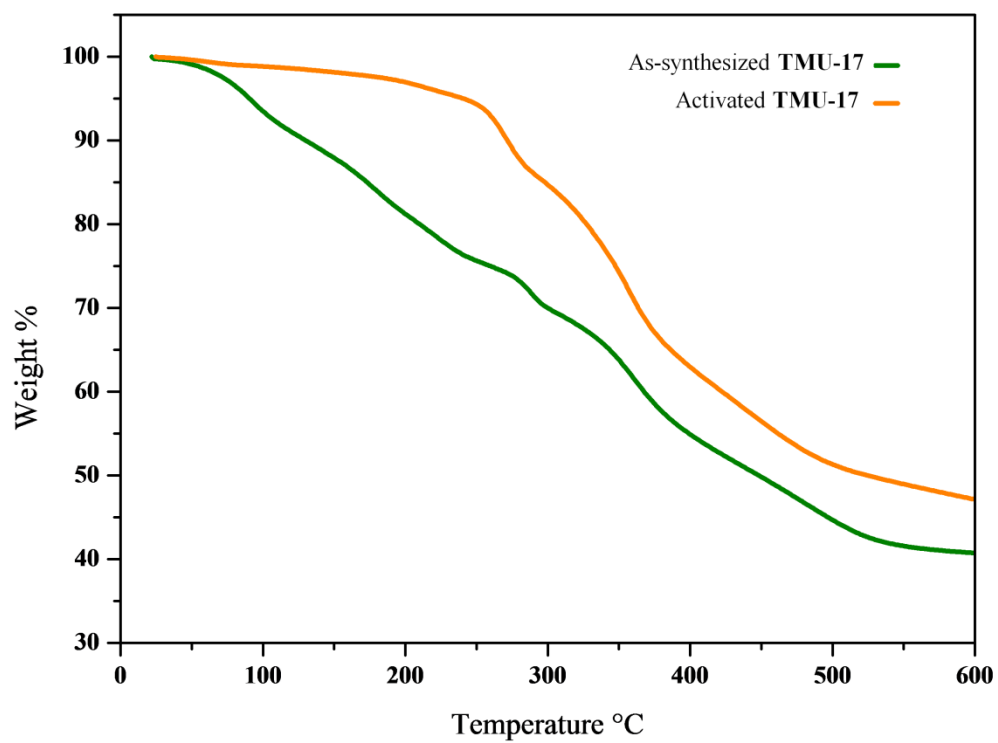
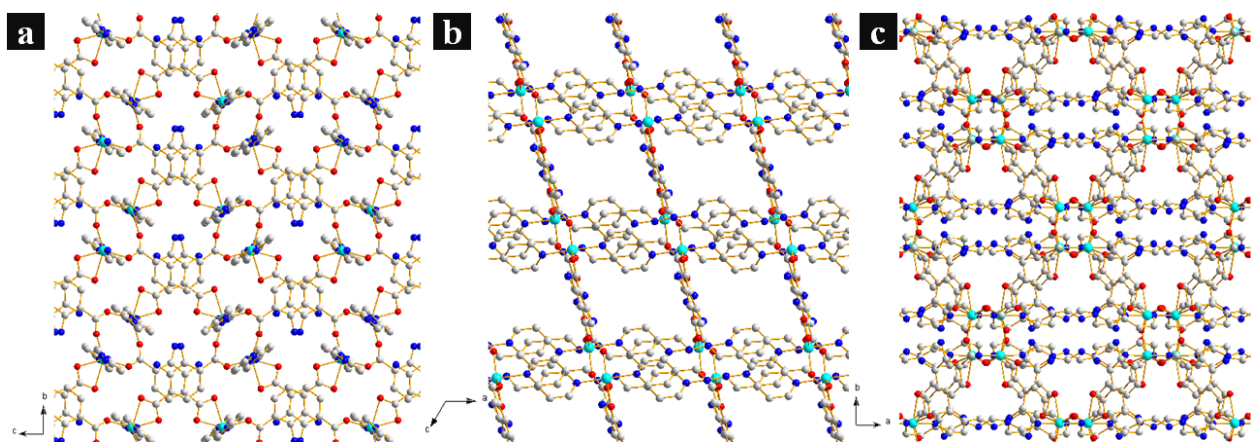
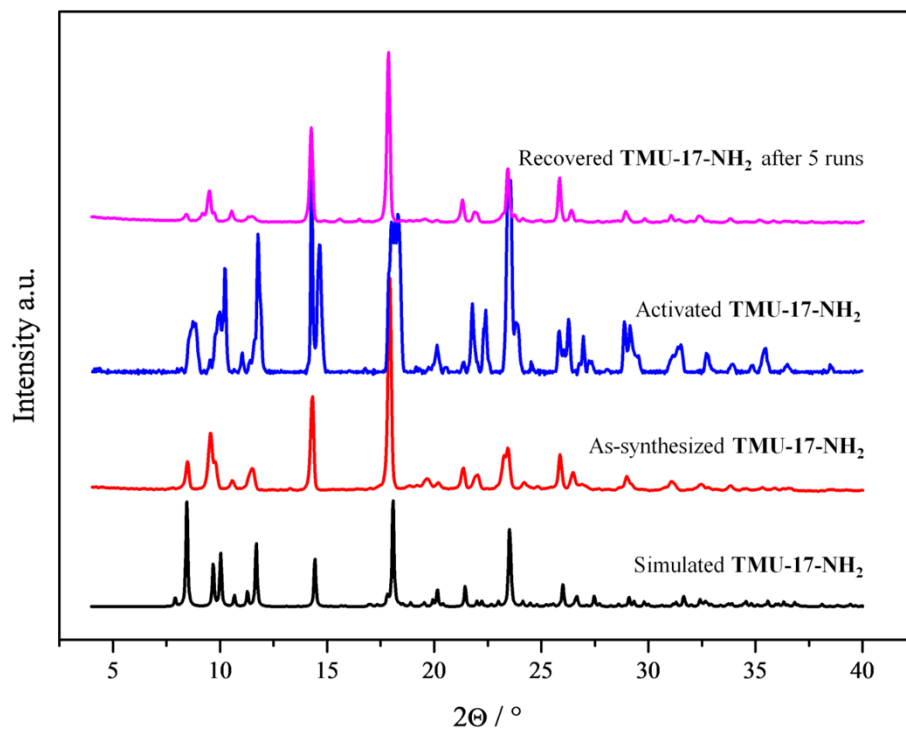


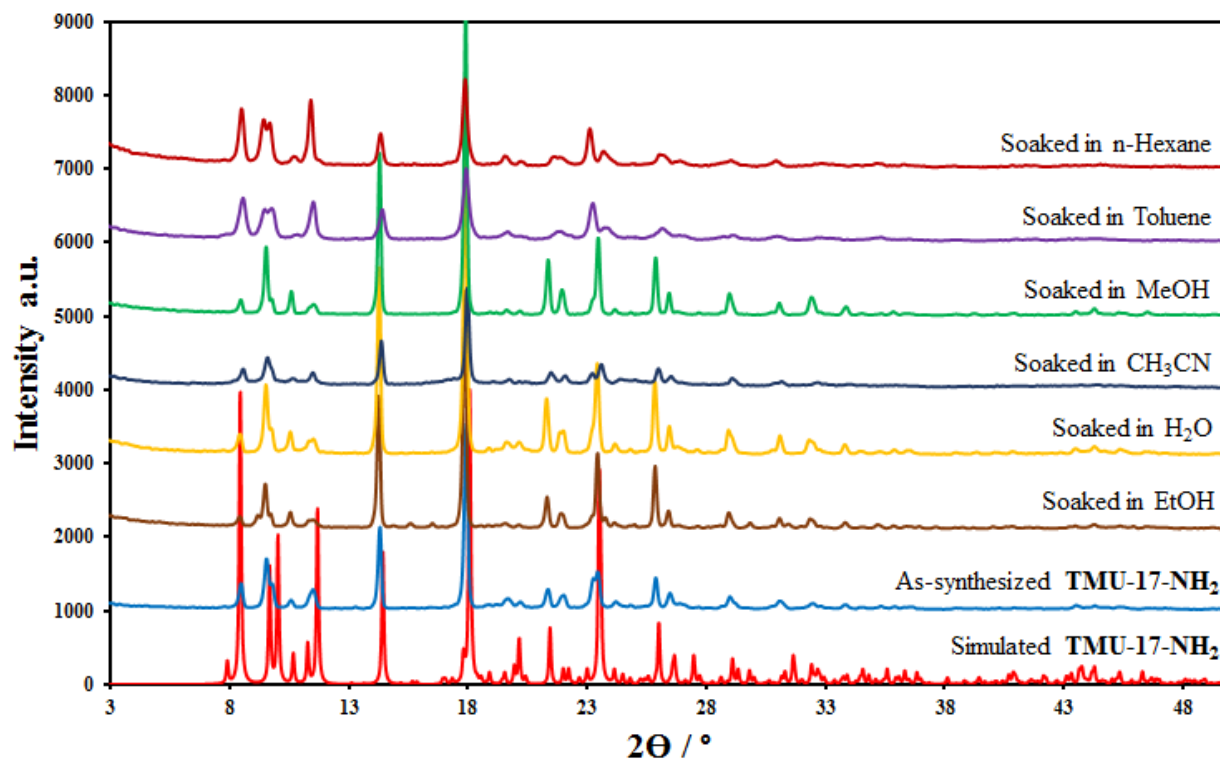
Fig. S2 TGA curve of as-synthesized (green) and activated (orange) TMU-17.



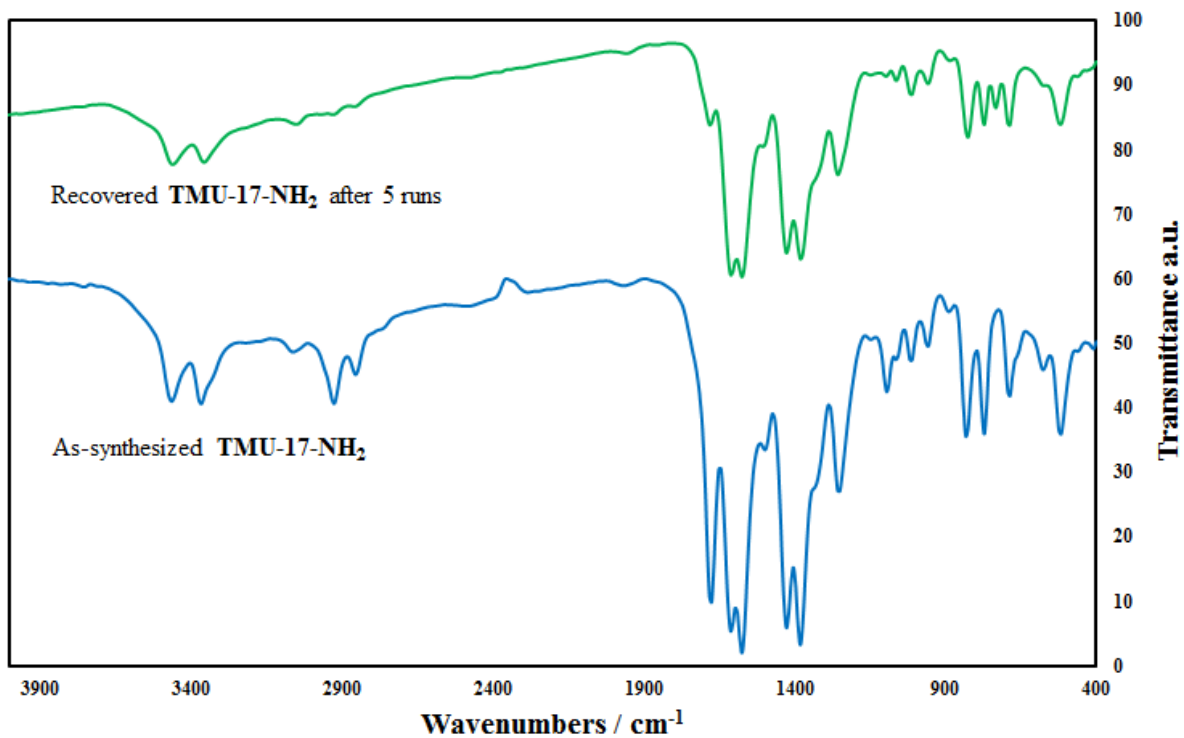
**Fig. S3** 3D two-fold interpenetrated **TMU-17-NH<sub>2</sub>**, viewed along (a) *a*, (b) *b* and (c) *c* axes. All hydrogen atoms and the disordered guest molecules in **TMU-17-NH<sub>2</sub>** are omitted for clarity.



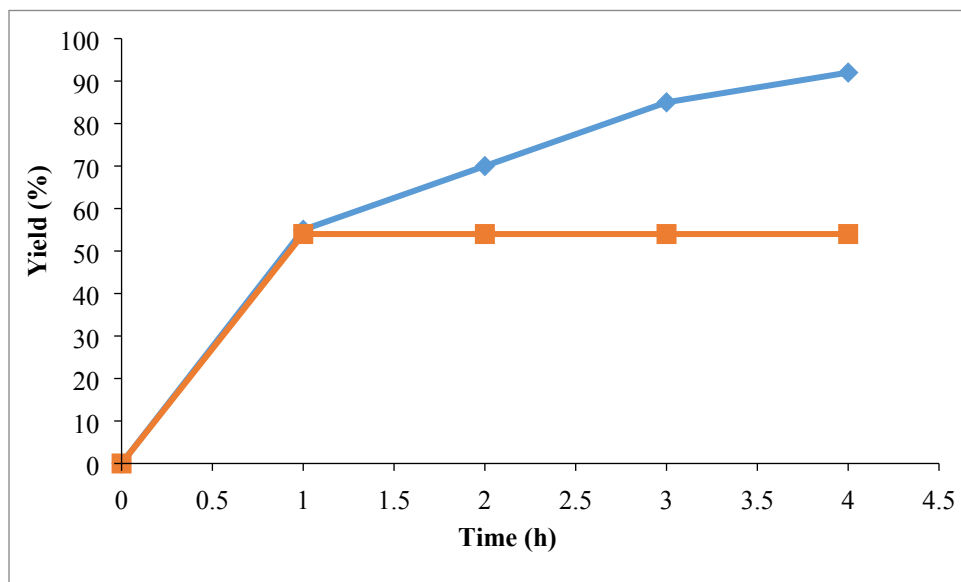
**Fig. S4** X-ray powder diffraction patterns of recovered TMU-17-NH<sub>2</sub> after five runs (purple); activated (blue); as-synthesized (red) and simulated (black) TMU-17-NH<sub>2</sub>.



**Fig. S5** PXRD patterns of simulated (red) and as-synthesized (blue) TMU-17-NH<sub>2</sub>; and immersed TMU-17-NH<sub>2</sub> in solvents for 2 hours.



**Fig. S6** FT-IR spectra of recovered TMU-17-NH<sub>2</sub> after five runs (green) and as-synthesized TMU-17-NH<sub>2</sub> (blue).

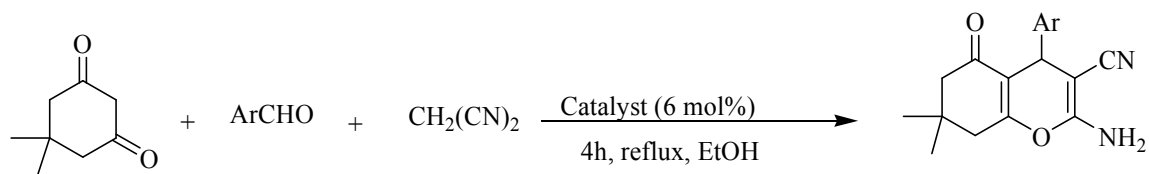


**Fig. S7.** Leaching test for TMU-17-NH<sub>2</sub>.

**Table S1** Preliminary crystallographic data for **TMU-17** in comparison with those of **TMU-17-NH<sub>2</sub>**.

Entry	MOF	Crystal System	<i>a</i> (Å)	<i>b</i> (Å)	<i>c</i> (Å)	β (deg)
<b>1</b>	TMU-17-NH <sub>2</sub>	monoclinic	15.655	16.561	21.200	120.567
<b>2</b>	TMU-17	monoclinic	15.939	16.354	21.511	122.001

**Table S2.** The reaction in the presence of **TMU-17-NH<sub>2</sub>** and **TMU-17**.



Entry	Catalyst	Yield (%)
<b>1</b>	TMU-17	43
<b>2</b>	TMU-17-NH <sub>2</sub>	92
<b>3</b>	-	0

**Table S3** Crystal data and structure refinements for compound **TMU-17-NH<sub>2</sub>**.

Identification code	TMU-17-NH <sub>2</sub>
Chemical formula	C <sub>23</sub> H <sub>22</sub> N <sub>6</sub> O <sub>5</sub> Zn
formula mass	527.84
<i>T</i> (K)	180(2)
Crystal syst	Monoclinic
Space group	<i>C</i> <sub>2</sub> / <i>c</i>
<i>a</i> (Å)	15.655(3)
<i>b</i> (Å)	16.561(3)
<i>c</i> (Å)	21.200(3)
β (deg)	120.567(9)
<i>V</i> (Å <sup>3</sup> )	4732.6(14)
<i>Z</i>	8
μ <sub>calcd.</sub> (g/m <sup>3</sup> )	1.085
F(000)	2176
Reflections collected	4534 / 2844
R(int)	0.0855
Goodness-of-fit on <i>F</i> <sup>2</sup>	1.083
R <sub>1</sub> <sup>a</sup> [ <i>I</i> > 2σ( <i>I</i> )]	0.0805
wR <sub>2</sub> <sup>b</sup>	0.1589

$${}^a R_1 = \frac{\sum ||F_o| - |F_c||}{\sum |F_o|}, \quad {}^b wR_2 = \frac{|\sum w(|F_o|^2 - |F_c|^2)|}{\sum w(F_o^2)^2}^{1/2}$$