Construction of structural diversity and fine-tuned porosity in acylamide

MOFs by synthetical approach

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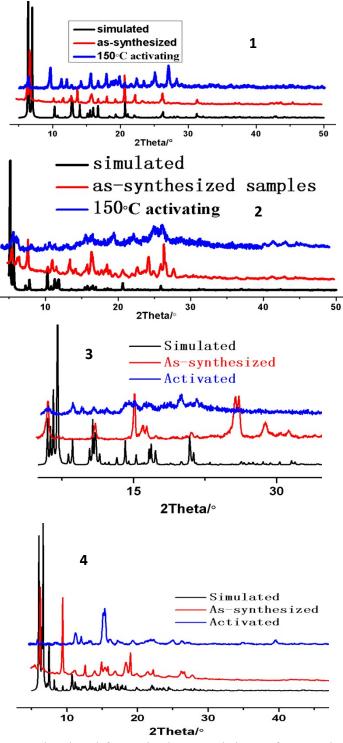


Fig. S1 The PXRD patterns simulated from single crystal data, of as-synthesized bulk samples and

after activating at 150°C for polymers1-4, respectively.

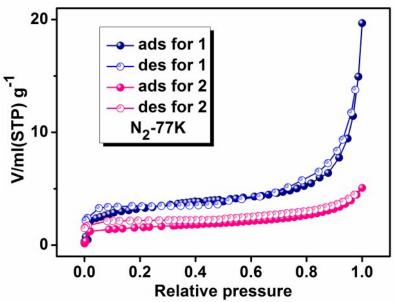


Fig. S2 The N_2 adsorption isotherms of 1 and 2 at 77K.