

Supporting Information

Synthesis of microporous organic polymers with high CO₂-over-N₂ selectivity and CO₂ adsorption

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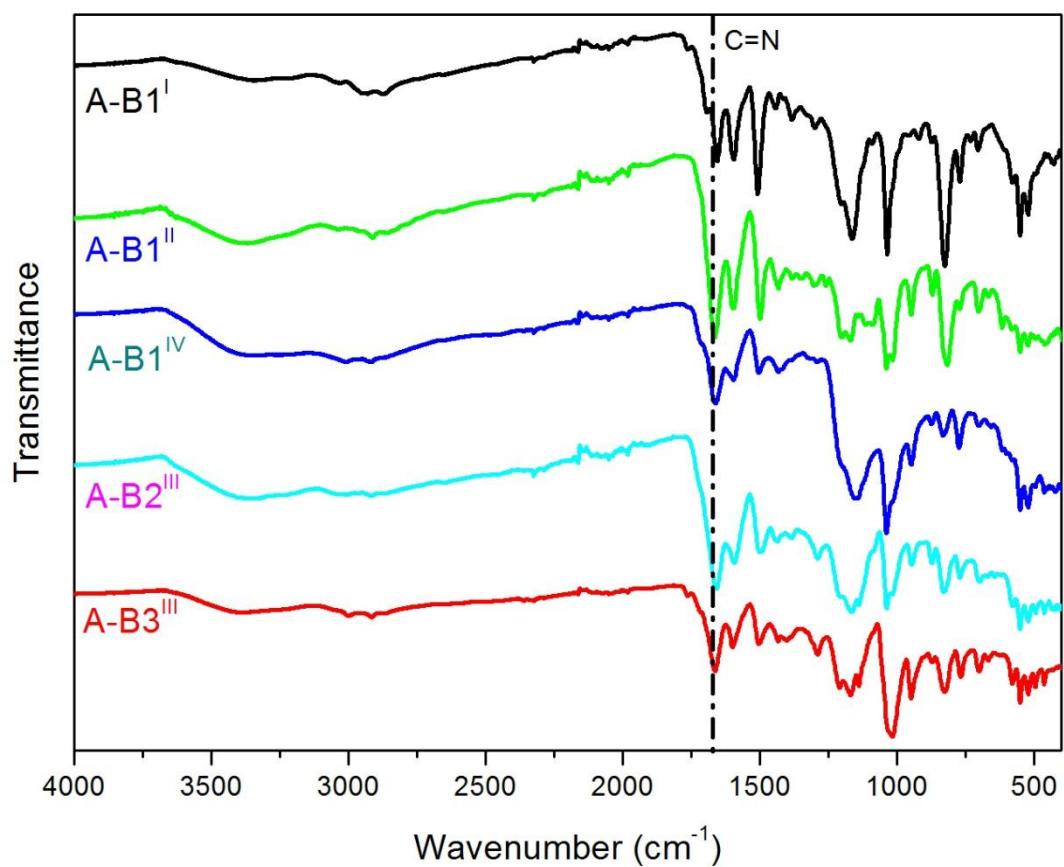


Fig. 1S Fourier transform infrared (FT-IR) spectra of as-synthesized MOPs.

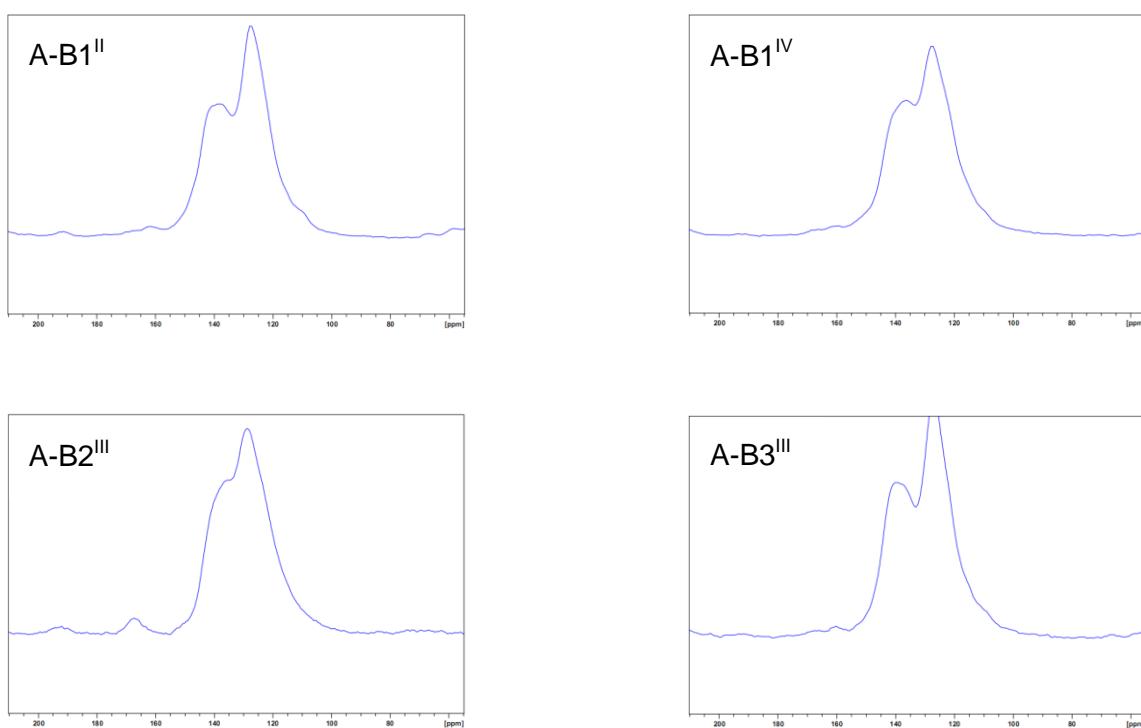


Fig. 2S Solid state ^{13}C CP/MAS NMR spectrum of MOPs after heated at 300°C

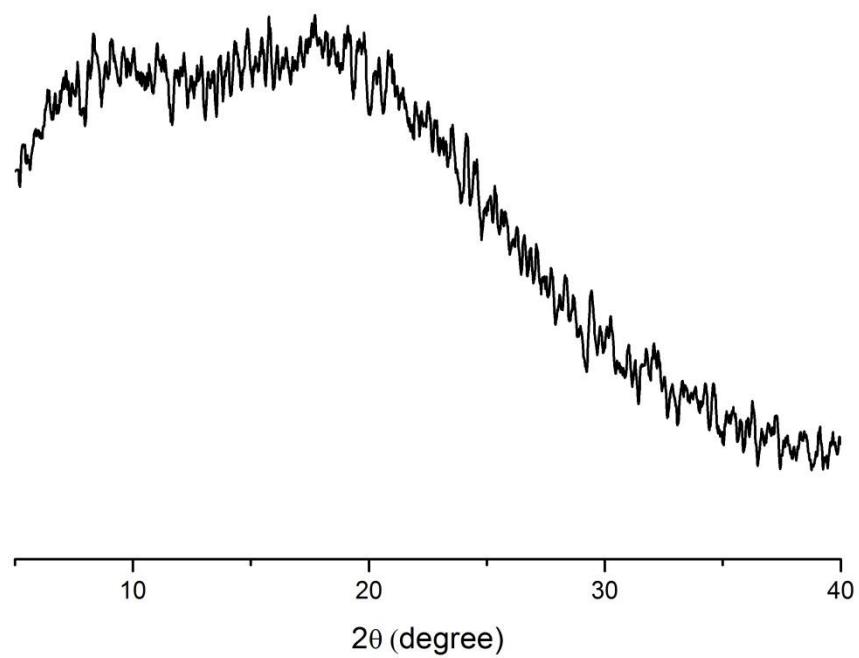


Fig. 3S Powder X-ray diffraction (PXRD) pattern of **A-B2^{III}**.

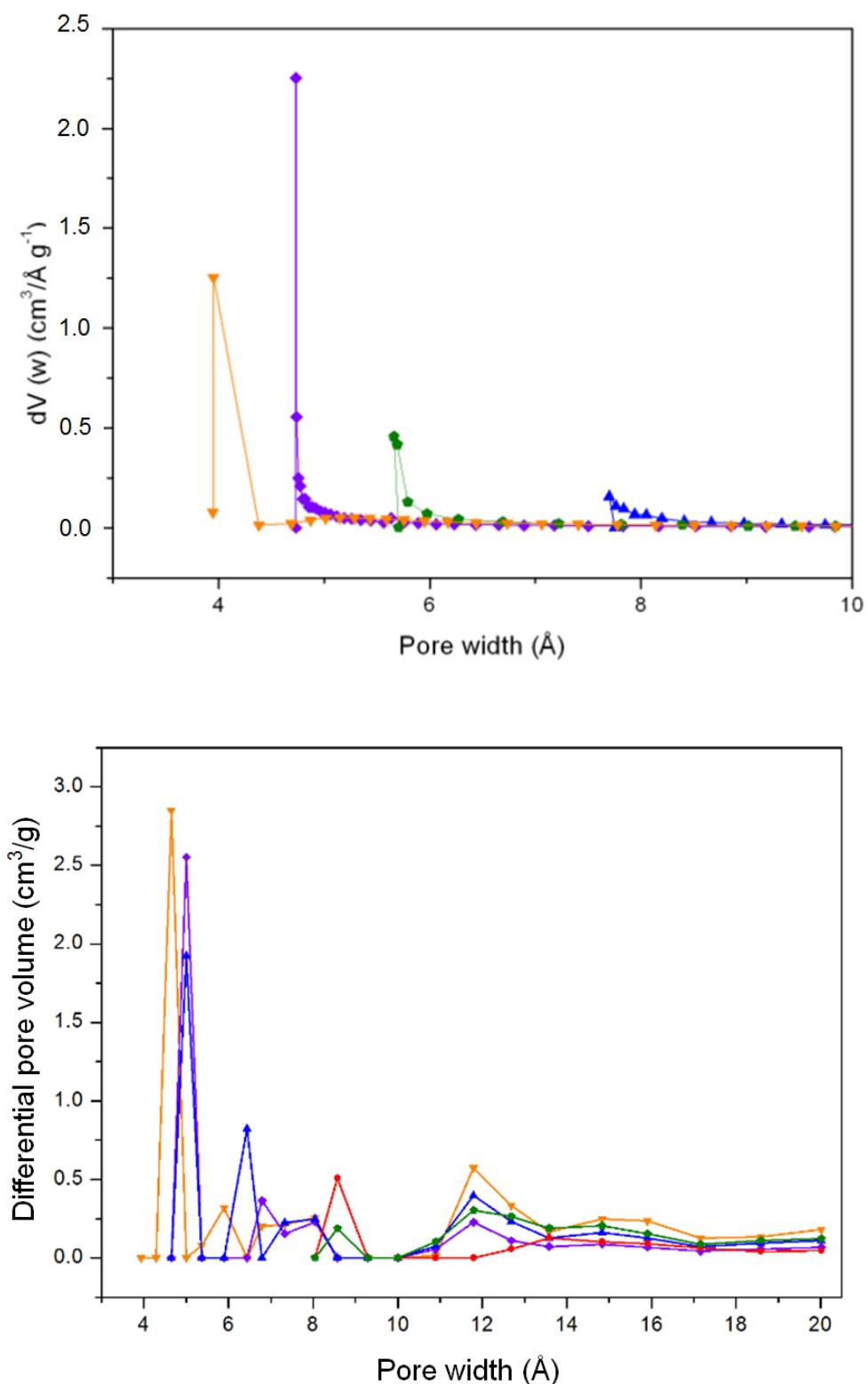


Fig. 4S Pore size distributions of microporous organic polymers by using N_2 adsorption isotherms at 77 K with Horvath-Kawazoe (top) and Density Functional Theory (bottom) method. **A-B1^{II}**, red line with circle; **A-B1^{III}**, violet line with diamond; **A-B1^{IV}**, blue line with triangle; **A-B2^{III}**, olive line with pentagon; **A-B3^{III}**, orange line with reverse triangle.

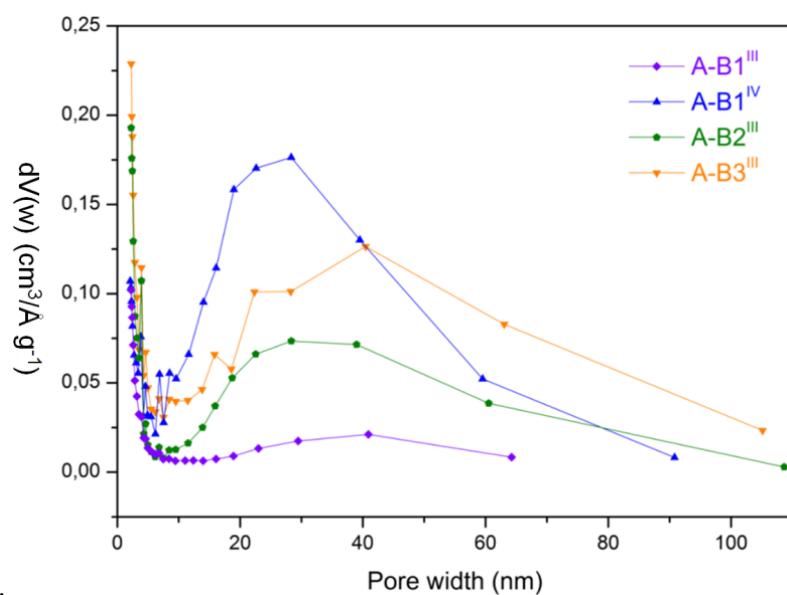


Fig. 5S Pore size distribution of microporous organic polymers (MOPs) by using N₂ desorption isotherms with Barrett–Joyner–Halenda (BJH) method.

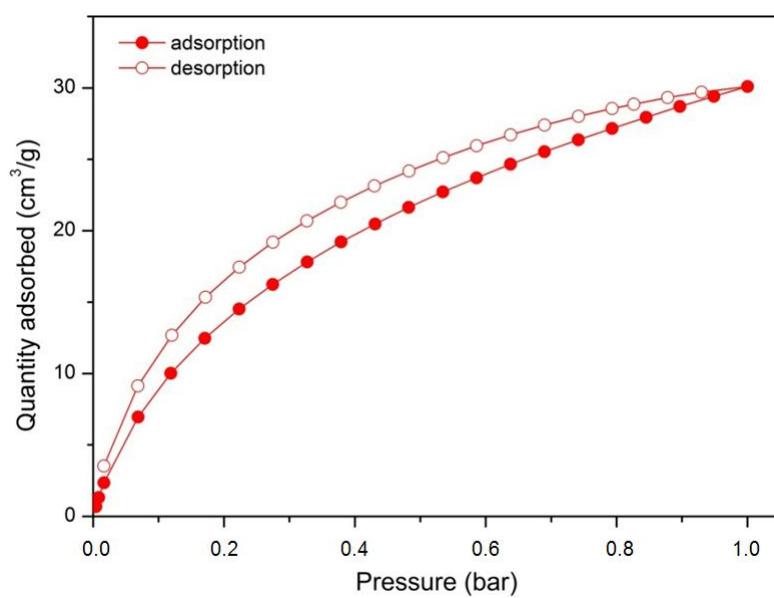


Fig. 6S CO₂ adsorption and desorption isotherm of A-B1^I at 273 K

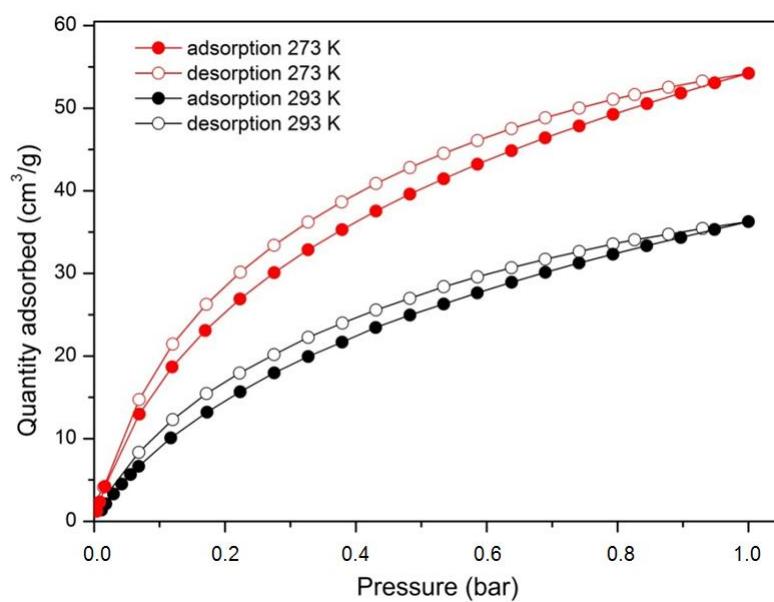


Fig. 7S CO₂ adsorption and desorption isotherm of A-B1^{II} at 273 K and 293 K.

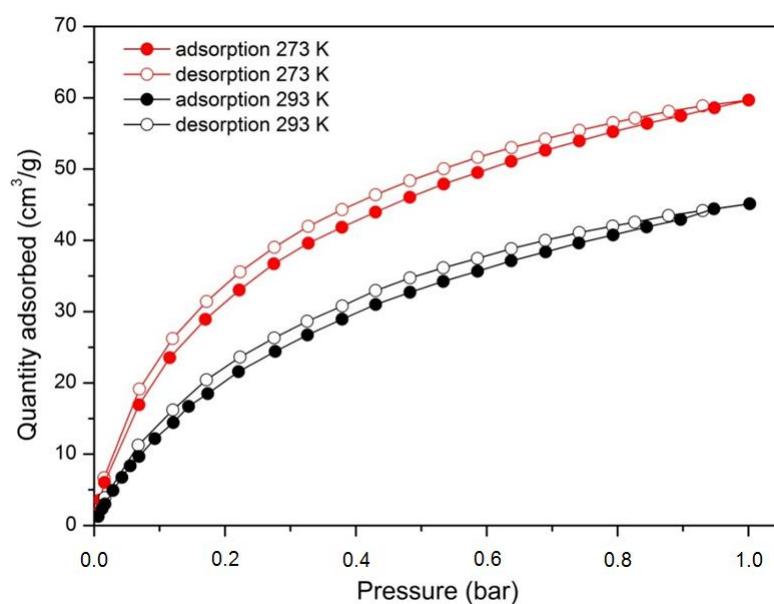


Fig. 8S CO₂ adsorption and desorption isotherm of A-B1^{III} at 273 K and 293 K.

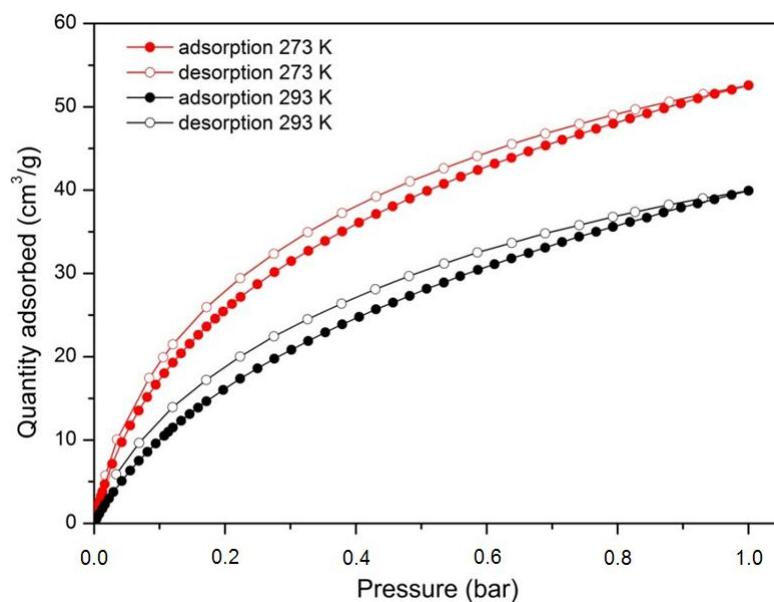


Fig. 9S CO₂ adsorption and desorption isotherm of A-B1^{IV} at 273 K and 293 K.

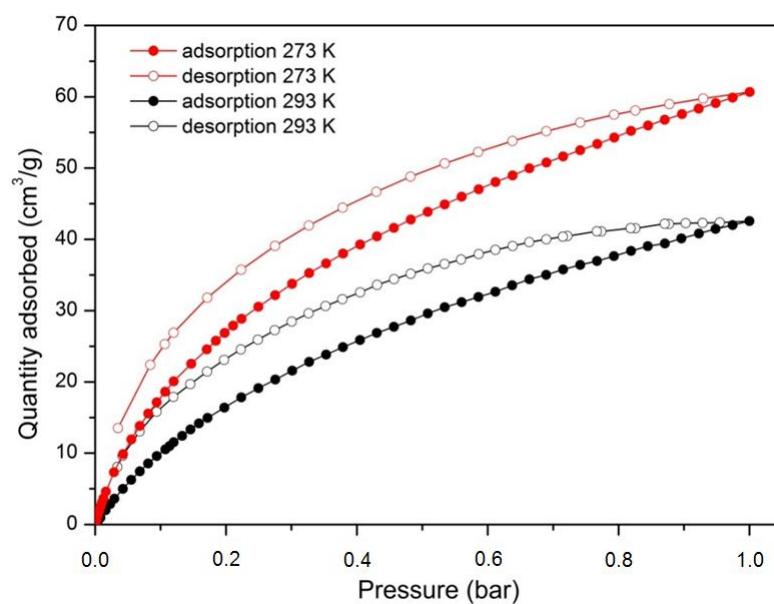


Fig. 10S CO₂ adsorption and desorption isotherm of A-B2^{III} at 273 K and 293 K.

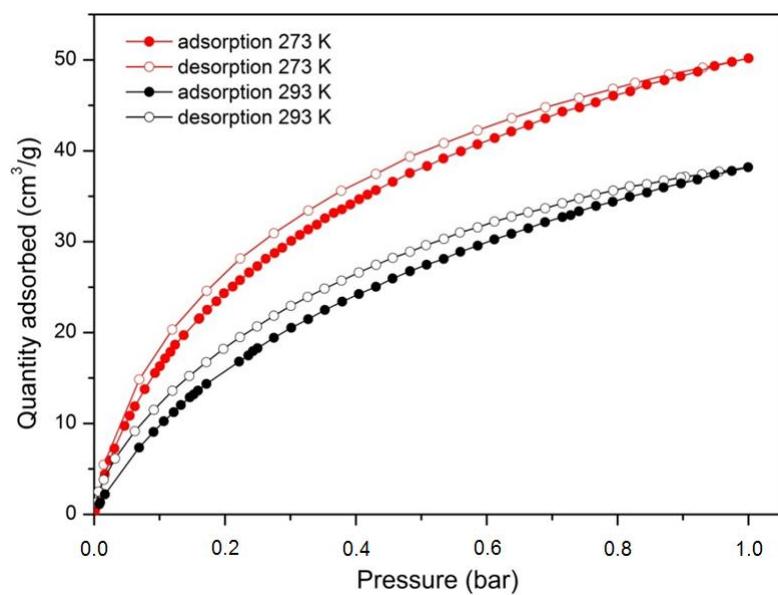


Fig. 11S CO₂ adsorption and desorption isotherm of A-B3^{III} at 273 K and 293 K.

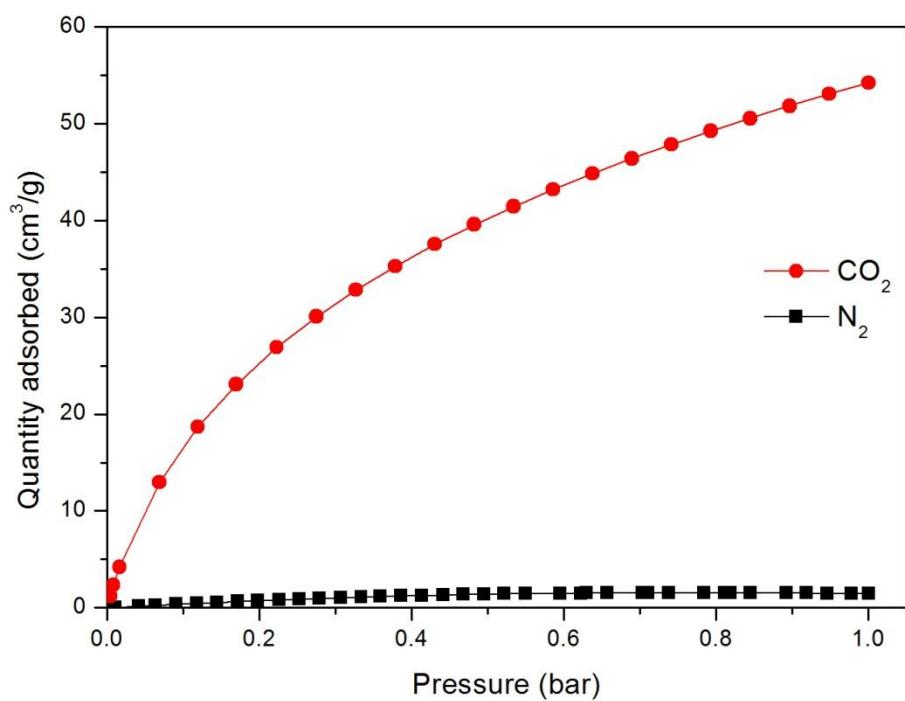


Fig. 12S CO₂ and N₂ adsorption capacity for A-B1^{II} at 273 K.

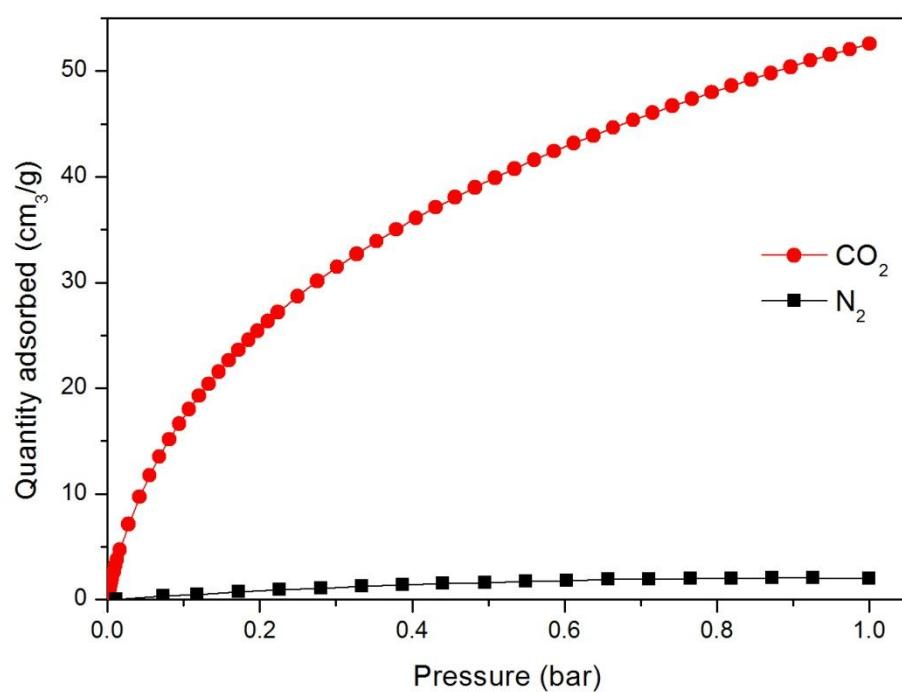


Fig. 13S CO₂ and N₂ adsorption capacity for **A-B1^{IV}** at 273 K.

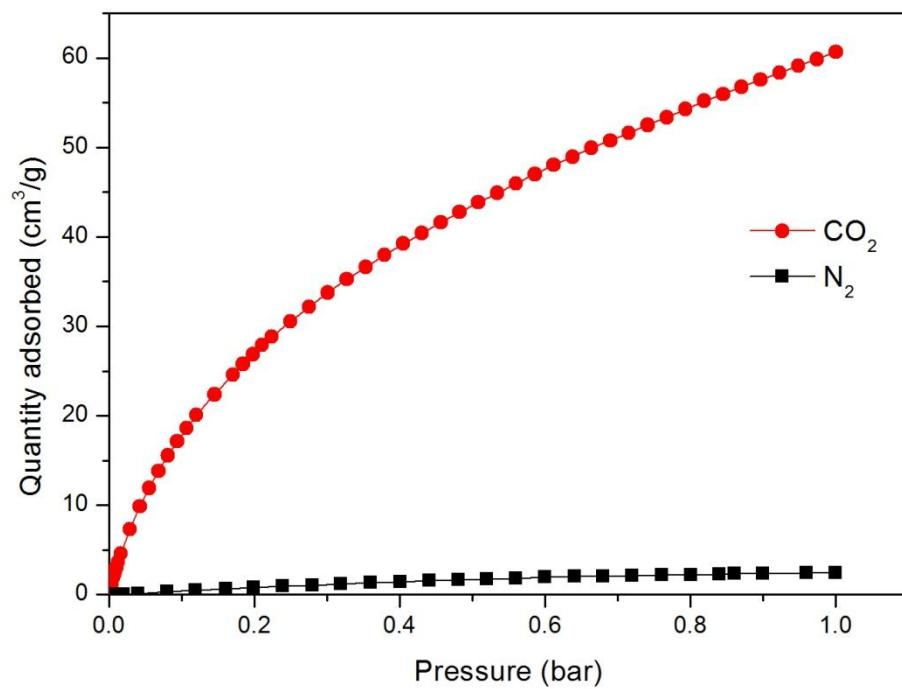


Fig. 14S CO₂ and N₂ adsorption capacity for **A-B2^{III}** at 273 K.

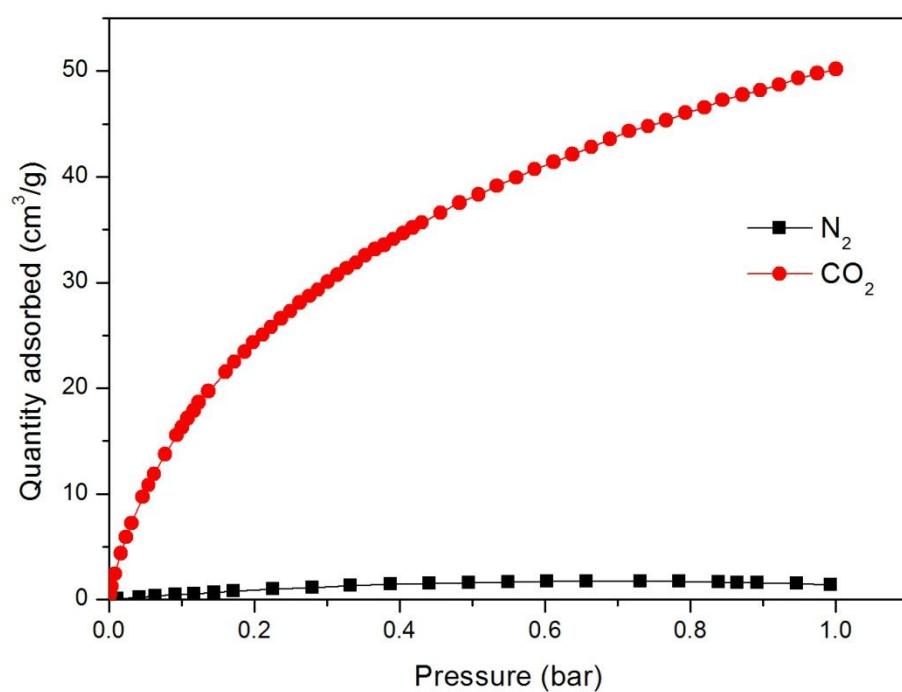


Fig. 15S CO₂ and N₂ adsorption capacity for **A-B3^{III}** at 273 K.

Table 1S. Summary of S_{BET} and CO_2 uptake for nitrogen containing microporous organic polymers/cages.

Material	S_{BET} (m^2/g)	CO ₂ uptake at			T (K)	Ref
		0.15 bar	1.0 bar			
A-B1 ^I	68	0.51	1.34		273	<i>a</i>
A-B1 ^{II}	124	0.95	2.42		273	<i>a</i>
		0.53	1.62		293	
A-B1 ^{III}	377	1.20	2.67		273	<i>a</i>
		0.76	2.01		293	
A-B1 ^{IV}	452	0.98	2.35		273	<i>a</i>
		0.60	1.78		293	
A-B2	614	1.02	2.71		273	<i>a</i>
		0.59	1.90		293	
A-B3	592	0.93	2.24		273	<i>a</i>
		0.58	1.71		293	
POF A1-B4	129	0.83	1.96		273	50
POF A1-B5	63	0.65	1.65		273	50
POF A1-B6	50	0.54	1.36		273	50
CMP-1-NH ₂	710	0.60	1.64		273	46
		0.22	0.95		298	
PI-1	506	0.87	2.10		273	51
		0.47	1.41		298	
PI-2	568	0.45	1.60		273	51
		0.24	1.00		298	
TBI-1	609	1.08	3.17		273	35
TBI-2	582	1.02	2.68		273	35
BILP-1	1172	1.63	4.27		273	62
		0.91	2.98		298	
BILP-2	708	1.45	3.39		273	52
		0.78	2.36		298	

Material	S_{BET} (m^2/g)	CO ₂ uptake at				Ref
		0.15 bar	1.0 bar	T (K)		
BILP-3	1306	1.67	5.11	273	60	
		0.82	3.30	298		
BILP-4	1135	1.99	5.34	273	52	
		1.01	3.59	298		
BILP-5	599	1.12	2.90	273	52	
		0.57	1.98	298		
BILP-6	1261	1.81	4.80	273	60	
		0.98	2.75	298		
BILP-7	1122	1.47	4.39	273	52	
		0.73	2.77	298		
PPN-6-CH ₂ -DETA	555	3.04	4.30	295	59	
PPN-6-CH ₂ -TAEA	663	2.35	3.54	295	59	
PPN-6-CH ₂ -TETA	634	2.04	3.25	295	59	
PPN-6-CH ₂ -EDA	1014	1.82	3.26	295	59	
CC1	23	0.56	1.27	275	65	
		0.40	1.03	289		
CC2	533	1.02	3.00	275	65	
		0.72	2.21	289		
CC3	624	0.90	2.47	275	65	
		0.61	1.94	289		
CC4	930	0.59	1.90	275	66	
CC5	1333	0.91	3.07	275	67	
CC6	99	0.20	0.90	273	68	
Cage3	1377	0.45	2.10	273	69	
Cage3a	744	0.52	2.70	298	70	

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