## Supplementary Information

## High and Selective CO<sub>2</sub> adsorption by a phthalocyanine nanoporous polymer

Venkata S. Pavan K. Neti,<sup>a</sup> Jun Wang,<sup>b</sup> Shuguang Deng,<sup>b</sup> and Luis Echegoyen<sup>\*a</sup>

<sup>a</sup> Department of Chemistry, University of Texas at El Paso, El Paso, TX 79968, United States <sup>b</sup> Department of Chemical Engineering, New Mexico State University, Las Cruces, NM, United States

<sup>a</sup> E-mail: echegoyen@utep.edu, Tel/Fax: +1 (915) 747-7573/(915) 747-8807

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## Langmuir model fits



Figure S1. Langmuir model fits for CO<sub>2</sub> (top), CH<sub>4</sub> (middle), and N<sub>2</sub> adsorption (bottom) of CPP at 273K.

Henry's constant by the product of Langmuir constants, that is K=a\*b. K1 (273K) and K2 (298K), ln K vs 1/T (below). Van't Hoff equation is used to get  $Q_{st}$  at zero coverage.



Figure S2. Van't Hoff plots of isosteric heat of adsorption for CH<sub>4</sub> (top) and CO<sub>2</sub> (bottom).

## Calculation of isosteric heat of adsorption

The adsorption enthalpy at zero coverage was calculated from Henry's constant using the Van't Hoff equation as

$$\ln K = -\frac{\Delta H}{RT} + \frac{\Delta S}{R}$$

K is the Henry's constant, T is the temperature, plotting *ln K* vs. 1000/T

**Surface Area Measurements** 



Figure S3. Cumulative (left) pore size distribution plot of CPP from the application of the NLDFT model to the  $N_2$  isotherm. BET plot (right) for CPP calculated from isotherm data.