

## **Supporting Information**

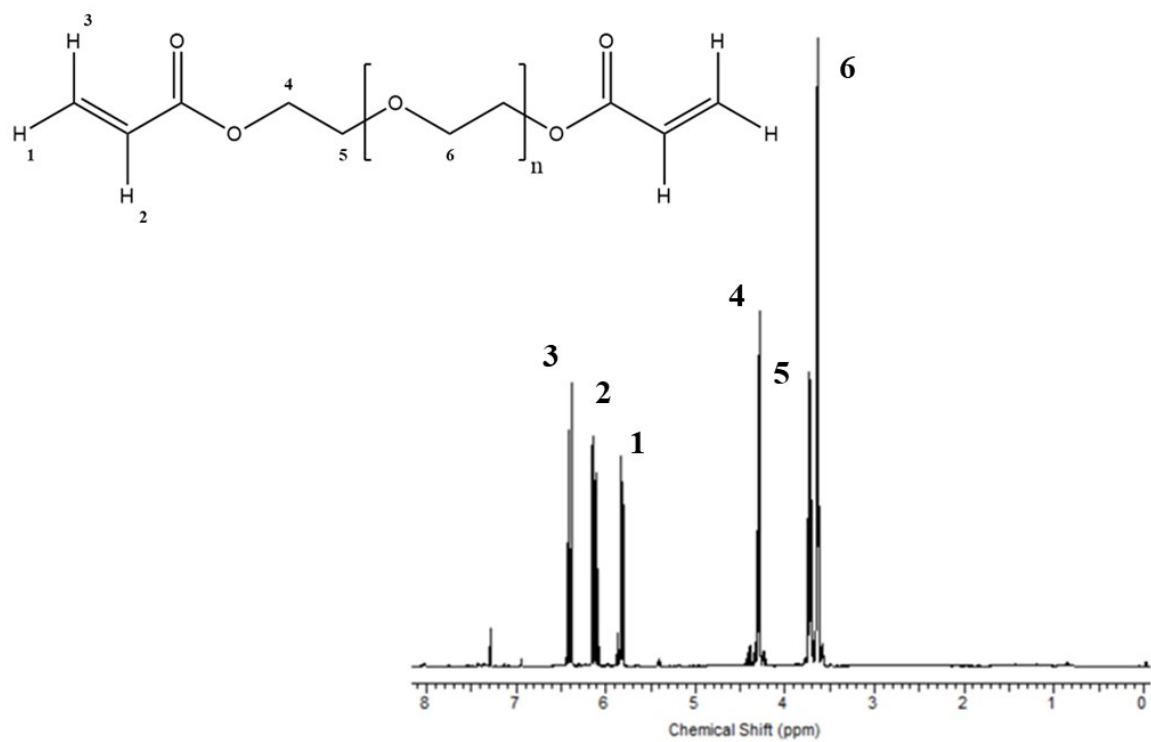
# Tetraalkylammonium based Dicationic Ionic Liquid (IL) for CO<sub>2</sub> Capture

Prashant S. Kulkarni<sup>1\*</sup>, Prathamesh Ranjane<sup>1</sup>, Karun Mishra<sup>1</sup>, Swati Sundararajan<sup>1</sup> and Sanjay Kamble<sup>2</sup>

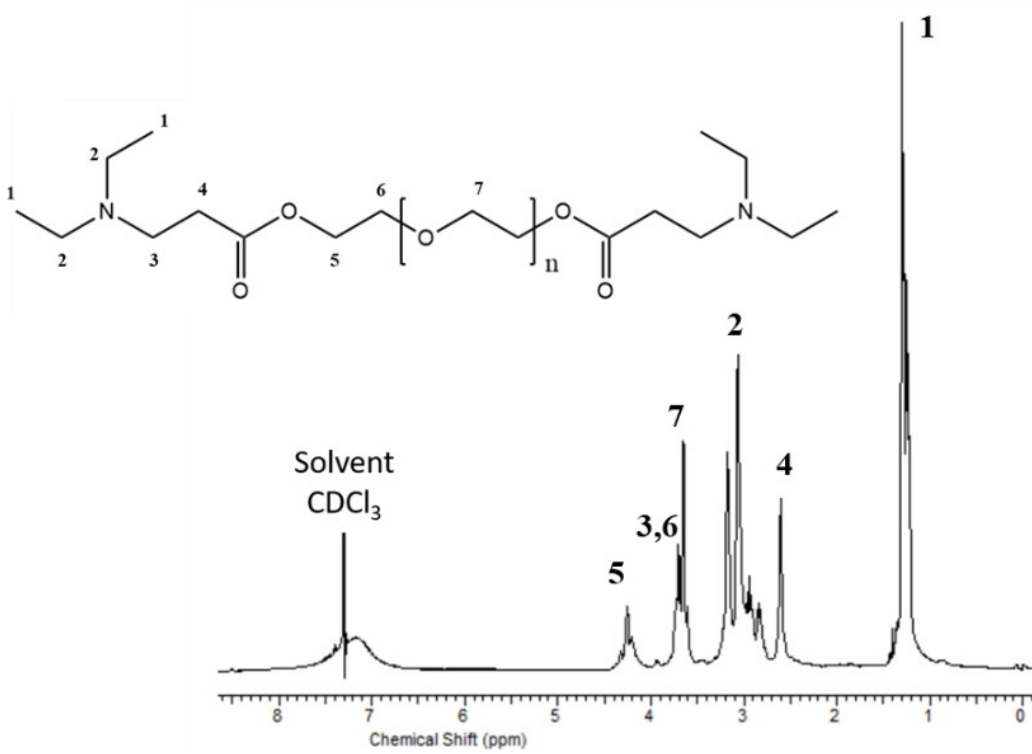
<sup>1</sup>Energy and Environment Laboratory, Department of Applied Chemistry, Defence Institute of Advanced Technology, Deemed University, Girinagar, Pune- 411 025, India

<sup>2</sup>Chemical Engineering and Process Development Laboratory, National Chemical Laboratory, Dr. Homi Bhabha Road, Pune 411008, India

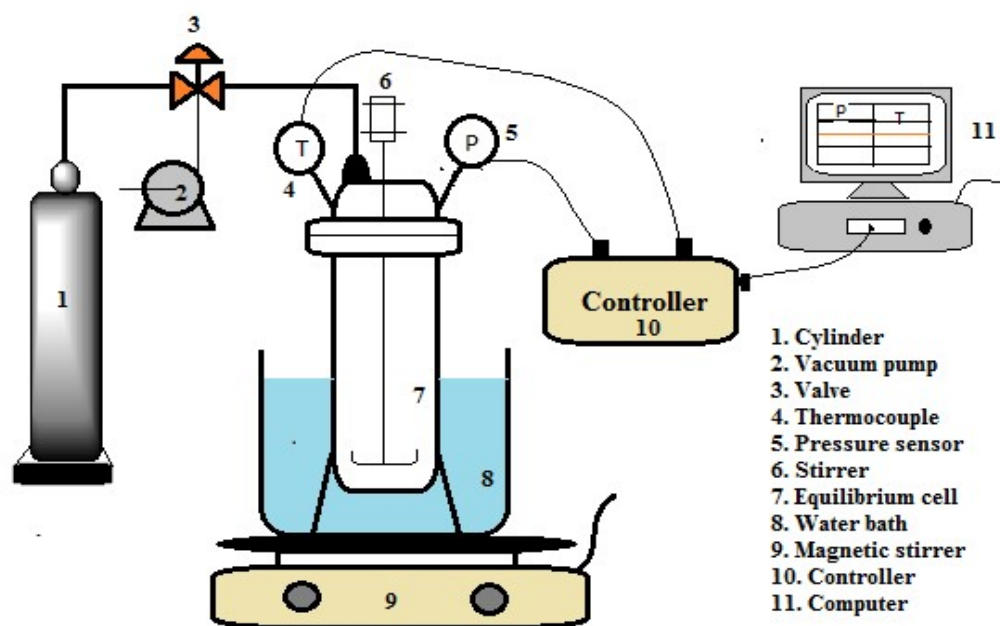
\* **Corresponding author:** Dr. Prashant S. Kulkarni, Phone: +91-02024604462,  
e-mail: [ps\\_kulkarni@rediffmail.com](mailto:ps_kulkarni@rediffmail.com) & [pskulkarni@diat.ac.in](mailto:pskulkarni@diat.ac.in)



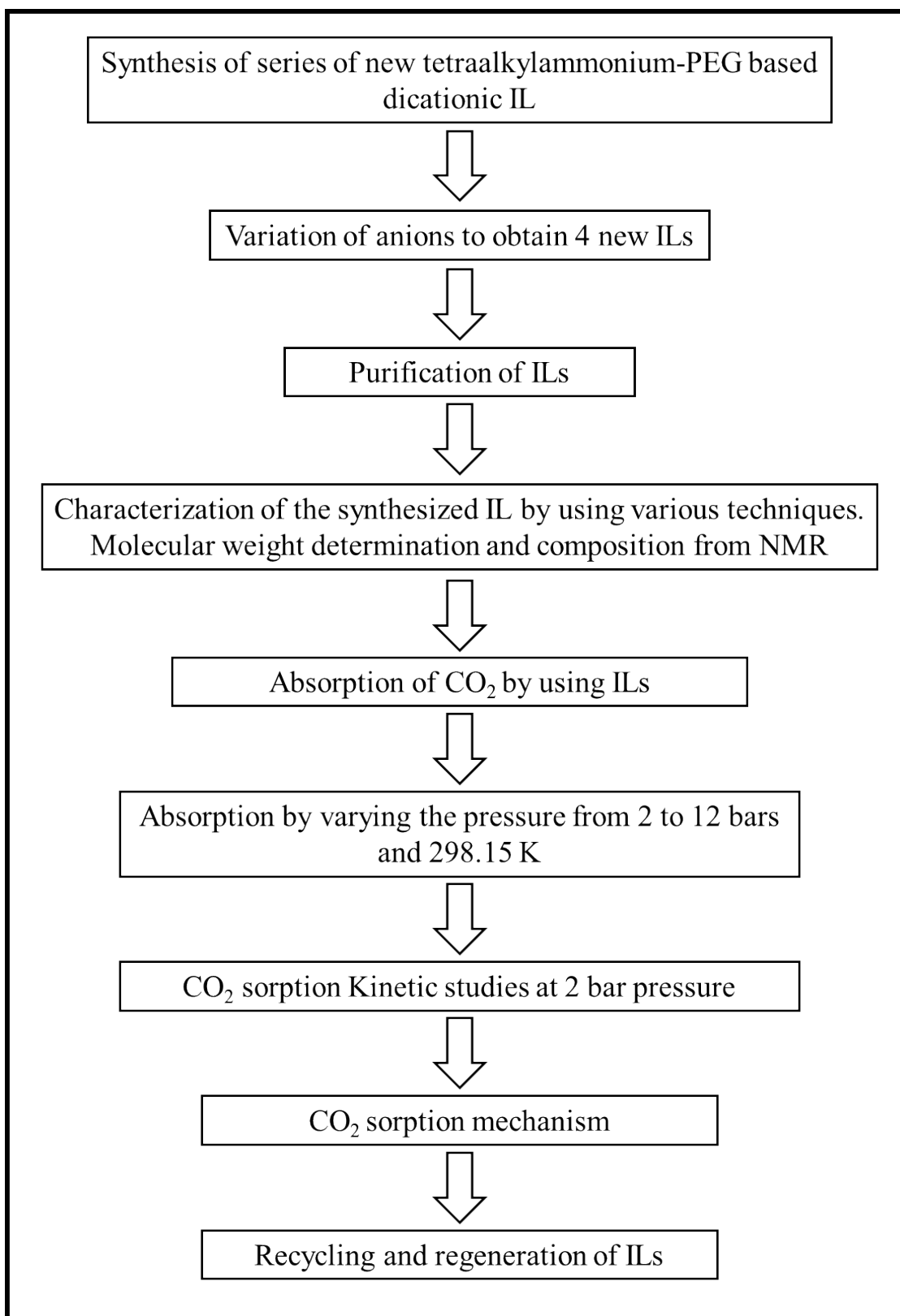
**Figure S1:** <sup>1</sup>H NMR spectra of PEGDA.



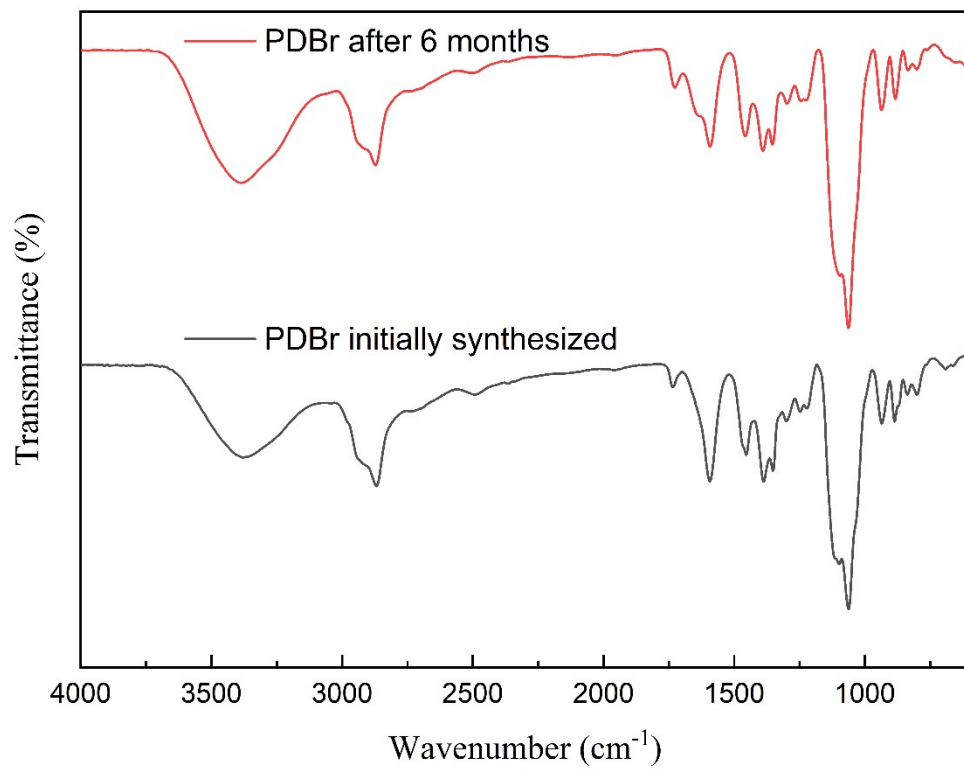
**Figure S2:** <sup>1</sup>H NMR spectra of PEGDA-DEA product (PD).



**Figure S3:** Schematic representation of fabricated CO<sub>2</sub> absorption apparatus.



**Figure S4:** Flowchart - absorption of CO<sub>2</sub> by using tetraalkylammonium dicationic IL.



**Figure S5:** Chemical stability of PDBr IL after 6 months, as investigated by FTIR.

**Table S1:** Design of experiment with input parameters and operating conditions.

Ionic Liquid (IL)	Operating Conditions					Outcomes	
	Pressure bar	Temperature K	Molecular Weight (g.mol <sup>-1</sup> )	Viscosity (mm <sup>2</sup> .s <sup>-1</sup> ) (1 bar and 298.15 K)	Density (g.cm <sup>-3</sup> ) (1 bar and 298.15 K)	$x_{CO_2}$	mol CO <sub>2</sub> /mol IL
PDBF <sub>4</sub>	2	298.15	764.48	198.40	1.09	0.128	0.147
	4					0.193	0.239
	6					0.222	0.287
	8					0.368	0.584
	10					0.559	1.273
	12					0.577	1.365
PDNTf <sub>2</sub>	2	298.15	1151.17	84.74	1.40	0.095	0.105
	4					0.211	0.268
	6					0.334	0.502
	8					0.385	0.626
	10					0.443	0.797
	12					0.471	0.892
PDPF <sub>6</sub>	2	298.15	880.80	44.66	1.29	0.167	0.201
	4					0.181	0.222
	6					0.227	0.294
	8					0.283	0.396
	10					0.421	0.727
	12					0.459	0.849
PDBr	2	298.15	750.67	80.33	1.23	0.039	0.041
	4					0.160	0.191
	6					0.232	0.303
	8					0.243	0.321
	10					0.334	0.503
	12					0.362	0.568