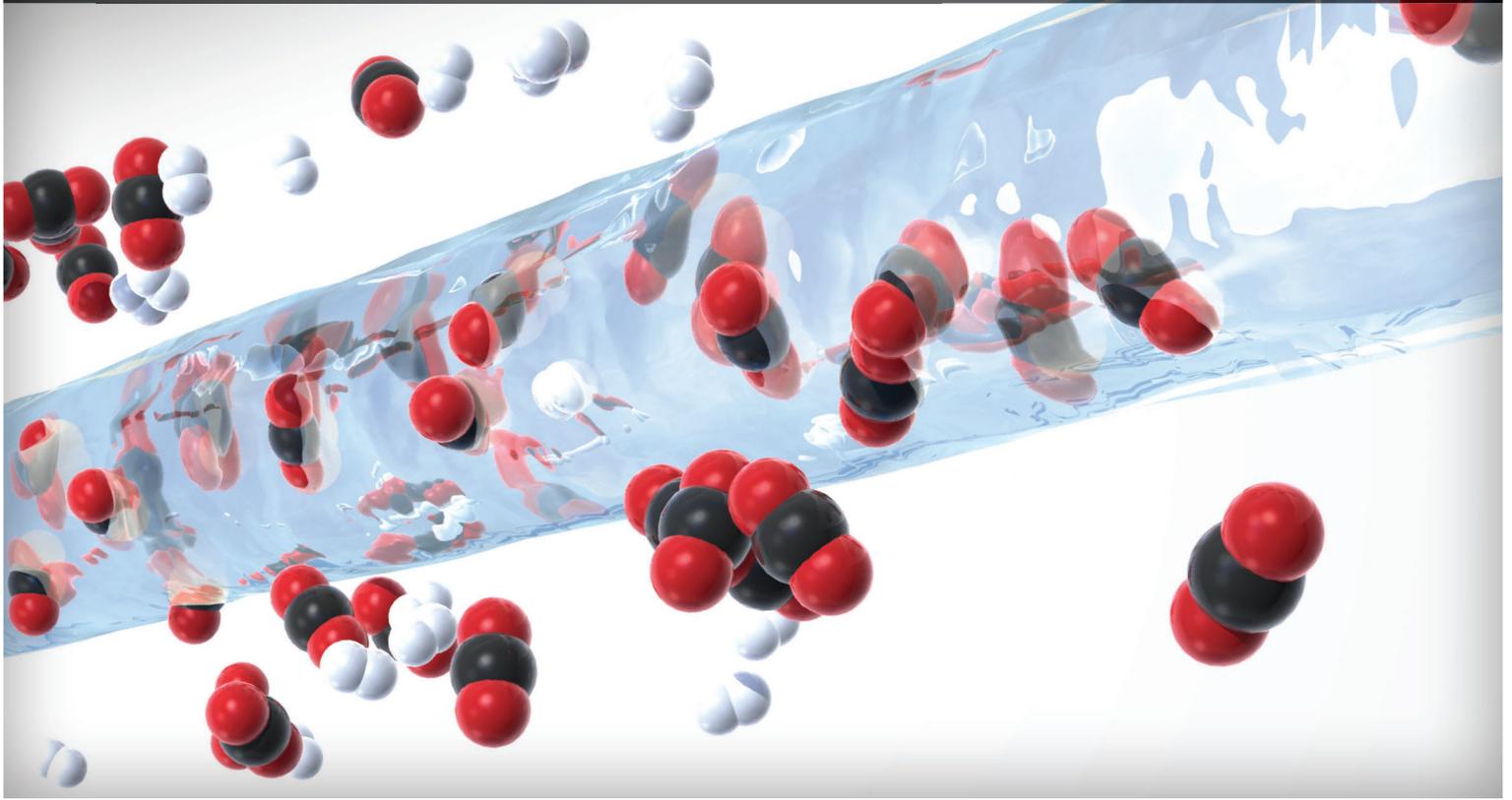


# PRE-COMBUSTION SOLVENTS FOR CARBON CAPTURE



# NETL

NATIONAL ENERGY TECHNOLOGY LABORATORY

## BACKGROUND

Carbon capture and storage from fossil-based power generation is a critical component of realistic strategies for preventing a further rise in atmospheric CO<sub>2</sub> concentrations. However, capturing meaningful amounts of CO<sub>2</sub> using current technology could result in a prohibitive rise in the cost of energy production. One strategy to minimize energy and cost penalties is to exploit industrial streams in which CO<sub>2</sub> is already at high partial pressures, such as the syngas exiting coal gasifiers at integrated gasification combined cycle (IGCC) power plants. In these high-pressure CO<sub>2</sub>-containing streams, one well-established approach to removing acid gases (CO<sub>2</sub>/H<sub>2</sub>S) from the syngas stream (which also contains H<sub>2</sub>/CO/CH<sub>4</sub>/N<sub>2</sub>) is the use of physical solvents. Selexol® (Union Carbide, Houston, Texas, United States) and Rectisol® (Lurgi AG, Frankfurt am Main, Germany) are the standard, commercially available physical solvents for CO<sub>2</sub> capture. Both are hydrophilic, have high vapor pressure, and can cause significant corrosion at elevated temperatures. To avoid this corrosive damage, the syngas temperature for both processes is typically lowered to sub-ambient conditions (10 °C for Selexol and -10 °C for Rectisol) and then raised back up to roughly 200 °C for combustion. This process is inefficient from both a cost and net electricity perspective. Instead, hydrophobic solvents with low vapor pressures could be operated at higher temperatures to minimize the energy and cost penalties associated with cooling the syngas to below ambient conditions.

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## APPROACH

The National Energy Technologies Laboratory (NETL) is developing low viscosity, hydrophobic solvents with low vapor pressures using a comprehensive approach that includes—

- Molecular design and optimization with assistance of computational chemistry
- Synthesis of novel solvents
- Experimental studies at lab/bench scale
- Process design aided by numerical simulations for industrial scale

The research leverages cutting-edge facilities, world-class scientists and engineers, and strategic collaborations to foster the discovery, development, and demonstration of efficient and economical approaches to carbon capture. By drawing on many different scientific and engineering disciplines, NETL researchers have created two new materials that may have superior performance to the existing commercial capture solvents. First, by chemically combining polyethylene glycol diethyl ether (PEGDME)—the active ingredient in Selexol—with hydrophobic silicone oil and polydimethyl siloxane (PDMS), NETL researchers have synthesized a new class of solvents with the best properties of each parent compound. An example from this PEGDME-PDMS class of synthesized solvents is shown in Figure 1 (left). Also shown in Figure 1 (right) is an example from the second class of synthesized solvents. Allyl pyridinium Tf<sub>2</sub>N is an ionic liquid with a low viscosity and extremely high CO<sub>2</sub>/H<sub>2</sub> selectivity. NETL researchers have fully characterized these solvents' CO<sub>2</sub> capture performance and CO<sub>2</sub>/H<sub>2</sub> selectivity between 25 °C and 100 °C and are also measuring their kinetics, mass transfer, regeneration energy, and stability. Each of these properties may be tuned through additives or minor changes to the solvents' molecular structure to optimize CO<sub>2</sub> separation energetics and ultimately to reduce CO<sub>2</sub> capture cost. NETL is also conducting system and economic studies to determine the precise impact of the new materials on the cost of pre-combustion capture from IGCC power plants. NETL-developed solvents have been tested in a slip stream of synthesis gas at the National Carbon Capture Center in Wilsonville, Alabama. Future work includes testing these solvents at a new pilot plant facility being built at NETL to test pre-combustion physical solvents in a continuously looping system consisting of a packed-bed absorber and three flash tanks (Figure 2).

## ACCOMPLISHMENTS

The PEGDME-PDMS hybrid solvents have shown similar or greater CO<sub>2</sub> capacity and CO<sub>2</sub>/H<sub>2</sub> selectivity compared to their parent compounds in capture performance testing. The ionic liquid solvent shown in Figure 1 also has shown exceptional performance in both CO<sub>2</sub> uptake and CO<sub>2</sub>/H<sub>2</sub> selectivity. Other important parameters such as viscosity, stability, and heat capacity are similar to the commercial materials. The primary advantages of these new solvents are their low vapor pressure, their lack of affinity for water, and their low corrosion on carbon steels, all of which allow these solvents to operate at higher temperatures than the conventional hydrophilic solvents.

A U.S. patent has been awarded for the class of materials shown in Figure 1 (left): Patent US 20150114226 A1. In addition, a patent application has been filed for the ionic liquid shown in Figure 1 (right) (U.S. Patent Application 15/045,201). Industrial partners are currently being sought for transfer, scale-up, and commercialization.

These solvents, and similar solvents being synthesized, are expected to be economically viable replacements for Selexol and Rectisol in pre-combustion capture of CO<sub>2</sub> without requiring the syngas to be cooled below ambient conditions.

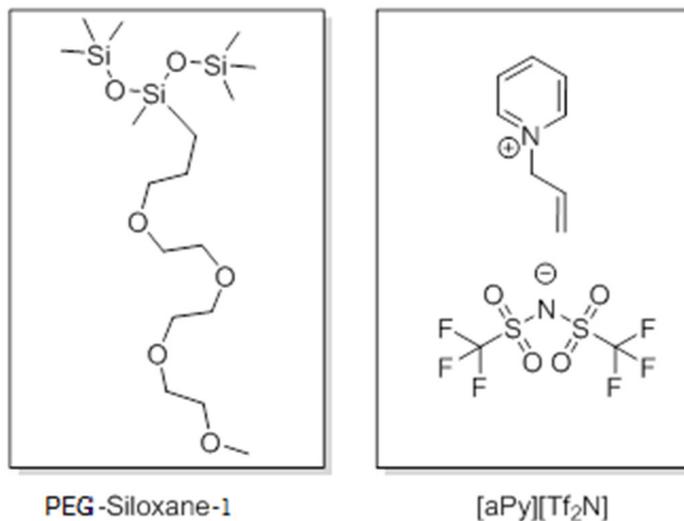


Figure 1. (left) Molecular structure of the hybrid PDMS-PEGDME solvent labeled PEG-Siloxane-1; (right) Molecular structure of the allyl pyridinium Tf<sub>2</sub>N ionic liquid solvent.

# PRE-COMBUSTION SOLVENTS FOR CARBON CAPTURE

Experimental and computational results on these solvents can be found within the following peer reviewed manuscripts:

1. W. Shi, N.S. Siefert, H.O. Baled, J.A. Steckel, D.P. Hopkinson, "Molecular Simulations of the Thermophysical Properties of Polyethylene Glycol Siloxane (PEGS) Solvent for Pre-combustion CO<sub>2</sub> Capture," J. Phys. Chem. C, 120 (36), pp 20158-20169 (2016).
2. N.S. Siefert, S. Agarwal, F. Shi, W. Shi, E.A. Roth, D. Hopkinson, V.A. Kusuma, R.L. Thompson, D.R. Luebke, H.B. and Nulwala, "Hydrophobic physical solvents for pre-combustion CO<sub>2</sub> capture: Experiments, Computational simulations, and Techno-economic analysis," International Journal of Greenhouse Gas Control, 49, 364-371 (2016).
3. W. Shi, N.S. Siefert, and B.D. Morreale, "Molecular Simulations of CO<sub>2</sub>, H<sub>2</sub>, H<sub>2</sub>O, and H<sub>2</sub>S Gas Absorption into Hydrophobic Poly(dimethylsiloxane) (PDMS) Solvent: Solubility and Surface Tension," J. Phys. Chem. C, 119 (33), pp 19253–19265 (2015).

## BENEFITS

This research will move toward the programmatic goal of capturing 90 percent of the CO<sub>2</sub> produced by an IGCC power plant at a cost of less than \$40/tonne CO<sub>2</sub>.

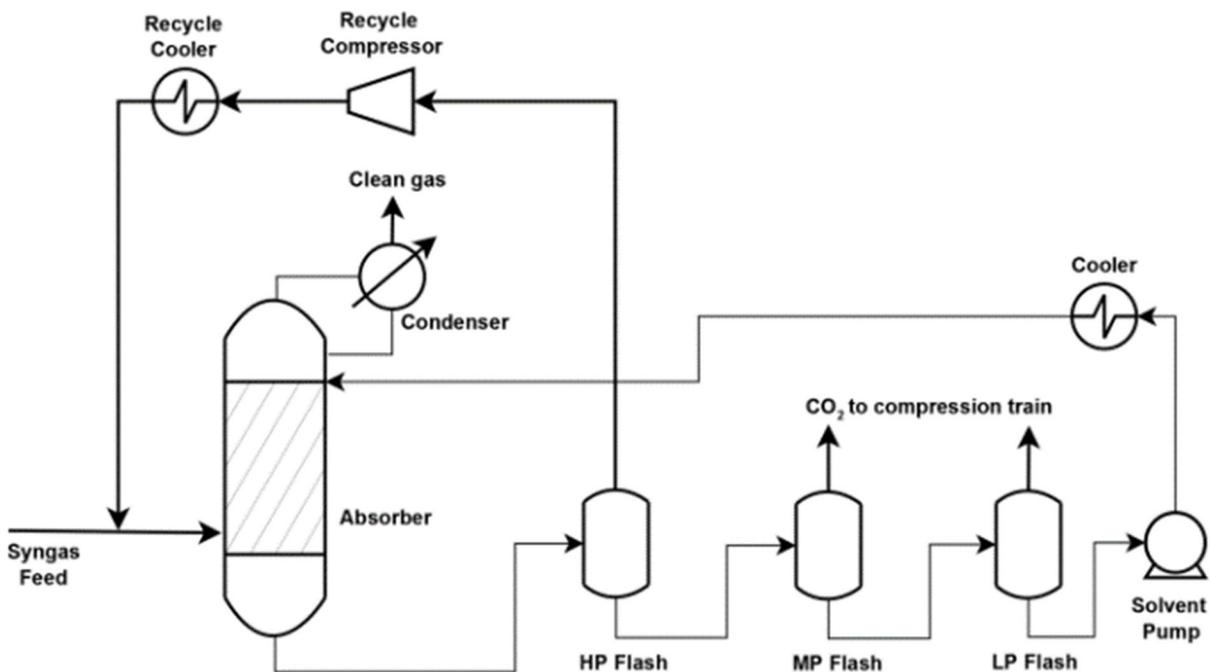


Figure 2. Process flow diagram for NETL's Continuously looping, precombustion CO<sub>2</sub> capture pilot plant facility



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