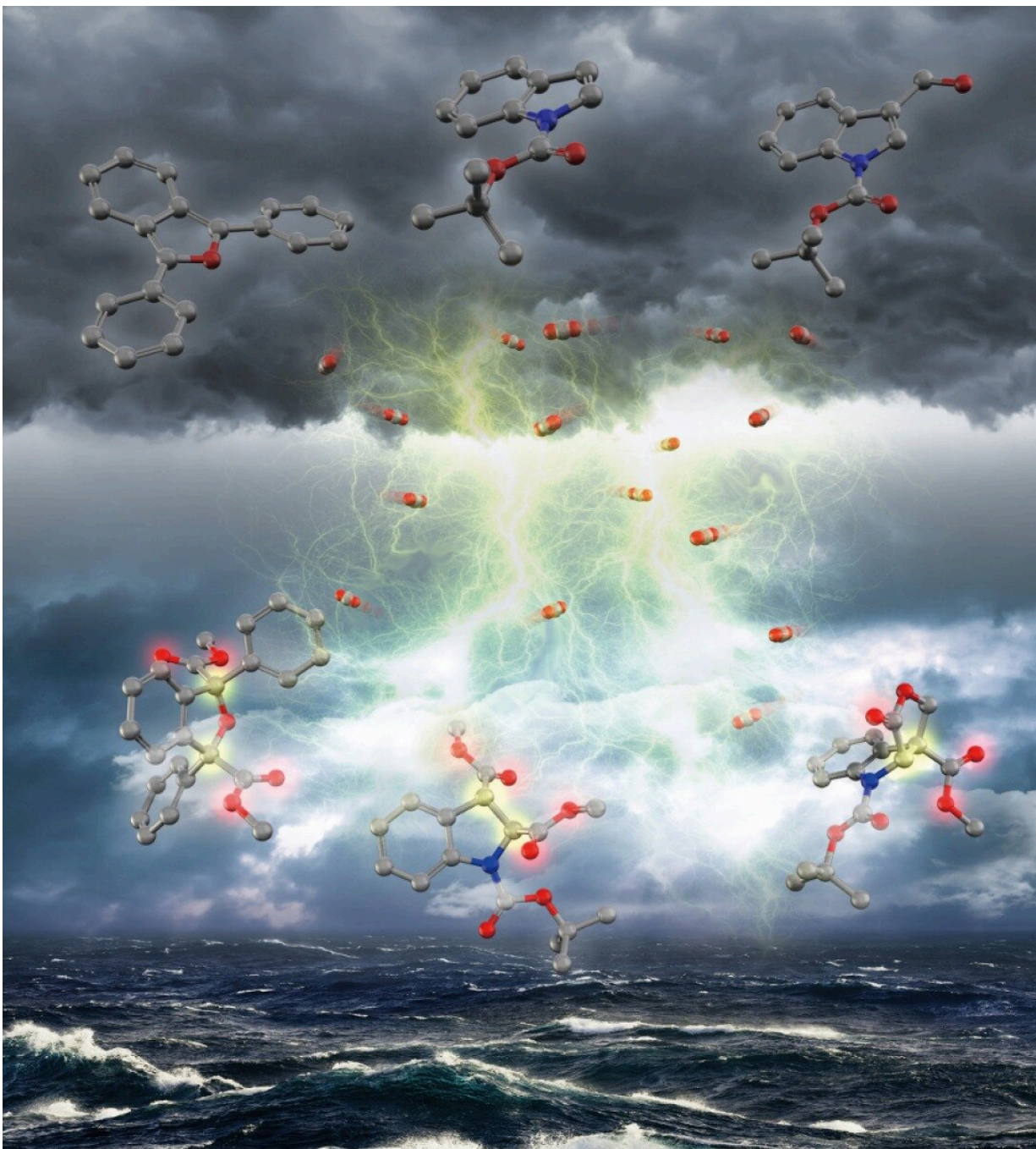


CO₂ recycling and efficient drug development: Tackling two problems with one reaction

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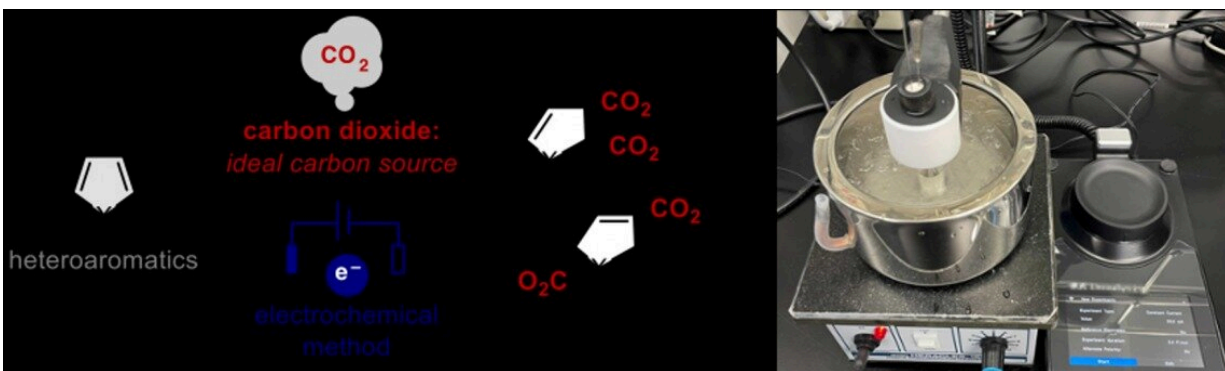
Electricity enabling the addition of CO₂ to heteroaromatic compounds. Credit: Tsuyoshi Mita

Using electricity, a new method offers the possibility of recycling CO₂ while also performing a notoriously difficult reaction, producing compounds potentially useful for drug development.

Scientists at the Institute for Chemical Reaction Design and Discovery (ICReDD) in Hokkaido University have developed a method that has the potential to help recycle waste CO₂ while also producing molecules useful for drug development.

In addition to the ever-more important demand for carbon-neutrality, chemists are increasingly interested in using [carbon dioxide](#) (CO₂) in syntheses since it is abundant, inexpensive, relatively nontoxic, and renewable. However, the reactivity of CO₂ is relatively low. To overcome this, the team led by Professor Tsuyoshi Mita utilized an electrochemical method in which an electron is added to either the CO₂ molecule or to the other molecule in the solution, making it far easier for them to react with each other.

This work marks an especially large breakthrough, since CO₂ is used to carry out a traditionally difficult type of transformation with unprecedented efficiency. When certain conditions are met, electrons can be shared between many atoms in a molecule by what is called an aromatic system. These systems are especially stable and difficult to break, but the new method developed at ICReDD is able to dearomatize, or break, these stable aromatic systems by adding CO₂ to the molecule with the help of electricity. This process has the potential to both recycle CO₂ while also producing high value-added [dicarboxylic acids](#) from simple starting materials, solving two problems at once.



(Left) General process of heteroaromatics undergoing dearomative dicarboxylation with CO₂. (Right) Reaction setup for electrochemical carboxylation. Credit: Yong You, et al., *Journal of the American Chemical Society*, Feb 21, 2022

Prior to the actual experiments, scientists from ICR_eDD screened various heteroaromatic [compounds](#) by calculating their reduction potentials, which is a measure of how a compound will react when subjected to an electric environment. The results enabled researchers to identify potentially reactive compounds and carry out targeted electrochemical experiments. They demonstrate that a wide variety of substrates that exhibit highly negative reduction potentials can very efficiently undergo this unprecedented dearomative addition of two CO₂ [molecules](#). The obtained dicarboxylic acids can be easily and cost-effectively modified into key intermediates for biologically active compounds, which could lead to more efficient and economical [drug development](#). Researchers involved in the study attribute the rapid development of this new process to their strategy of first performing computational analyses that informed their experimental choices in the lab.

"I started to learn computational chemistry when I joined ICR_eDD.

Within one year, I was able to utilize advanced calculation techniques, which was very useful for guiding my decisions in the lab," said first author Dr. Yong You. "It took only eight months to complete the research and publish the paper, which is much faster than a conventional project involving experiments. Significant research time is saved because a computer can reliably predict the feasibility of the reactant structures and possible reaction pathways," commented Tsuyoshi Mita, who led this project.

More information: Yong You et al, Electrochemical Dearomative Dicarboxylation of Heterocycles with Highly Negative Reduction Potentials, *Journal of the American Chemical Society* (2022). [DOI: 10.1021/jacs.1c13032](https://doi.org/10.1021/jacs.1c13032)

Provided by Hokkaido University

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