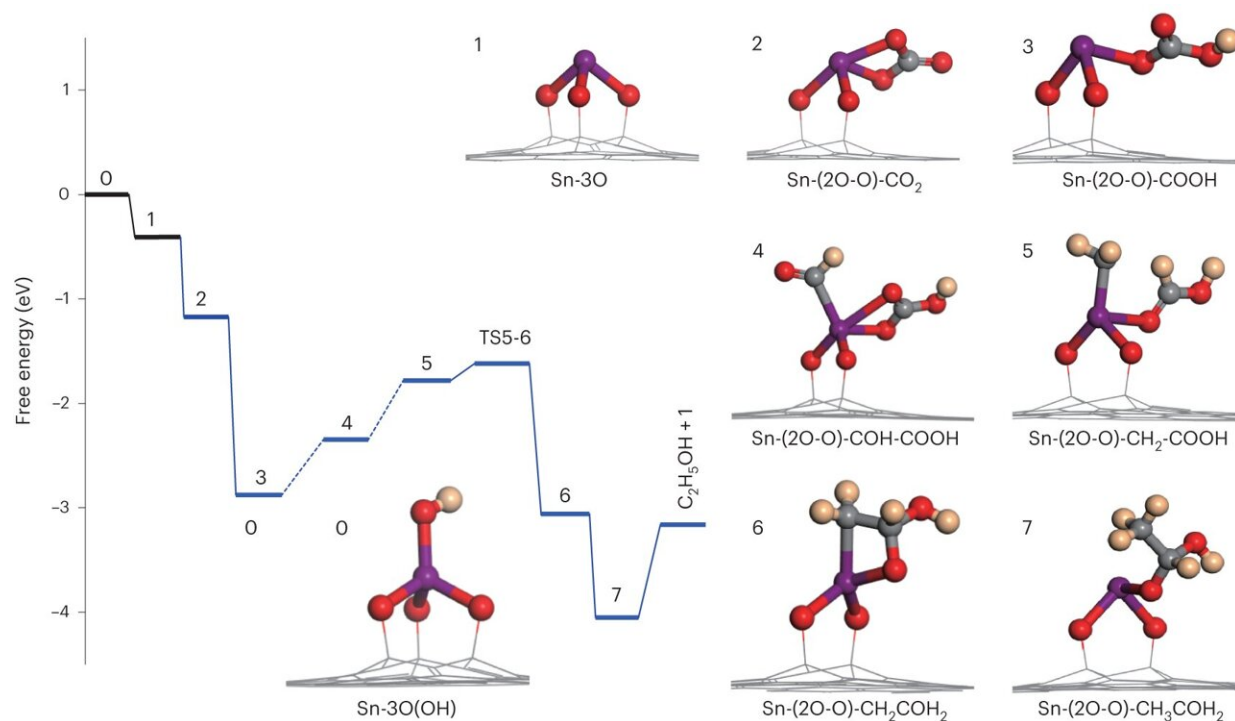


Tandem single-atom electrocatalyst realizes reduction of CO₂ to ethanol

November 13 2023, by Li Yuan



C–C bond formation via a formyl-bicarbonate coupling pathway (with a key intermediate 4). Reaction energy profiles and the corresponding intermediate structures (0 to 7) for the formation of ethanol via the CO₂RR on the Sn₁-O₃G catalyst. Credit: *Nature Energy* (2023). DOI: 10.1038/s41560-023-01389-3

The electrochemical CO₂ reduction reaction (CO₂RR) into carbon-based fuels provides a promising strategy to mitigate CO₂ emission and

promotes the utilization of renewable energy.

The C_n ($n \geq 2$) liquid products are desirable because of their high energy densities and ease of storage. However, manipulation of C–C coupling pathway remains a challenge due to the limited mechanistic understanding.

Recently, a research group led by Profs. Zhang Tao and Huang Yanqiang from the Dalian Institute of Chemical Physics (DICP) of the Chinese Academy of Sciences (CAS) has developed a Sn-based tandem electrocatalyst ($\text{SnS}_2@Sn_1\text{-O3G}$), which could reproducibly yield ethanol with a Faradaic efficiency of up to 82.5% at $-0.9 V_{\text{RHE}}$ and a geometric current density of 17.8 mA/cm^2 .

The study was [published](#) in *Nature Energy* on Oct. 30.

The researchers fabricated the $\text{SnS}_2@Sn_1\text{-O3G}$ through solvothermal reaction of SnBr_2 and thiourea on a three-dimensional carbon foam. The electrocatalyst comprised SnS_2 nanosheets and atomically dispersed Sn atoms ($\text{Sn}_1\text{-O3G}$).

Mechanistic study showed that this $\text{Sn}_1\text{-O3G}$ could respectively adsorb $*\text{CHO}$ and $*\text{CO}(\text{OH})$ intermediates, therefore promoting C–C bond formation through an unprecedented formyl-bicarbonate coupling pathway.

Moreover, by using isotopically labeled reactants, the researchers traced the [pathway](#) of C atoms in the final C_2 product formed over the catalyst of $\text{Sn}_1\text{-O3G}$. This analysis suggested that the methyl C in the product comes from [formic acid](#) whereas the methylene C was from CO_2 .

"Our study provides an alternative platform for C–C bond formation for [ethanol](#) synthesis and offers a strategy for manipulating CO_2 reduction

pathways towards desired products," said Prof. Huang.

More information: Jie Ding et al, A tin-based tandem electrocatalyst for CO₂ reduction to ethanol with 80% selectivity, *Nature Energy* (2023). [DOI: 10.1038/s41560-023-01389-3](https://doi.org/10.1038/s41560-023-01389-3)

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