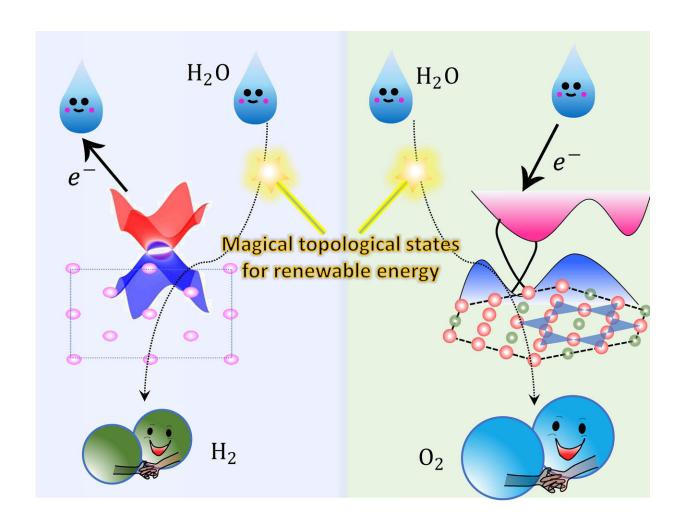


Speeding up hydrogen production by the magic topological surface states

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Topological non-trivial surface states can accept or donate electrons during the water electrolysis process. Credit: MPI CPfS



The hydrogen economy is considered to be one of the best options for providing renewable energy and, thereby, contributing to mitigating today's environmental challenges. The energy density of hydrogen is between 120-142 MJ/kg, which is much greater than that of chemical, fossil, and biofuels. More importantly, water is the only byproduct when hydrogen is used to produce electricity.

Water electrolysis could provide high-quality hydrogen gas that can be used in fuel cells directly. However, since noble metals, such as platinum and iridium, are currently needed to initiate such a reaction, the cost is very high. "Obviously, catalysts that are low-cost with high-activity are needed to make hydrogen energy more competitive with traditional technologies," says Guowei Li at the Max Planck Institute for Chemical Physics of Solids, who studied the surface reactions of several topological materials.

It was obviously a great challenge to find alternatives beyond noble metals. "Topology may be the key to unlocking the barrier in the search for ideal catalysts," says Prof. Claudia Felser, the director of Max Planck Institute for Chemical Physics of Solids. "We studied the surface properties of materials with topological order, from topological insulators to topological semimetals and metals, all these materials have non-trivial surface states that are protected by symmetries."

"In other words, these surface states are very stable and robust against surface modifications such as impurity scattering and even oxidation: the question we're asking is can we find such a perfect system that combines topological order, lost-cost, high efficiency, and high stability."

The team from the Max Planck Institute Chemical Physics of Solids, Dresden, together with colleagues from the TU Dresden and the Max Planck Institute for Microstructure Physics and Max-Planck-Institut für Kohlenforschung, Mülheim published a breakthrough result in *Science*



Advances concerning a topological material, namely a magnetic Weyl semimetal, that is a superior oxygen evolution reaction (OER) <u>catalyst</u>. The magnetic weyl semimetal that the team identified is $Co_3Sn_2S_2$, a Kagome-lattice Shandite compound.

High-quality bulk single crystals of $Co_3Sn_2S_2$ with sizes of up to centimeters can be exfoliated into thin layers with defined crystal surfaces. The team showed that these surfaces act as superior catalysts for water splitting, even though the surface area is several orders of magnitudes smaller than today's conventional nano-structured catalysts. In collaboration with Yan Sun's theory group from the Max Planck Institute of Chemical Physics of Solids, they found that there are cobaltderived topological surface states just above the Fermi level. In the water oxidation process, these surface states can accept electrons from the reaction intermediates, acting as an electron channel whose resistance is not affected by the harsh electrochemical environment.

Inspired by this strategy, the team then investigated the catalytic performance of a Dirac nodal arc semimetal $PtSn_4$, a compound that has a much lower percentage of expensive <u>platinum</u>. Such crystals showed superior electrocatalytic stability for periods of time exceeding one month.

"The work serves as an interesting lens into the chemistry of these reaction processes and could be a pathway towards understanding the chemistry itself by clear knowledge of the topological nature of the semimetal catalyst," says one of the expert reviewers of the paper.

More information: Guowei Li et al, Surface states in bulk single crystal of topological semimetal $Co_3Sn_2S_2$ toward water oxidation, *Science Advances* (2019). <u>DOI: 10.1126/sciadv.aaw9867</u>

Guowei Li et al. Dirac Nodal Arc Semimetal PtSn4 : An Ideal Platform



for Understanding Surface Properties and Catalysis for Hydrogen Evolution, *Angewandte Chemie International Edition* (2019). DOI: 10.1002/anie.201906109

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