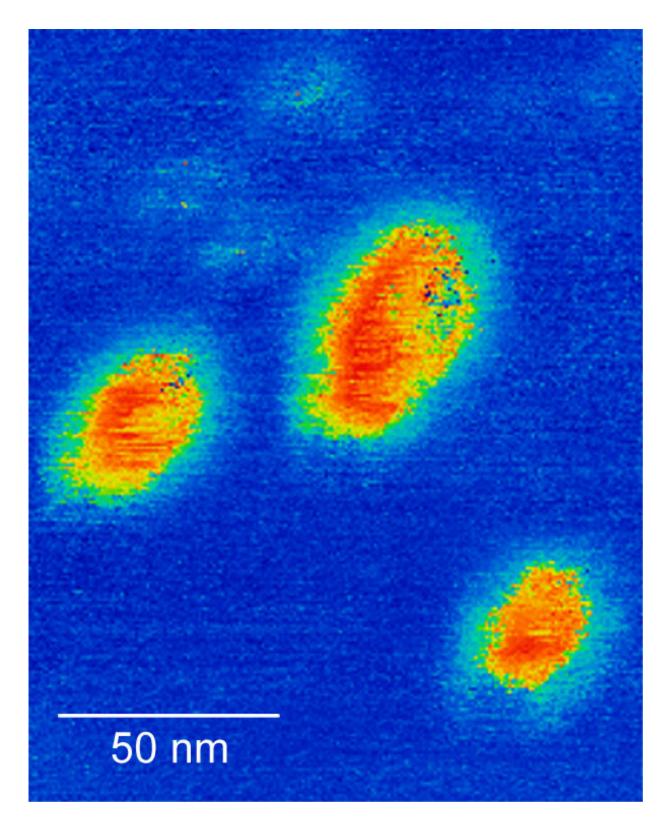


# First observation of native ferroelectric metal

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Ferroelectric domains in a  $WTe_2$  single crystal (PFM imaging). Credit: FLEET



In a paper released today in *Science Advances*, Australian researchers describe the first observation of a native ferroelectric metal: a native metal with bistable and electrically switchable spontaneous polarization states—the hallmark of ferroelectricity. The study found coexistence of native metallicity and ferroelectricity in bulk crystalline tungsten ditelluride (WTe<sub>2</sub>) at room temperature. A van-der-Waals material that is both metallic and ferroelectric in its bulk crystalline form at room temperature has potential for nano-electronics applications.

The study represents the first example of a native metal with bistable and electrically switchable spontaneous polarization states—the hallmark of ferroelectricity.

"We found coexistence of native metallicity and ferroelectricity in bulk crystalline tungsten ditelluride (WTe<sub>2</sub>) at room temperature," explains study author Dr. Pankaj Sharma.

"We demonstrated that the ferroelectric state is switchable under an external electrical bias and explain the mechanism for 'metallic ferroelectricity' in WTe<sub>2</sub> through a systematic study of the crystal structure, electronic transport measurements and theoretical considerations."

"A van der Waals material that is both metallic and ferroelectric in its bulk <u>crystalline form</u> at room temperature has potential for new nanoelectronics applications," says author Dr. Feixiang Xiang.

## Ferroelectric backgrounder

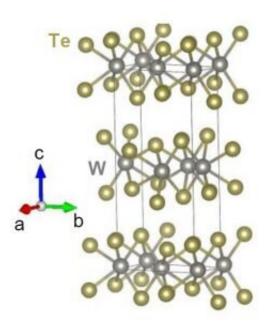
Ferroelectricity can be considered an analogy to ferromagnetism. A ferromagnetic material displays permanent magnetism, and in layperson's terms, is simply, a 'magnet' with north and south pole. Ferroelectric material likewise displays an analogous electrical property



called a permanent electric polarisation, which originates from electric dipoles consisting of equal, but oppositely charged ends or poles. In ferroelectric materials, these electric dipoles exist at the unit cell level and give rise to a non-vanishing permanent electric dipole moment.

This spontaneous electric dipole moment can be repeatedly transitioned between two or more equivalent states or directions upon application of an external electric field—a property utilised in numerous ferroelectric technologies, for example nano-electronic computer memory, RFID cards, medical ultrasound transducers, infrared cameras, submarine sonar, vibration and pressure sensors, and precision actuators.

Conventionally, <u>ferroelectricity</u> has been observed in materials that are insulating or semiconducting rather than metallic, because conduction electrons in metals screen-out the static internal fields arising from the dipole moment.



Model of tungsten ditelluride WTe<sub>2</sub> crystals in a layered, orthorhombic structure. Credit: FLEET



#### The study

A <u>room-temperature</u> ferroelectric semimetal was published in *Science Advances* in July 2019.

Bulk single-crystalline tungsten ditelluride (WTe<sub>2</sub>), which belongs to a class of materials known as transition metal dichalcogenides (TMDCs), was probed by spectroscopic electrical transport measurements, conductive-atomic force microscopy (c-AFM) to confirm its metallic behaviour, and by piezo-response force microscopy (PFM) to map the polarisation, detecting lattice deformation due to an applied electric field.

Ferroelectric domains—ie, the regions with oppositely oriented direction of polarization—were directly visualised in freshly-cleaved WTe<sub>2</sub> single crystals.

Spectroscopic-PFM measurements with top electrode in a capacitor geometry was used to demonstrate switching of the ferroelectric polarization.

The study was supported by funding from the Australian Research Council through the ARC Centre of Excellence in Future Low-Energy Electronics Technologies (FLEET), and the work was performed in part using facilities of the NSW Nodes of the Australian National Fabrication Facility, with the assistance of the Australian Government Research Training Program Scholarship scheme.

First-principles density functional theory (DFT) calculations (University of Nebraska) confirmed the experimental findings of the electronic and structural origins of the ferroelectric instability of WTe<sub>2</sub>, supported by



the National Science Foundation.

#### Ferroelectric studies at FLEET

Ferroelectric materials are keenly studied at FLEET (the ARC Centre of Excellence in Future Low-Energy Electronics Technologies) for their potential use in low-energy electronics, 'beyond CMOS' technology.

The switchable electric dipole moment of ferroelectric materials could for example be used as a gate for the underlying 2-D electron system in an artificial topological insulator.

In comparison with conventional semiconductors, the very close (subnanometre) proximity of a ferroelectric's electron dipole moment to the electron gas in the atomic crystal ensures more effective switching, overcoming limitations of conventional semiconductors where the conducting channel is buried tens of nanometres below the surface.

Topological <u>materials</u> are investigated within FLEET's Research theme 1, which seeks to establish ultra-low resistance electronic paths with which to create a new generation of ultra-low energy electronics.

**More information:** "A room-temperature ferroelectric semimetal" *Science Advances*, <u>DOI: 10.1126/sciadv.aax5080</u>, advances.sciencemag.org/content/5/7/eaax5080

### Provided by FLEET

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