

because the non-metallic [element](#) carbon prefers covalent, directional bonds that effectively cause structural planarity. Metallic bonds are less directional, whereby metals often coalesce into compact clusters. However, recent experiments suggest that, by using pores in other nanostructures, even metals can be stabilized into atomically thin planes.

Inspired by these experimental indications, researchers at the Nanoscience Center, University of Jyväskylä, used computer simulations to predict systematically the properties of atomically thin structures made exclusively from [metallic elements](#).

"We made a new opening in material research, which was basic research, but highly rewarding as such," says postdoctoral researcher Janne Nevalaita. "One could say that we hit on an untouched estate, bulldozed it and created a foundation. Now others can build solid scientific structures based on that foundation," he continues.

According to the simulations, the flat metals inherit their properties predictably from conventional three-dimensional [metal](#) structures. The study appeared in *Physical Review B*, a journal published by the American Physical Society. It was selected as an Editor's suggestion and highlighted as an esteemed Physics synopsis.

More information: Janne Nevalaita et al. Atlas for the properties of elemental two-dimensional metals, *Physical Review B* (2018). [DOI: 10.1103/PhysRevB.97.035411](https://doi.org/10.1103/PhysRevB.97.035411)

Provided by University of Jyväskylä

Citation: New research opening for atomically thin metal nanostructures (2018, January 12) retrieved 10 April 2024 from

<https://phys.org/news/2018-01-atomically-thin-metal-nanostructures.html>

This document is subject to copyright. Apart from any fair dealing for the purpose of private study or research, no part may be reproduced without the written permission. The content is provided for information purposes only.