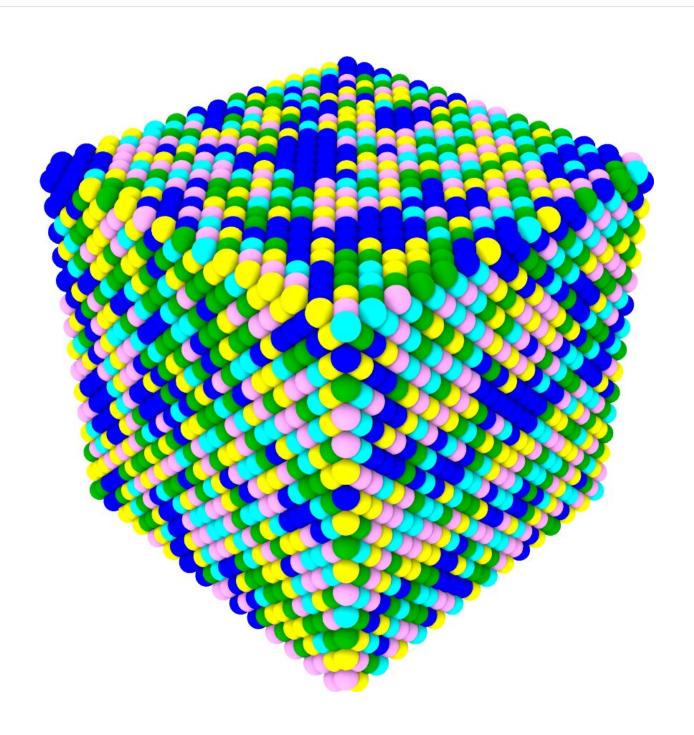


Atomic-level imaging could offer roadmap to metals with new properties

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This schematic illustration of the new palladium-containing high entropy allow shows how new alloy contains large palladium clusters (blue atoms). Credit: Ting Zhu

High-entropy alloys, which are made from nearly equal parts of several primary metals, could hold great potential for creating materials with superior mechanical properties.

But with a practically unlimited number of possible combinations, one challenge for metallurgists is figuring out where to focus their research efforts in a vast, unexplored world of metallic mixtures.

In a study published Oct. 9 in the journal *Nature*, the researchers described using energy-dispersive X-ray spectroscopy to create maps of individual metals in two high-entropy alloys. This spectroscopy technique, used in conjunction with transmission electron microscopy, detects X-rays emitted from a sample during bombardment by an electron beam to characterize the elemental composition of an analyzed sample. The maps show how <u>individual atoms</u> arrange themselves within the alloy, allowing researchers to look for patterns that could help them design alloys emphasizing individual properties.

For example, the maps could give researchers clues to understand why substituting one metal for another could make an alloy stronger or



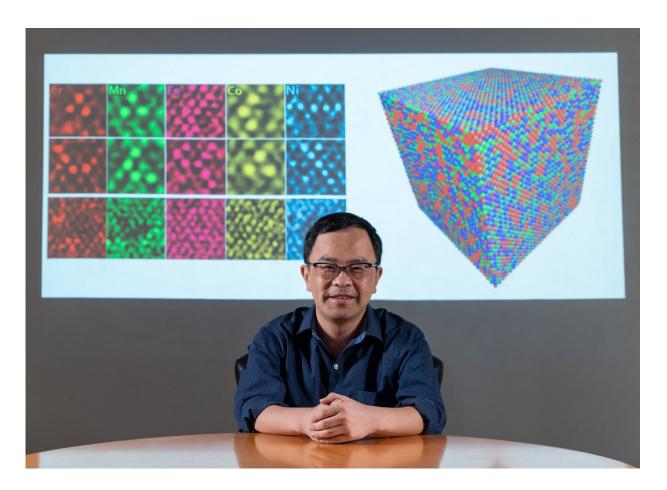
weaker, or why one metal outperforms others in extremely cold environments.

"Most alloys used in engineering applications have only one primary metal, such as iron in steel or nickel in nickel-based super alloys, with relatively small amounts of other metals," said Ting Zhu, a professor in the George W. Woodruff School of Mechanical Engineering at Georgia Tech. "These new alloys that have relatively high concentrations of five or more metals open up the possibility of unconventional alloys that may have unprecedented properties. But this is a new compositional space that has not been explored, and we still have a very limited understanding of this class of materials."

The name "high entropy" refers to the lack of uniformity in the mixture of metals as well as how many different and somewhat random ways the atoms from the metals can be arranged as they are combined.

The new maps could help researchers determine whether there are any unconventional atomic structures that such alloys take that could be leveraged for engineering applications, and how much control researchers could have over the mixtures in order to "tune" them for specific traits, Zhu said.





Ting Zhu, a professor in the George W. Woodruff School of Mechanical Engineering at Georgia Tech, has helped develop a new process to gain insights into individual high-entropy alloys and help characterize their properties. Credit: Rob Felt

To test the new imaging approach, the research team compared two highentropy alloys containing five metals. One was a mixture of chromium, iron, cobalt, nickel, and manganese, a combination commonly referred to as a "Cantor" alloy. The other was similar but substituted palladium for the manganese. That one substitution resulted in much different behavior in how the atoms arranged themselves in the mixture.

"In the Cantor alloy, the distribution of all five elements is consistently



random," Zhu said. "But with the new alloy containing palladium, the elements show significant aggregations due to the much different atomic size of palladium atoms as well as their difference in electronegativity compared to the other elements."

In the new alloy with palladium, the mapping showed that palladium tended to form large clusters while cobalt seemed to collect in places where iron was in low concentrations.

Those aggregations, with their sizes and spacings in the range of a few nanometers, provide strong deformation resistance and could explain the differences in mechanical properties from one high-entropy alloy to another. In straining tests, the alloy with palladium showed higher yield strength while keeping similar strain hardening and tensile ductility as the Cantor alloy.

"The atomic scale modulation of element distribution produces the fluctuation of lattice resistance, which strongly tunes dislocation behaviors," said Qian Yu, a coauthor of the paper and a professor in Zhejiang University. "Such modulation occurs at a scale that is finer than precipitation hardening and is larger than that of traditional solid solution strengthening. And it provides understanding for the intrinsic character of high-entropy alloys."

The findings could enable researchers to custom design <u>alloys</u> in the future, leveraging one property or another.

The team also included researchers from the University of Tennessee, Knoxville; Tsinghua University; and the Chinese Academy of Sciences.

More information: Tuning element distribution, structure and properties by composition in high-entropy alloys, *Nature* (2019). <u>DOI:</u> 10.1038/s41586-019-1617-1, nature.com/articles/s41586-019-1617-1



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