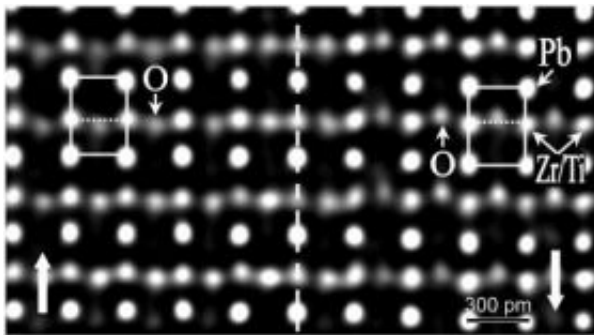


# Electron microscopy enters the picometer scale

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Using electron microscope methods of a hitherto unknown accuracy, scientists from Forschungszentrum Juelich have succeeded in locally demonstrating polarization in the ferroelectric  $\text{PbZr}_{0.2}\text{Ti}_{0.8}\text{O}_3$  and measuring it atom by atom. The broken line forms the boundary of two areas with different electrical polarization marked by the arrows. This is due to the fact that the atoms (Pb: lead; Z: zircon; Ti: titanium; O: oxygen) are displaced from their positions and therefore their electrical charges cannot compensate for each other. On the left, the oxygen atoms are displaced 38 pm downwards, and on the right to the same degree upwards out of the zircon/titanium atomic row. This row itself is displaced vertically by 10 pm from the center line between the lead atoms. In order to write information in applications for data storage, the boundary between these two areas of different polarization directions is displaced to the left or to the right so that only one polarization direction exists in the material. Image: Forschungszentrum Juelich

Jülich scientists have succeeded in precisely measuring atomic spacings down to a few picometres using new methods in ultrahigh-resolution

electron microscopy. This makes it possible to find out decisive parameters determining the physical properties of materials directly on an atomic level in a microscope. Knut Urban from Forschungszentrum Jülich, a member of the Helmholtz Association, reports on this in the latest issue (25 July) of the scientific high-impact journal *Science*.

Progress in research in the area of physics is very frequently connected to an increase in the accuracy of measurements, which help researchers to track natural phenomena. With the aid of new methods in electron optics, researchers were able to microscopically measure atomic displacements precisely to a few picometres. A picometre corresponds to a billionth of a millimetre a distance that is one hundred times smaller than the diameter of an atom.

This is one of the highlights that Knut Urban, director of the Ernst Ruska-Centre in Jülich, reports on in *Science* as part of a review of ten years of electron microscopy with aberration-corrected lenses.

Jülich scientists investigated, for example, the configuration of atoms in orthogonal grain boundaries of the oxide superconductor  $\text{YBa}_2\text{Cu}_3\text{O}_7$ . These atoms mark the boundary between two areas of the crystalline material with atomic structures that are tilted at an angle of exactly  $90^\circ$  to each other. From microscopic images taken under different conditions, the physicists succeeded in using computers to calculate the quantum-mechanical wave function of the electrons, which served as a basis for determining the exact position of the atoms.

In doing so, it became apparent that the relatively heavy atomic species barium, copper and yttrium are systematically displaced a few picometres from their ideal position in the grain boundary and that the lighter oxygen atoms follow this displacement. This provides an explanation for the attenuation of superconducting properties, which can be observed when electric current flows over such a grain boundary. This

phenomenon is undesired if the superconductor is intended to be used for a loss-free current transport. However, it is useful for the construction of so-called SQUIDs (superconducting quantum interference devices), which exploit the magnetic field dependency of this disturbance to measure smallest magnetic fields, for example, to measure brain waves (magnetoencephalography).

Displacements of a few picometres decide on a whole number of physical properties, which are of eminent importance for technology. Another example is the ferroelectricity of titanates materials. Here, the electrical charges of the individual types of atoms inside the building blocks of crystals, the unit cells, cannot fully compensate for each other as they are not arranged in the necessary symmetry.

Therefore, electric dipoles are formed inside the unit cells, which add up over a larger crystal area to form the so-called polarisation. This is used to write information bits. An example is  $\text{PbZr}_{0.2}\text{Ti}_{0.8}\text{O}_3$  which is used in chip cards for data storage. With the aid of new electron optical methods, atomic displacements can be measured atom by atom thus making it possible to determine local polarisation for the first time.

Knut Urban explains: "This is the beginning of a new physics of materials which enables researchers to determine physical parameters and properties in the nano range through highly precise measurements of the atomic spacings. This will also provide clues on how these properties may be manipulated in order to gain new functions and better functional performance."

Source: Helmholtz Association of German Research Centres

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