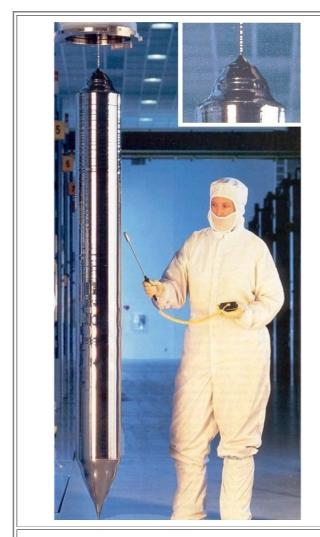
Single Crystal Silicon

Here are a few pictures of the biggest and most perfect crystal that is known on this side of Jupiter.



Freshly grown Si single crystal with 200 mm diameter

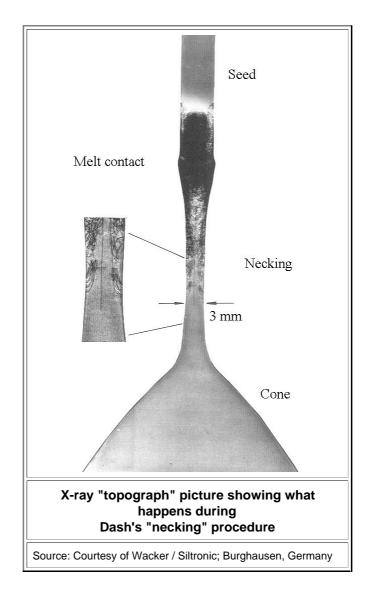
The inset shows the top part of the crystal, showing that the \approx 150 kg heavy crystal is hanging on a seed crystal just a few mm in diameter.

Some Smithsonian magazine (?)

The thin neck is absolutely essential for growing a dislocation-free crystal. Below is a rather amazing picture where we look with X-rays into the neck region of a silicon crystal in such a way that we see dislocations as thin dark lines.

When you dip your already quite thin seed crystal into the hot melt (1410 °C; 2570°F), the thermal shock generates stress, and stress generates dislocations at this temperature quite easily. You can't avoid that. What you do then is: First pull rather fast. The diameter of the growing crystal will then decrease to a few mm. This is the "Dash process", invented by W.C. Dash in 1958, "just in time". The trick, in a nutshell, is to pull faster than the dislocations can move. That demands a very thin crystal, and as the picture shows, it works! The dislocations move out to the surface and eventually the growing crystal is dislocation free. Then you slow down and increase the diameter to whatever you need, e.g. 200 mm or 300 mm.

This "necking procedure", or the "Dash necking process" is one of the many little marvels that enabled silicon technology. Of course, the <u>first law of economics</u> still applies. You end up with a huge and heavy (up 250 kg) crystal hanging on a thread or better threat! If the thin neck breaks while you are pulling, and your already grown crystal plunges into the hot melt, you better start running fast. You must outrun either the exploding machine or your boss, who will try to get you for destroying equipment worth millions!



Here is one of the humongous 300 mm diameter Si crystals mass-produced today (2011):



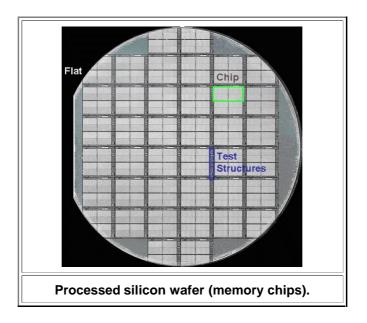
Iron, Steel and Swords script - Page 2

Silicon single crystal with 300 mm diameter

Source: Courtesy MEMC.

The cylindrical part of these crystals will be cut into discs that are subsequently ground and polished to the ubiquitous **silicon wafers** as shown below.

Now you have the starting material for all kinds of "micro"-electronic nano devices, like all your electronic stuff, power managing components, solar cells, and not to forget, *MEMS* products like the acceleration sensors and gyros in your car and your i-phone, or the light-processing chips with millions of tiny mirrors in your beamer.



- These silicon single crystals do not contain a single <u>dislocation</u>. They are so far the only big crystals that we can make completely free of dislocations. Germanium (Ge) crystals can be pretty good, too, but are not quite as perfect as silicon crystals. They are also far more expensive since "raw" germanium is expensive.
 - That is an extraordinary occurrence! Microelectronics needs materials with some special electronic properties and rather perfect, large, dislocation-free crystals of that material. Germanium doesn't meet the electronic specifications (not to mention cost considerations). Gallium arsenide crystals or other "III-V's" good enough for making electronic components like LED's or Lasers are rather expensive but far less perfect than silicon and never dislocation free.
 - It is a kind of amazing coincidence that the basic properties of silicon meets all electronic specifications admirably *and* that it is the *only* material that can be grown into huge and *nearly* perfect single crystals! Those crystals are also extremely clean. Besides a few doping atoms (As, P, or B) incorporated intentionally at precisely determined concentrations, they only contain some dissolved oxygen (a few ppm), and somewhat less than 1 ppm carbon. Those impurities are tolerated or even helpful. Everything else is far below the ppt level
- By the way, there cannot be such thing as a fully perfect Si crystal. Even the most perfect ones must by necessity contain a low density of small clusters of vacancies and self-interstitials.
 Why? Consult this link to find out.