

INDIA

## Scientists decry killings of secularists

Indian academy members condemn intolerance.

BY T. V. PADMA

Indian scientists are voicing concerns over religious intolerance and the killings of three noted advocates of rational thinking. The actions are unusual in a country where scientists rarely comment on political issues, says physicist Shri Krishna Joshi, a member of India's Inter-Academy Panel on Ethics in Science.

Anti-superstition activist Narendra Dabholkar was killed in 2013, communist politician Govind Pansare in February this year and literature scholar Malleshappa Kalburgi in August. All three deaths have been blamed on members of extreme right-wing Hindu groups.

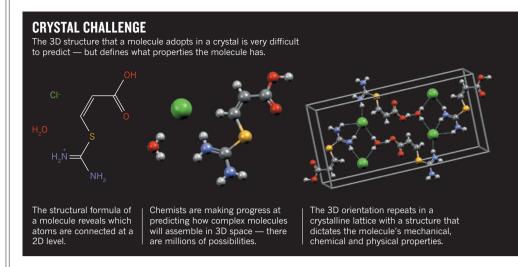
On 22 October, scientists launched an online petition to India's president, Pranab Mukherjee, protesting against the killings. "The government has failed to check or discourage the anti-rational environment," says petition signatory Naresh Dadhich, a physicist at the Inter-University Centre for Astronomy and Astrophysics in Pune, India.

The petition was followed on 27 October by a statement from the Inter-Academy Panel on Ethics in Science, set up by the Indian National Science Academy in New Delhi; the Indian Academy of Sciences, Bangalore; and the National Academy of Sciences in Allahabad. The Indian constitution mandates that "its citizens abide by and uphold reason and scientific temper", the statement said. Several statements and actions "run counter to this constitutional requirement," it notes.

Indira Nath, a member of the panel and an immunologist at the Indian National Science Academy, says that the panel wants to "bring back rationality and scientific thinking to the mainstream".

Also last week, more than 100 scientists from leading Indian institutes, including national award winners, three fellows of the Royal Society in London, and a foreign associate of the US National Academy of Sciences, signed a second statement expressing deep concern over the "climate of intolerance".

Pushpa Mittra Bhargava, former director of the Centre for Cellular and Molecular Biology in Hyderabad, says that he plans to return a national award in protest. "Science is about reason and rationality. If three rationalists can be killed, scientists too can be killed." ■



CHEMISTRY

## Software predicts crystal structures

Chemists have succeeded at a fiendish task — forecasting how complex molecules will assemble in 3D.

## BY ELIZABETH GIBNEY

Sketch the structure of an organic molecule on a napkin and it may not be apparent that there are millions of possible ways that it could assemble as a 3D crystal. Now, a collaboration of dozens of chemists and computer programmers has successfully predicted the crystal structure of five, complex, 'drug-like' organic molecules — using nothing but a 2D map showing which atoms connect to which.

The achievement, announced on 27 October at a workshop in Cambridge, UK, paves the way for software that would cut the cost of the design and manufacture of drugs and other chemical products, as well as further our understanding of fundamental chemistry.

A molecule's crystal structure determines its properties (see 'Crystal challenge'). In 1998, the US pharmaceutical firm Abbott Laboratories learned this the hard way when it had to pull production of the capsule form of the HIV treatment ritonavir because the drug had started switching to an unexpected structure during manufacture. The crystal structure that a molecule adopts is generally the one with the lowest energy, but predicting what this is for any particular molecule is "fiendishly difficult", says Colin Groom, executive director of the Cambridge Crystallographic Data Centre (CCDC).

Even when chemists know which atoms are connected to which, the atoms can still be in different orientations because the bonds that connect them can bend and rotate in myriad ways. There are also multiple options for how molecules can pack together. "It is like looking for a needle in an unimaginably big haystack," says Anthony Reilly, a structural chemist at the CCDC.

Since 1999, the CCDC has organized six challenges known as the Blind Test of Organic Crystal Structure Prediction Methods. Rather than a contest, organizers see the challenge as a large collaborative attempt to compare the strengths of the latest techniques. "The groups participating represent pretty much the entire crystal-structure prediction community, and the methods used are the very best developed," says Groom.

The challenge typically takes place over a year, and sets two major problems. First, teams must come up with a list of all possible arrangements in which the molecules could form a crystal. Some teams do a rough calculation of the energy of each to whittle down the list, burning up hundreds of thousands of hours of computing time; others start with pure guesses and iteratively 'breed' the most stable to derive possible candidates more quickly. In the second stage, teams take the shortlists — sometimes assembled by a different group — and do more-precise