

properties in self assembled supramolecular materials, considering host-guest intergrowth crystals. A host urea molecule forms hydrogen-bonded hexagonal channels of 5.5 Å diameter in which long guest chains are densely packed in a one-dimensional arrangement [1]. This simple paradigm crystal offers a unique opportunity to address at a fundamental level the question of the nature and the roles of interactions in self-organized architectures [2]. Original physical properties in these host-guest materials are related to their aperiodic feature which, unambiguously, appears in their superspace diffraction pattern [3,4]. Here, we will present a first evidence of a superspace symmetry breaking involving only the internal variable of the superspace in these materials. These observations force a total reconsideration of the interactions in these self-organized compounds.

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Structural Study of Single Crystalline i-Zn-Mg-Dy at high Pressure and high Temperature

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In order to understand the origin of the structural stability of quasicrystals, it is important to investigate the pressure and temperature dependence of their structures. Quasicrystalline phases in the Zn-Mg-R (R = rare-earth and Y) alloys are classified as the Frank-Kasper type. Both the icosahedral and decagonal phases have been obtained in the alloys, which are thermodynamically stable at ambient conditions and reveal a high structural perfection [1]. The icosahedral Zn-Mg-Y quasicrystal is found to be stable at high pressures up to 70GPa at room temperature [2]. The hardness seems to be primarily governed by the complexity of the structure and the bonding strength. At room temperature, Zn-Mg-R quasicrystals exhibit a strong indentation size effect with the hardness increasing with decreasing load. This effect becomes inverted at higher temperatures [3]. Therefore it is interesting to perform structural studies at HP/HT.

We will report on the results of an in-situ single crystal x-ray diffraction study on i-Zn-Mg-Dy up to ca 12GPa and 873K using a heatable diamond anvil cell. The icosahedral quasicrystal is found to be essentially stable within the experimental framework.

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Framework Structures for Quasicrystal Models based on Dense Icosahedral Sphere Packings

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Dense icosahedral sphere packings (DISP's) are known to be obtained by placing spheres on subset vertices of three-dimensional Penrose tiling (3DPT) which are called the twelve-fold packing sites [1]. The DISP's give the best known quasiperiodic cluster packings with *b*- and *c*-links along their twofold and threefold directions, respectively ($b=2.75a$ and $c=2.38a$, where a is the edge length of the 3DPT). However the shapes of the possible interstices are unknown.

Instead of spheres, we consider to place rhombic triacontahedra (RT's), having an edge length of a , on the sites. As a result, the RT's

share a rhombus face and an obtuse rhombohedron with *b*-link and *c*-link, respectively. These framework structures are well described by the section method in six-dimension, once occupation domains for the DISP's are specified. It is shown that the interstices can be described by acute and obtuse rhombohedra. Application of the present framework structures to p-type icosahedral quasicrystals is discussed.

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