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MS-17.01.07 NEW APPROACH TO REPRESENTATIONS OF SPACE GROUPS AND SUBPERIODIC GROUPS. By V. Kopský, Department of Physics, University of the South Pacific, POBox 1168, Suva, Fiji.

It has been recently found that subperiodic groups appear as factor groups of reducible space groups (Kopský, V. (1989). Acta Cryst. A45, 805-815 and 815-823). This has immediate and useful consequences in the theory of representations of both space and subperiodic groups analogous to known relationship between space groups and corresponding point groups. The point group G of a space group G is isomorphic to the factor group G/T_G where T_G is the full translation subgroup of G and there exists a homomorphism $\sigma: G \longrightarrow G$ with Ker $\sigma = T_G$. The lattice of equitranslational subgroups of G is isomorphic to the lattice of subgroups of G and representations of G engender those representations of G, the kernel of which contains T_G . These are exactly the representations which correspond to wavevector $\mathbf{k} = 0$ and the kernels are the equitranslational subgroups of G.

Let us now consider a reducible space group \mathcal{G} , the translation subgroup of which splits into a direct sum $T_G = T_{G1} \oplus T_{G2}$ of G- invariant subgroups $T_{G1} = T(\mathbf{a}, \mathbf{b})$ and $T_{G2} = T(\mathbf{c})$. These two groups are normal in \mathcal{G} and the corresponding factor groups $\mathcal{G}/T(\mathbf{c})$, $\mathcal{G}/T(\mathbf{a}, \mathbf{b})$ have the structure of a layer group \mathcal{L} with translation subgroup $T(\mathbf{c})$, respectively. Both the layer group \mathcal{L} and the rod group \mathcal{R} belong to the same geometric class G as the space group \mathcal{G} . There exist homomorphisms $\sigma_1: \mathcal{G} \longrightarrow \mathcal{L}$ with Ker $\sigma_1 = T(\mathbf{c})$ and $\sigma_2: \mathcal{G} \longrightarrow \mathcal{R}$ with Ker $\sigma_2 = T(\mathbf{a}, \mathbf{b})$.

The lattice of subgroups of the layer group \mathcal{L} is isomorphic to the lattice of those subgroups of \mathcal{G} which contain $T(\mathbf{c})$ and representations of \mathcal{L} engender those representations of \mathcal{G} , the kernel of which contains $T(\mathbf{c})$.

The lattice of subgroups of the rod group \mathcal{R} is isomorphic to the lattice of those subgroups of \mathcal{G} which contain $T(\mathbf{a}, \mathbf{b})$ and representations of \mathcal{R} engender those representations of \mathcal{G} , the kernel of which contains $T(\mathbf{a}, \mathbf{b})$.

The reciprocal space \widetilde{V} also splits into a direct sum of G-invariant subspaces $\widetilde{V}_1 = \widetilde{V}(\widetilde{\mathbf{a}}, \widetilde{\mathbf{b}})$ and $\widetilde{V}_2 = \widetilde{T}(\widetilde{\mathbf{c}})$. Representations of \mathcal{L} , \mathcal{R} and respective engendered representations of \mathcal{G} correspond to wavevectors $\mathbf{k}_1 \in \widetilde{V}_1$ and $\mathbf{k}_2 \in \widetilde{V}_2$.

The best way to record ireps (irreducible representations) is to give their kernels and mapping of cosets onto matrix images. Kernels of ireps generate by intersections the lattice of normal subgroups which contains implicitly important information about mode interactions (Kopský, V. (1988). Comput. Math. Applic. 16, 493-505). These lattices are especially simple for irreducible space and subperiodic groups (Kopský, V. (1987). Czech. J. Phys. B37, 785-808) from which the systematic theory and recording develops. They have been actually derived already time ago by Fuksa & Kopský but so far unpublished; it is necessary to develop first a standard system of symbols for space and subperiodic groups which will include the specification of the origin, of nonstandard orientation and of translation subgroup.

It would be desirable to develop correlated standards of ireps for space and subperiodic groups to replace existing systems which are not quite compatible. Standards of subperiodic groups (pending Vol E of International Tables for Crystallography) correlated with standards of space groups will provide necessary background for this project including system of Hermann-Mauguin symbols which will suit the purpose.

PS-17.01.08 PERIODIC CLOSE PACKINGS OF IDENTICAL ELLIPSES. By Takeo Matsumoto*, Department of Earth Sciences, Faculty of Science, Kanazawa University, Kakuma-machi, Kanazawa 920-11, Japan.

In contrast with the closest packing of circles, with plane group p6mm, packings of ellipses with 6 contacting neighbours do not always have the maximum density $p = \pi/2\sqrt{3}$. Nowacki (1948) indicated five different densest packings of ellipses in 4 different plane groups: p2, c2mm, p2gg and p31m. In all of these packings, every ellipse is in contact with 6 neighbour ellipses. These packings could be considered to be the densest among the possible forms of ellipse packing.

Matsumoto (1968) and Matsumoto and Nowacki (1966) have shown that the first two of the above densest packings of ellipses, p2 and c2mm, always attain the above maximum density. That of the third, p2gg, cannot exceed this maximum density. Namely, the densest p2 and c2mm packings of identical ellipses are derived from the densest packing of circles by affine transformation, while the p2gg packing of ellipses can never attain the above maximum density ρ.

Tanemura and Matsumoto (1992) have shown that the density of p31m packings of ellipses, the fourth one of the above list, never exceeds the above maximum density ρ . Grünbaum and Shephard (1987) have published a fascinating book, in which 58 types of periodic patterns of ellipses are listed. Among them, 54 types are due to Nowacki (1948), and the other four are new packings, three of them with 6 contacting neighbours. Whether or not the density of these packings is limited to the value $\pi/2\sqrt{3}$ still remains to be established.

PS-17.01.09 DEFECTS OF HETEROSTRUCTURES CAUSED BY THE SYMMETRY DISCREPANCY (AN EXAMPLE OF HTS-LAYER ON PEROV-SKITE-LIKE SUBSTRATES). By A.N. Efimov* and A.O. Lebedev, Ioffe Physical-Technical Institute, Petersburg, Russia. Types of structural defects in heterostructures "layer on non-isomorphic substrate", causes of their origin and their density as a function of techniques and crys-

Indexes of	Expected types of defects	Table 1 Parameters determining the density of defects
{111}	90° -twins (an angle between	density of nucleation
{001}	"c"-axes of twins is 45° -twins (axes "a" and "b" of replace each other)	density of nucleation twins
	coherent polysynthetic twins with interface on (110) or (110), incoherent interfaces between them.	
	antiphase domains	angular deflection of the substrate plane from singular one
{110} {hhl}	antiphase domains 45° -twins	density of nucleation density of nucleation
	coherent polysynthetic twins with interface on	
	(110) or (110)	on
	antiphase domains	density of nucleation and angular deflection
		of the substrate plane from (001)
{0k1}, {hk1}	antiphase domains	density of nucleation and angular deflection of the substrate plane from (001)

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tallographic constraints (Pond R.C., J. Crystal Growth, 1986,79,946; Efimov A.N., Thesis, Leningrad, 1986) are presented for the case of negligible metric misfit. As example an analysis of epitaxial defects was carried out for the structure "rhombic yttrium HTS 1-2-3 on SrTiO3". Result is shown in table 1. All these types of defects was observed experimentally by using electron microscope technique.

Possibilities of single crystalline growth for rhombic yttrium HTS -crystal 1-2-3 on substrate with structure of distorted perovskite (Fesenko E.G., Perovskite family and ferroelectricity, Moscow, Atomizdat, 1972) are illustrated by table 2 (only planes (001), (010), (100) of reduced unit cell are considered).

Table 2.

Symmetry of reduced perovskite (001) (010) (100) unit cell Cubic twinning twinning twinning Tetragonal twinning single crystal single crystal Rhombic 1 single crystal single crystal single crystal Rhombic 2 single crystal twinning* single crystal twinning* single crystal twinning* twinning* Rhombohedral twinning* Monoclinic single crystal twinning single crystal

* Polysynthetic twinning without incoherent interfaces is possible.

It is shown that defect-free heterocomposition should be created provided the translation groups of layer and substrate are in agreement meanwhile defects caused by the discrepancy between the point groups ought to be removed by the angular deflection of substrate plane from singular one.

PS-17.01.10 DETERMINATION OF LAUE SYMMETRY USING CONVERSE-TRANSFORMATION ANALYSIS. Vicky Lynn Karen* and Alan Mighell, Reactor Radiation Division, National Institute of Standards and Technology, Gaithersburg, MD 20899 USA.

Our newly-developed theory of Converse-Transformation analysis (J. Appl. Cryst. (1991) 24, 1076) permits symmetry, as a theoretical concept, to be abstracted directly from the experimental data. In practice, the Converse-Transformation matrices are calculated as soon as any unit cell has been determined. This new approach permits the collection of experimental data, the evaluation of experimental errors, the determination of metric symmetry and the assignment of the Laue symmetry with respect to ANY basis, no matter how skewed. The metric symmetry is deduced simply by counting; the greater the number of matrices, the higher the symmetry. All possible symmetries and pseudo-symmetries within any specified tolerance are immediately apparent. This is especially important in preliminary structure work and in protein crystallography where the experimental errors may be large. The Laue symmetry is assigned by using the Converse-Transformation matrices to generate sets of equivalent hkl's, then measuring and evaluating the intensities of the equivalent reflections. Mistakes in Laue symmetry are avoided since it is not necessary to do key experimental steps out of order. With this approach, all the data required to assign the Laue symmetry are collected before a conventional cell is determined. In contrast, procedures currently used in diffractometry are based on the risky practice of assuming a conventional cell and symmetry, and then collecting data to verify the assumption. This erroneous strategy may lead to the assignment of a Laue group of too low symmetry. This new method has many experimental applications including the automation of diffractometers and the editorial review of manuscripts prior to publication. Software for analyzing symmetry and for determining general lattice relationships may be obtained from Dr. Vicky Lynn Karen.

PS-17.01.11 FREE SPACE CRYSTALLOGRAPHY AND THE VOLUME TERM OF GIBBS ENERGY: COMPARISON OF THREE PATTERNS OF HEPTAGONS!

By D. Leuschner, Dresden, Germany

In many cases space filling is of interest in crystallography (Paufler, Leuschner 1975). In this connection the study of gaps is relevant. Here the free space in three patterns of heptagons (generalization from 2-D nets to 3-D lattices is also performed) is investigated. Geometrical and physical peculiarities are considered. For instance the effect in volume term of the Gibbs energy (Leuschner 1979, 1989) is estimated.