

4. DIFFUSE SCATTERING AND RELATED TOPICS

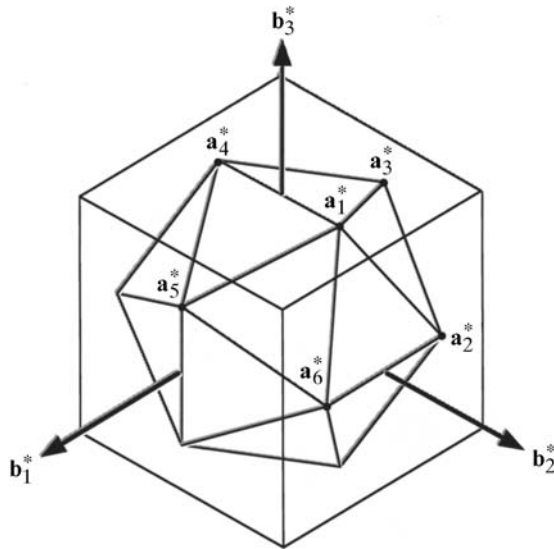


Fig. 4.6.3.32. Perspective parallel-space view of the two alternative reciprocal bases of the 3D Penrose tiling: the cubic and the icosahedral setting, represented by the bases \mathbf{b}_i^* , $i = 1, \dots, 3$, and \mathbf{a}_i^* , $i = 1, \dots, 6$, respectively.

basis vectors \mathbf{e}_i along the twofold axes) (Bancel *et al.*, 1985). The reciprocal basis is

$$\begin{pmatrix} \mathbf{a}_1^* \\ \mathbf{a}_2^* \\ \mathbf{a}_3^* \\ \mathbf{a}_4^* \\ \mathbf{a}_5^* \\ \mathbf{a}_6^* \end{pmatrix} = \frac{a^*}{(1 + \tau^2)^{1/2}} \begin{pmatrix} 0 & 1 & \tau \\ -1 & \tau & 0 \\ -\tau & 0 & 1 \\ 0 & -1 & \tau \\ \tau & 0 & 1 \\ 1 & \tau & 0 \end{pmatrix} \begin{pmatrix} \mathbf{e}_1^C \\ \mathbf{e}_2^C \\ \mathbf{e}_3^C \end{pmatrix}.$$

An alternate way of indexing is based on a 3D set of cubic reciprocal-basis vectors \mathbf{b}_i^* , $i = 1, \dots, 3$ (setting 2) (Fig. 4.6.3.32):

$$\begin{pmatrix} \mathbf{b}_1^* \\ \mathbf{b}_2^* \\ \mathbf{b}_3^* \end{pmatrix} = \frac{1}{2} \begin{pmatrix} 0 & \bar{1} & 0 & 0 & 0 & 1 \\ 1 & 0 & 0 & \bar{1} & 0 & 0 \\ 0 & 0 & 1 & 0 & 1 & 0 \end{pmatrix} \begin{pmatrix} \mathbf{a}_1^* \\ \mathbf{a}_2^* \\ \mathbf{a}_3^* \\ \mathbf{a}_4^* \\ \mathbf{a}_5^* \\ \mathbf{a}_6^* \end{pmatrix} \\ = \frac{a^*}{(1 + \tau^2)^{1/2}} \begin{pmatrix} \mathbf{e}_1^C \\ \mathbf{e}_2^C \\ \mathbf{e}_3^C \end{pmatrix}.$$

The Cartesian C basis is related to the V basis by a $\theta/2$ rotation around $[100]_C$, yielding $[001]_V$, followed by a $\pi/10$ rotation around $[001]_C$:

$$\begin{pmatrix} \mathbf{e}_1^C \\ \mathbf{e}_2^C \\ \mathbf{e}_3^C \end{pmatrix} = \begin{pmatrix} \cos(\pi/10) & \sin(\pi/10) & 0 \\ -\cos(\theta/2) \sin(\pi/10) & \cos(\theta/2) \cos(\pi/10) & \sin(\theta/2) \\ \sin(\theta/2) \sin(\pi/10) & -\sin(\theta/2) \cos(\pi/10) & \cos(\theta/2) \end{pmatrix} \begin{pmatrix} \mathbf{e}_1^V \\ \mathbf{e}_2^V \\ \mathbf{e}_3^V \end{pmatrix}.$$

Thus, indexing the diffraction pattern of an icosahedral phase with integer indices, one obtains for setting 1 $\mathbf{H} = \sum_{i=1}^6 h_i \mathbf{a}_i^*$, $h_i \in \mathbb{Z}$. These indices $(h_1 h_2 h_3 h_4 h_5 h_6)$ transform into the second setting to $(h/h' k/k' l/l')$ with the fractional cubic indices $h_1^c = h + h'\tau$, $h_2^c = k + k'\tau$, $h_3^c = l + l'\tau$. The transformation matrix is

$$\begin{pmatrix} h \\ h' \\ k \\ k' \\ l \\ l' \end{pmatrix}_C = \begin{pmatrix} 0 & \bar{1} & 0 & 0 & 0 & 1 \\ 0 & 0 & \bar{1} & 0 & 1 & 0 \\ 1 & 0 & 0 & \bar{1} & 0 & 0 \\ 0 & 1 & 0 & 0 & 0 & 1 \\ 0 & 0 & 1 & 0 & 1 & 0 \\ 1 & 0 & 0 & 1 & 0 & 0 \end{pmatrix} \begin{pmatrix} h_1 \\ h_2 \\ h_3 \\ h_4 \\ h_5 \\ h_6 \end{pmatrix}_D = \begin{pmatrix} h_6 - h_2 \\ h_5 - h_3 \\ h_1 - h_4 \\ h_6 + h_2 \\ h_5 + h_3 \\ h_1 + h_4 \end{pmatrix}_D.$$

4.6.3.3.3.2. Diffraction symmetry

The diffraction symmetry of icosahedral phases can be described by the Laue group $K = m\bar{3}5$. The set of all vectors \mathbf{H} forms a Fourier module $M^* = \{\mathbf{H}^{\parallel} = \sum_{i=1}^6 h_i \mathbf{a}_i^* | h_i \in \mathbb{Z}\}$ of rank 6 in physical space. Consequently, it can be considered as a projection from a 6D reciprocal lattice, $M^* = \pi^{\parallel}(\Sigma^*)$. The parallel and perpendicular reciprocal-space sections of the 3D Penrose tiling decorated with equal point scatterers on its vertices are shown in Figs. 4.6.3.33 and 4.6.3.34. The diffraction pattern in perpendicular space is the Fourier transform of the triacontahedron. All Bragg reflections within $10^{-4}|F(\mathbf{0})|^2 < |F(\mathbf{H})|^2 < |F(\mathbf{0})|^2$ are depicted. Without intensity-truncation limit, the diffraction pattern would be densely filled with discrete Bragg reflections.

The 6D icosahedral space groups that are relevant to the description of icosahedral phases (six symmorphic and five non-symmorphic groups) are listed in Table 4.6.3.2. These space groups are a subset of all 6D icosahedral space groups fulfilling the condition that the 6D point groups they are associated with are isomorphic to the 3D point groups $\frac{2}{m}\bar{3}5$ and 235 describing the diffraction symmetry. From 826 6D (analogues to) Bravais groups (Levitov & Rhyner, 1988), only three fulfil the condition that the projection of the 6D hypercubic lattice upon the 3D physical space is compatible with the icosahedral point groups $\frac{2}{m}\bar{3}5$, 235: the primitive hypercubic Bravais lattice P , the body-centred Bravais lattice I with translation $1/2(111111)$, and the face-centred Bravais lattice F with translations $1/2(110000) + 14$ further cyclic permutations. Hence, the I lattice is twofold primitive (*i.e.* it contains two vertices per unit cell) and the F lattice is 16-fold primitive. The orientation of the symmetry elements in the 6D space is defined by the isomorphism of the 3D and 6D point groups. The action of the fivefold rotation, however, is different in the subspaces \mathbf{V}^{\parallel} and \mathbf{V}^{\perp} : a rotation of $2\pi/5$ in \mathbf{V}^{\parallel} is correlated with a rotation of $4\pi/5$ in \mathbf{V}^{\perp} . The reflection and inversion operations are equivalent in both subspaces.

4.6.3.3.3.3. Structure factor

The structure factor of the icosahedral phase corresponds to the Fourier transform of the 6D unit cell,

$$F(\mathbf{H}) = \sum_{k=1}^N f_k(\mathbf{H}^{\parallel}) T_k(\mathbf{H}^{\parallel}, \mathbf{H}^{\perp}) g_k(\mathbf{H}^{\perp}) \exp(2\pi i \mathbf{H} \cdot \mathbf{r}_k),$$

with 6D diffraction vectors $\mathbf{H} = \sum_{i=1}^6 h_i \mathbf{d}_i^*$, parallel-space atomic scattering factor $f_k(\mathbf{H}^{\parallel})$, temperature factor $T_k(\mathbf{H}^{\parallel}, \mathbf{H}^{\perp})$, and perpendicular-space geometric form factor $g_k(\mathbf{H}^{\perp})$. $T_k(\mathbf{H}^{\parallel}, \mathbf{0})$ is equivalent to the conventional Debye–Waller factor and $T_k(\mathbf{0}, \mathbf{H}^{\perp})$ describes random fluctuations in perpendicular space. These fluctuations cause characteristic jumps of vertices (*phason flips*) in the physical space. Even random phason flips map the vertices onto positions that can still be described by physical-space vectors of the type $\mathbf{r} = \sum_{i=1}^6 n_i \mathbf{a}_i$. Consequently, the set $M = \{\mathbf{r} = \sum_{i=1}^6 n_i \mathbf{a}_i | n_i \in \mathbb{Z}\}$ of all possible vectors forms a \mathbb{Z} module. The shape of the atomic surfaces corresponds to a selection rule for the positions actually occupied. The geometric form factor $g_k(\mathbf{H}^{\perp})$ is equivalent to the Fourier transform of the *atomic surface*, *i.e.* the 3D perpendicular-space component of the 6D *hyperatoms*.

For the example of the canonical 3D Penrose tiling, $g_k(\mathbf{H}^{\perp})$ corresponds to the Fourier transform of a triacontahedron:

4.6. RECIPROCAL-SPACE IMAGES OF APERIODIC CRYSTALS

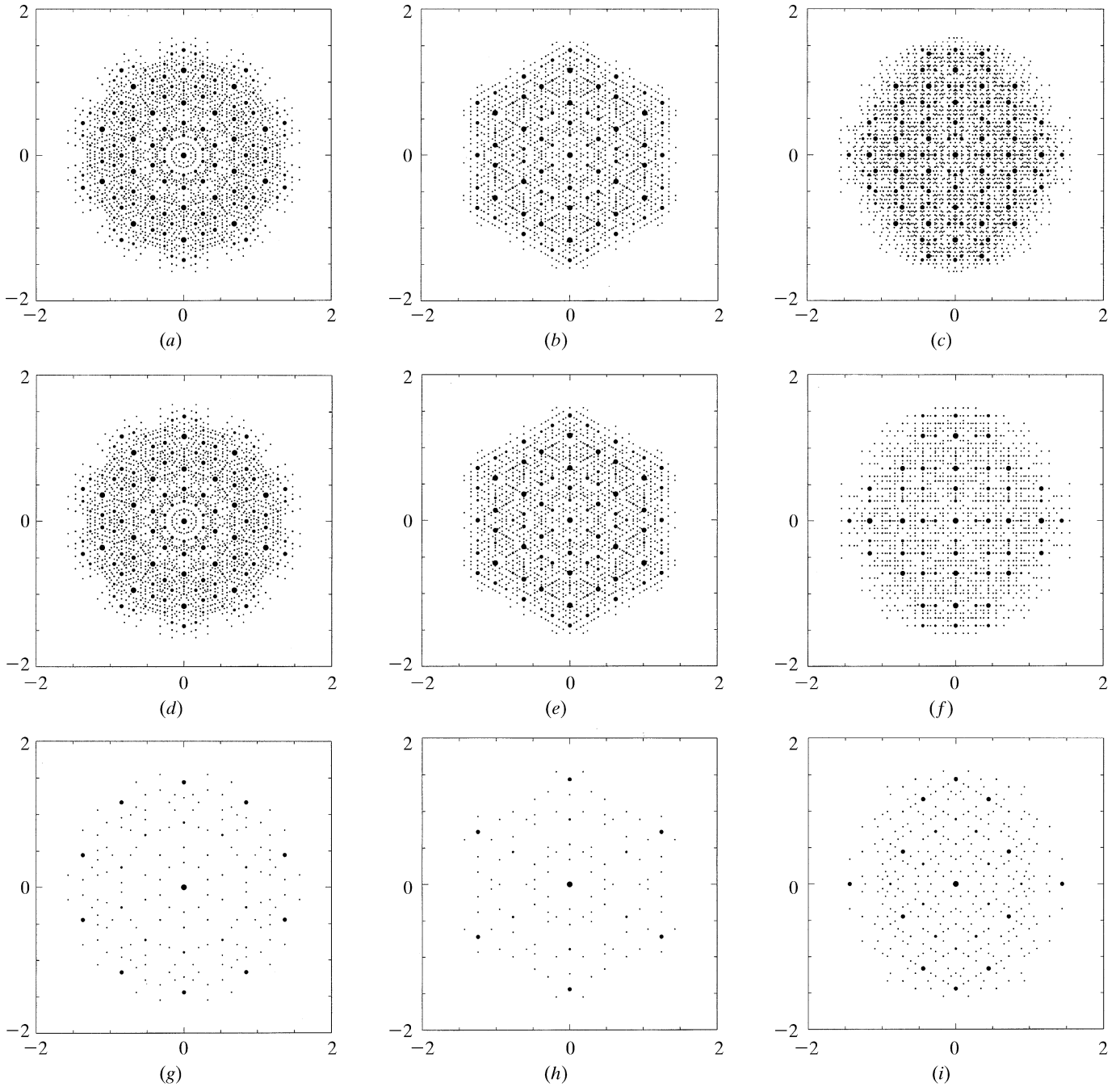


Fig. 4.6.333. Physical-space diffraction patterns of the 3D Penrose tiling decorated with point atoms (edge lengths of the Penrose unit rhombohedra $a_r = 5.0$ Å). Sections with five-, three- and twofold symmetry are shown for the primitive 6D analogue of Bravais type P in (a, b, c), the body-centred 6D analogue to Bravais type I in (d, e, f) and the face-centred 6D analogue to Bravais type F in (g, h, i). All reflections are shown within $10^{-4}|F(\mathbf{0})|^2 < |F(\mathbf{H})|^2 < |F(\mathbf{0})|^2$ and $-6 \leq h_i \leq 6, i = 1, \dots, 6$.

$$g_k(\mathbf{H}^\perp) = (1/A_{\text{UC}}^\perp) \int_{A_k} \exp(2\pi i \mathbf{H}^\perp \cdot \mathbf{r}) \, d\mathbf{r},$$

where A_{UC}^\perp is the volume of the 6D unit cell projected upon \mathbf{V}^\perp and A_k is the volume of the triacontahedron. A_{UC}^\perp and A_k are equal in the present case and amount to the volumes of ten prolate and ten oblate rhombohedra: $A_{\text{UC}}^\perp = 8a_r^3 [\sin(2\pi/5) + \sin(\pi/5)]$. Evaluating the integral by decomposing the triacontahedron into trigonal pyramids, each one directed from the centre of the triacontahedron to three of its corners given by the vectors $\mathbf{e}_i, i = 1, \dots, 3$, one obtains

$$g(\mathbf{H}^\perp) = (1/A_{\text{UC}}^\perp) \sum_R g_k(R^T \mathbf{H}^\perp),$$

with $k = 1, \dots, 60$ running over all site-symmetry operations R of the icosahedral group,

$$g_k(\mathbf{H}^\perp) = -iV_r [A_2 A_3 A_4 \exp(iA_1) + A_1 A_3 A_5 \exp(iA_2) + A_1 A_2 A_6 \exp(iA_3) + A_4 A_5 A_6] \times (A_1 A_2 A_3 A_4 A_5 A_6)^{-1},$$

$A_j = 2\pi \mathbf{H}^\perp \cdot \mathbf{e}_j, j = 1, \dots, 3, A_4 = A_2 - A_3, A_5 = A_3 - A_1, A_6 = A_1 - A_2$ and $V_r = \mathbf{e}_1 \cdot (\mathbf{e}_2 \times \mathbf{e}_3)$ the volume of the parallelepiped defined by the vectors $\mathbf{e}_i, i = 1, \dots, 3$ (Yamamoto, 1992b).

4.6.3.3.4. Intensity statistics

The radial structure-factor distributions of the 3D Penrose tiling decorated with point scatterers are plotted in Fig. 4.6.335 as a function of parallel and perpendicular space. The distribution of $|F(\mathbf{H})|$ as a function of their frequencies clearly resembles a centric