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## Vector algebra and crystallography

JEAN SIVARDIÈRE at CEA/Département de Recherche Fondamentale sur la Matière Condensée/MRS, 38054 Grenoble CEDEX 9, France

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## Abstract

The volumes of a crystal unit cell and of its reciprocal cell, the relationships between direct and reciprocal interaxial angles and the coordinates of the reciprocal vectors in the direct basis are rederived in a concise way by means of an elementary formula in vector algebra. The product of two rotations is also considered.

The volumes of the direct and reciprocal cells of a crystal and the relationships between direct- and reciprocal-lattice quantities are classically derived using the algebra of determinants and spherical crystallography or advanced vector methods such as Lagrange formulas for products of four vectors. We present here a simple vector formula from which all these quantities are then rederived. Finally, the relationships between direct- and reciprocal-lattice quantities are applied to spherical trigonometry and the product of two rotations.

Consider three linearly independent vectors  $\mathbf{a}$ ,  $\mathbf{b}$ ,  $\mathbf{c}$  and introduce the coordinates x, y, z of  $\mathbf{c}$  in the  $\mathbf{a}$ ,  $\mathbf{b}$ ,  $\mathbf{a} \times \mathbf{b}$  basis:

$$\mathbf{c} = x\mathbf{a} + y\mathbf{b} + z\mathbf{a} \times \mathbf{b}$$
.

Taking the scalar product of this relation successively by  $\mathbf{a}$ ,  $\mathbf{b}$  and  $\mathbf{a} \times \mathbf{b}$ , we find:

$$\mathbf{a} \cdot \mathbf{c} = xa^2 + y\mathbf{a} \cdot \mathbf{b}$$
$$\mathbf{b} \cdot \mathbf{c} = x\mathbf{a} \cdot \mathbf{b} + yb^2$$
$$(\mathbf{a}, \mathbf{b}, \mathbf{c}) = z|\mathbf{a} \times \mathbf{b}|^2.$$

Using the Lagrange identity

$$|\mathbf{a} \times \mathbf{b}|^2 = a^2b^2 - (\mathbf{a} \cdot \mathbf{b})^2$$

we get finally the 'basic' formula

$$|\mathbf{a} \times \mathbf{b}|^2 \mathbf{c} = [b^2 (\mathbf{a} \cdot \mathbf{c}) - (\mathbf{a} \cdot \mathbf{b})(\mathbf{b} \cdot \mathbf{c})]\mathbf{a} + [a^2 (\mathbf{b} \cdot \mathbf{c}) - (\mathbf{a} \cdot \mathbf{b})(\mathbf{a} \cdot \mathbf{c})]\mathbf{b} + (\mathbf{a}, \mathbf{b}, \mathbf{c})\mathbf{a} \times \mathbf{b}. \quad (1)$$

(a) Suppose that the three vectors  $\mathbf{a}$ ,  $\mathbf{b}$ ,  $\mathbf{c}$  define a crystal unit cell: the cell edges are a, b, c and the interaxial angles are  $\alpha$ ,  $\beta$ ,  $\gamma$ . The volume V of the cell is equal to the triple scalar product  $(\mathbf{a}, \mathbf{b}, \mathbf{c}) = \mathbf{a} \times \mathbf{b} \cdot \mathbf{c} = \mathbf{a} \cdot \mathbf{b} \times \mathbf{c}$ . Taking the scalar product of (1) by  $\mathbf{c}$ , we get the classical expression (Buerger, 1942; Carpenter, 1969; Neustadt & Cagle, 1968; Woolfson, 1970).

$$v^{2} = a^{2}b^{2}c^{2}(1 + 2\cos\alpha\cos\beta\cos\gamma - \cos^{2}\alpha - \cos^{2}\beta - \cos^{2}\gamma),$$
(2)

which is generally derived using the theory of determinants. An interesting equivalent expression is (Donnay & Donnay, 1959)

$$v^2 = 4a^2b^2c^2[\sin s\sin(s-\alpha)\sin(s-\beta)\sin(s-\gamma)]$$

with  $2s = \alpha + \beta + \nu$ .

(b) Consider now the reciprocal vectors  $\mathbf{a}^*$ ,  $\mathbf{b}^*$ ,  $\mathbf{c}^*$ . We want to determine the reciprocal-cell edges  $a^*$ ,  $b^*$ ,  $c^*$ , the interaxial angles  $\alpha^*$ ,  $\beta^*$ ,  $\gamma^*$  and the coordinates of the reciprocal vectors in the  $\mathbf{a}$ ,  $\mathbf{b}$ ,  $\mathbf{c}$  basis. Since  $v\mathbf{c}^* = \mathbf{a} \times \mathbf{b}$ , we have  $c^* = ab \sin \gamma/v$ . Equation (1) gives immediately the expression for  $c^*$  in the direct basis:

$$v^{2}\mathbf{c}^{*} = [(\mathbf{a} \cdot \mathbf{b})(\mathbf{b} \cdot \mathbf{c}) - b^{2}(\mathbf{a} \cdot \mathbf{c})]\mathbf{a}$$
$$+ [(\mathbf{a} \cdot \mathbf{b})(\mathbf{a} \cdot \mathbf{c}) - a^{2}(\mathbf{b} \cdot \mathbf{c})]\mathbf{b} + |\mathbf{a} \times \mathbf{b}|^{2}\mathbf{c}$$
(3)

or

$$v^{2}\mathbf{c}^{*} = ab^{2}c(\cos\alpha\cos\gamma - \cos\beta)\mathbf{a}$$
$$+ a^{2}bc(\cos\beta\cos\gamma - \cos\alpha)\mathbf{b} + a^{2}b^{2}\sin^{2}\gamma\mathbf{c}. \quad (3')$$

Comparing this with the general expression

$$\mathbf{c}^* = (\mathbf{a}^* \cdot \mathbf{c}^*)\mathbf{a} + (\mathbf{b}^* \cdot \mathbf{c}^*)\mathbf{b} + (\mathbf{c}^* \cdot \mathbf{c}^*)\mathbf{c} \tag{4}$$

and equating the coefficients of  $\mathbf{b}$  in (3) and (4), we get the angle  $\alpha^*$ :

$$\cos \alpha^* = (\cos \beta \cos \gamma - \cos \alpha) / \sin \beta \sin \gamma \tag{5}$$

Combining (3') and (5) gives

$$v^{2}\mathbf{c}^{*} = a^{2}bc\sin\alpha\sin\gamma\cos\beta^{*}\mathbf{a}$$
$$+ab^{2}c\sin\beta\sin\gamma\cos\alpha^{*}\mathbf{b} + a^{2}b^{2}\sin^{2}\gamma\mathbf{c}. \tag{6}$$

Calculating  $\sin \alpha^*$  from (5) and (2) gives

$$v = abc \sin \beta \sin \gamma \sin \alpha^* \tag{7}$$

and the sine relation

$$\frac{\sin \alpha}{\sin \alpha^*} = \frac{\sin \beta}{\sin \beta^*} = \frac{\sin \gamma}{\sin \gamma^*}.$$
 (8)

The volume  $\nu^*$  of the reciprocal cell is given by an expression similar to (7) and, from (8), one gets  $\nu\nu^* = 1$ .

(c) We translate now some of the preceding results into the language of spherical trigonometry.  $\cos \gamma^*$  is given by a relation similar to (5) and the reciprocal relation is

$$\cos \gamma^* = \cos \alpha^* \cos \beta^* - \sin \alpha^* \sin \beta^* \cos \gamma. \tag{9}$$

Using (7) in the form  $v = abc \sin \alpha \sin \beta \sin \gamma^*$  and (8), we may write (6) as

$$abc \sin \alpha^* \sin \beta^* \mathbf{a} \times \mathbf{b} = a^2 bc \sin \alpha^* \cos \beta^* \mathbf{a} + ab^2 c \sin \beta^* \cos \alpha^* \mathbf{b} + a^2 b^2 \sin \gamma^* \mathbf{c}.$$
(10)

Suppose that the extremities A, B and C of the three vectors a, b, c are on a unit sphere centered at the origin:

© 1996 International Union of Crystallography Printed in Great Britain – all rights reserved Acta Crystallographica Section A ISSN 0108-7673 © 1996 a=b=c=1. The side lengths of the spherical triangle are equal to the interaxial angles  $\alpha$ ,  $\beta$ ,  $\gamma$  and its dihedral angles are  $\pi-\alpha^*$ ,  $\pi-\beta^*$ ,  $\pi-\gamma^*$ . Using now the classical notations a, b, c for the side lengths and  $\alpha$ ,  $\beta$ ,  $\gamma$  for the dihedral angles, we get from (9) and (10), respectively,

$$\cos \gamma = -\cos \alpha \cos \beta + \sin \alpha \sin \beta \cos c \tag{11}$$

 $\sin \gamma \mathbf{c} = \sin \alpha \cos \beta \mathbf{a} + \cos \alpha \sin \beta \mathbf{b} + \sin \alpha \sin \beta \mathbf{a} \times \mathbf{b}.$ 

(12)

These formulas were found by Altmann by a more sophisticated procedure (Altmann, 1986). They can be used to derive the product of two rotations around intersecting axes from the Euler-Rodrigues-Hamilton theorem (Altmann, 1986; Sivardière, 1994, 1995). According to this theorem, a rotation of angle  $\theta_1 = 2\alpha$  around  $\mathbf{u}_1 = \mathbf{a}$  followed by a rotation of angle  $\theta_2 = 2\beta$  around  $\mathbf{u}_2 = \mathbf{b}$  is a rotation of angle  $\theta_3 = 2(\pi - \gamma)$  around the unit vector  $\mathbf{u}_3 = \mathbf{c}$ . Introducing the Euler vectors  $\mathbf{R}_i = \sin(\theta_i/2) \mathbf{u}_i$ , we obtain

$$\cos(\theta_3/2) = \cos(\theta_1/2)\cos(\theta_2/2) - \mathbf{R}_1 \cdot \mathbf{R}_2$$
$$\mathbf{R}_3 = \cos(\theta_2/2) \mathbf{R}_1 + \cos(\theta_1/2) \mathbf{R}_2 - \mathbf{R}_1 \times \mathbf{R}_2.$$

## References

Altmann, S. L. (1986). Rotations, Quaternions and Double Groups. Oxford: Clarendon Press.

Buerger, M. J. (1942). X-ray Crystallography. New York: John Wiley.

Carpenter, G. B. (1969). Principles of Crystal Structure Determination. New York: Benjamin.

Donnay, J. D. H. & Donnay, G. (1959). In *International Tables for X-ray Crystallography*, Vol. II. Birmingham: Kynoch Press.

Neustadt, R. J. & Cagle, F. W. (1968). Acta Cryst. A24, 247-248.

Sivardière, J. (1994). Am. J. Phys. 62, 737-743.

Sivardière, J. (1995). La Symétrie en Mathématiques, Physique et Chimie. Grenoble: Presses Universitaires de Grenoble.

Woolfson, M. M. (1970). An Introduction to X-ray Crystallography. Cambridge University Press.