

3.1. STRUCTURAL PHASE TRANSITIONS

Table 3.1.3.1 (cont.)

R-irep $\Gamma_\eta$	Standard variables	Ferroic symmetry			Principal tensor parameters	Domain states		
		$F_1$	$n_F$			$n_f$	$n_a$	$n_e$
<b>Parent symmetry G: <math>6_2 2_x 2_y D_6</math></b>								
$A_2$	$X_2$	$6_z$	$C_6$	1	$P_3$	2	1	2
$B_1$	$X_3$	$3_z 2_x$	$D_{3x}$	1	$d_{11} = -d_{12} = -d_{26}$	2	1	0
$B_2$	$X_4$	$3_z 2_y$	$D_{3y}$	1	$d_{22} = -d_{21} = -d_{16}$	2	1	0
$E_2$	$(x_2, 0)$	$2_x 2_y 2_z$	$D_2$	3	$\delta u_1 = -\delta u_2$	3	3	0
(La, Li)	$(x_2, y_2)$	$2_z$	$C_{2z}$	1	$(u_1 - u_2, 2u_6)$	6	6	2
$E_1$	$(x_1, 0)$	$2_x$	$C_{2x}$	3	$P_1; u_4$	6	6	6
	$(0, y_1)$	$2_y$	$C_{2y}$	3	$P_2; u_5$	6	6	6
(Li)	$(x_1, y_1)$	1	$C_1$	1	$(P_1, P_2); (u_4, -u_5)$	12	12	12
<b>Parent symmetry G: <math>6_2 m_x m_y C_{6v}</math></b>								
$A_2$	$X_2$	$6_z$	$C_6$	1	$\varepsilon; g_1 = g_2, g_3; d_{14} = -d_{25}$	2	1	1
$B_2$	$X_3$	$3_z m_x$	$C_{3vx}$	1	$d_{22} = -d_{21} = -d_{16}$	2	1	1
$B_1$	$X_4$	$3_z m_y$	$C_{3vy}$	1	$d_{11} = -d_{12} = -d_{26}$	2	1	1
$E_2$	$(x_2, 0)$	$m_x m_y 2_z$	$C_{2vz}$	3	$\delta u_1 = -\delta u_2$	3	3	1
(La)	$(x_2, y_2)$	$2_z$	$C_{2z}$	1	$(u_1 - u_2, 2u_6)$	6	6	1
$E_1$	$(x_1, 0)$	$m_x$	$C_{sx}$	3	$P_2; u_4$	6	6	6
	$(0, y_1)$	$m_y$	$C_{sy}$	3	$P_1; u_5$	6	6	6
	$(x_1, y_1)$	1	$C_1$	1	$(P_2, -P_1); (u_4, -u_5)$	12	12	12
<b>Parent symmetry G: <math>\bar{6}_2 2_x m_y D_{3h}</math></b>								
$A'_2$	$X_2$	$\bar{6}_z$	$C_{3h}$	1	$d_{22} = -d_{21} = -d_{16}$	2	1	0
$A''_1$	$X_3$	$3_z 2_x$	$D_{3x}$	1	$\varepsilon; g_1 = g_2, g_3; d_{14} = -d_{25}$	2	1	0
$A''_2$	$X_4$	$3_z m_y$	$C_{3vy}$	1	$P_3$	2	1	2
$E'$	$(x_2, 0)$	$2_x m_y m_z$	$C_{2vx}$	3	$P_1; \delta u_1 = -\delta u_2$	3	3	3
(La)	$(x_2, y_2)$	$m_z$	$C_{sz}$	1	$(P_1, -P_2); (u_1 - u_2, 2u_6)$	6	6	6
$E''$	$(x_1, 0)$	$2_x$	$C_{2x}$	3	$u_4$	6	6	3
	$(0, y_1)$	$m_y$	$C_{sy}$	3	$u_5$	6	6	6
	$(x_1, y_1)$	1	$C_1$	1	$(u_4, -u_5)$	12	12	12
<b>Parent symmetry G: <math>\bar{6}_2 m_x 2_y \bar{D}_{3h}</math></b>								
$A'_2$	$X_2$	$\bar{6}_z$	$C_{3h}$	1	$d_{11} = -d_{12} = -d_{26}$	2	1	0
$A''_2$	$X_3$	$3_z m_x$	$C_{3vx}$	1	$P_3$	2	1	2
$A'_1$	$X_4$	$3_z 2_y$	$D_{3y}$	1	$\varepsilon; g_1 = g_2, g_3; d_{14} = -d_{25}$	2	1	0
$E'$	$(x_2, 0)$	$m_x 2_y m_z$	$C_{2vy}$	3	$P_2; \delta u_1 = -\delta u_2$	3	3	3
(La)	$(x_2, y_2)$	$m_z$	$C_{sz}$	1	$(P_2, P_1); (u_1 - u_2, 2u_6)$	6	6	6
$E''$	$(x_1, 0)$	$m_x$	$C_{sx}$	3	$u_4$	6	6	6
	$(0, y_1)$	$2_y$	$C_{2y}$	3	$u_5$	6	6	3
	$(x_1, y_1)$	1	$C_1$	1	$(u_4, -u_5)$	12	12	12
<b>Parent symmetry G: <math>6_z/m_x m_y D_{6h}</math></b>								
$A_{2g}$	$X_2^+$	$6_z/m_z$	$C_{6h}$	1	$A_{31} = A_{32}, A_{33}, A_{15} = A_{24}$	2	1	0
$B_{1g}$	$X_3^+$	$3_z m_x$	$D_{3dx}$	1	$A_{11} = -A_{12} = -A_{26}$	2	1	0
$B_{2g}$	$X_4^+$	$3_z m_y$	$D_{3dy}$	1	$A_{22} = -A_{21} = -A_{16}$	2	1	0
$A_{1u}$	$X_1^-$	$6_2 2_x 2_y$	$D_6$	1	$\varepsilon; g_1 = g_2, g_3; d_{14} = -d_{25}$	2	1	0
$A_{2u}$	$X_2^-$	$6_z m_x m_y$	$C_{6v}$	1	$P_3$	2	1	2
$B_{1u}$	$X_3^-$	$6_z 2_x m_y$	$D_{3h}$	1	$d_{11} = -d_{12} = -d_{26}$	2	1	0
$B_{2u}$	$X_4^-$	$6_z m_x 2_y$	$\bar{D}_{3h}$	1	$d_{22} = -d_{21} = -d_{16}$	2	1	0
$E_{2g}$	$(x_2^+, 0)$	$m_x m_y m_z$	$D_{2h}$	3	$\delta u_1 = -\delta u_2$	3	3	0
(La)	$(x_2^+, y_2^+)$	$2_z/m_z$	$C_{2hz}$	1	$(u_1 - u_2, 2u_6)$	6	6	0
$E_{1g}$	$(x_1^+, 0)$	$2_x/m_x$	$C_{2hx}$	3	$u_4$	6	6	0
	$(0, y_1^+)$	$2_y/m_y$	$C_{2hy}$	3	$u_5$	6	6	0
	$(x_1^+, y_1^+)$	1	$C_i$	1	$(u_4, -u_5)$	12	12	0
$E_{1u}$	$(x_1^-, 0)$	$2_x m_y m_z$	$C_{2vx}$	3	$P_1$	6	3	6
	$(0, y_1^-)$	$m_x 2_y m_z$	$C_{2vy}$	3	$P_2$	6	3	6
	$(x_1^-, y_1^-)$	$m_z$	$C_{sz}$	1	$(P_1, P_2)$	12	6	12
$E_{2u}$	$(x_2^-, 0)$	$2_x 2_y 2_z$	$D_2$	3	$\delta g_1 = -\delta g_2; d_{36}, \delta d_{14} = \delta d_{25}$	6	3	0
	$(0, y_2^-)$	$m_x m_y 2_z$	$C_{2vz}$	3	$g_6; d_{32} = -d_{31}, d_{24} = -d_{15}$	6	3	2
	$(x_2^-, y_2^-)$	$2_z$	$C_{2z}$	1	$(g_1 - g_2, 2g_6); (2d_{36}, d_{32} - d_{31}), (d_{14} + d_{25}, d_{24} - d_{15})$	12	6	2

In tensor distinction of domains, the secondary tensor parameters play a secondary role in a sense that some but not all ferroic domain states exhibit different values of the secondary tensor parameters. This property forms a basis for the concept of partial ferroic phases (Aizu, 1970): A ferroic phase is a *partial ferroelectric (ferroelastic)* one if some but not all domain states differ in spontaneous polarization (spontaneous strain). A non-ferroelectric phase denotes a ferroic phase which is either non-polar or which possesses a unique polar direction available

already in the parent phase. A non-ferroelastic phase exhibits no spontaneous strain.

3.1.3.3. Tables of equitranslational phase transitions associated with irreducible representations

The first systematic symmetry analysis of Landau-type phase transitions was performed by Indenbom (1960), who found all equitranslational phase transitions that can be accomplished

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Table 3.1.3.1 (cont.)

(g) Cubic parent groups

Covariants with standardized labels and conversion equations:

$$\begin{aligned}
 u_{3x} &= u_{3x}^+ = u_3 - a(u_1 + u_2); & u_{3y} &= u_{3y}^+ = b(u_1 - u_2) \\
 \delta u_1 &= -\frac{1}{3}u_{3x}^+ + \frac{1}{\sqrt{3}}u_{3y}^+; & \delta u_2 &= -\frac{1}{3}u_{3x}^+ - \frac{1}{\sqrt{3}}u_{3y}^+; & \delta u_3 &= \frac{2}{3}u_{3x}^+ \\
 g_1^- &= g_1 + g_2 + g_3; & g_{3x}^- &= g_3 - a(g_1 + g_2); & g_{3y}^- &= b(g_1 - g_2) \\
 g_1 &= \frac{1}{3}g_1^- - \frac{1}{3}g_{3x}^- + \frac{1}{\sqrt{3}}g_{3y}^-; & g_2 &= \frac{1}{3}g_1^- - \frac{1}{3}g_{3x}^- - \frac{1}{\sqrt{3}}g_{3y}^-; & g_3 &= \frac{1}{3}g_1^- + \frac{2}{3}g_{3x}^- \\
 d_1^- &= d_{14} + d_{25} + d_{36}; & d_{3x}^- &= b(d_{14} - d_{25}), & d_{3y}^- &= a(d_{14} + d_{25}) - d_{36} \\
 d_{14} &= \frac{1}{3}d_1^- + \frac{1}{\sqrt{3}}d_{3x}^- + \frac{1}{3}d_{3y}^-; & d_{25} &= \frac{1}{3}d_1^- - \frac{1}{\sqrt{3}}d_{3x}^- + \frac{1}{3}d_{3y}^-; & d_{36} &= \frac{1}{3}d_1^- - \frac{2}{3}d_{3y}^- \\
 d_{1x} &= d_{13} - d_{12}; & d_{1y} &= d_{21} - d_{23}; & d_{1z} &= d_{32} - d_{31} \\
 d_{2x} &= d_{13} + d_{12}; & d_{2y} &= d_{21} + d_{23}; & d_{2z} &= d_{32} + d_{31} \\
 d_{13} &= \frac{1}{2}(d_{1x} + d_{2x}); & d_{21} &= \frac{1}{2}(d_{1y} + d_{2y}); & d_{32} &= \frac{1}{2}(d_{1z} + d_{2z}) \\
 d_{12} &= \frac{1}{2}(d_{2x} - d_{1x}); & d_{23} &= \frac{1}{2}(d_{2y} - d_{1y}); & d_{31} &= \frac{1}{2}(d_{2z} - d_{1z})
 \end{aligned}$$

$$a = \frac{1}{2}, b = \frac{\sqrt{3}}{2}, \pi_{\mu\nu}^a = (\pi_{\mu\nu} - \pi_{\nu\mu}), \mu = 1, 2, \dots, 6, \nu = 1, 2, \dots, 6.$$

R-irep $\Gamma_\eta$	Standard variables	Ferroic symmetry			Principal tensor parameters	Domain states		
		$F_1$		$n_F$		$n_f$	$n_a$	$n_e$
<b>Parent symmetry G: 23 T</b>								
$E$ (La)	$(x_3, y_3)$	$2_x 2_y 2_z$	$D_2$	1	$[u_3 - a(u_1 + u_2), b(u_1 - u_2)]$ $\delta u_1 + \delta u_2 + \delta u_3 = 0$	3	3	0
$T$ (La, Li)	$(0, 0, z_1)$ $(x_1, x_1, x_1)$ $(x_1, y_1, z_1)$	$2_z$ $3_p$ 1	$C_{2z}$ $C_{3p}$ $C_1$	3 4 1	$P_3; u_6$ $P_1 = P_2 = P_3; u_4 = u_5 = u_6$ $(P_1, P_2, P_3); (u_4, u_5, u_6)$	6 4 12	6 4 12	6 4 12
<b>Parent symmetry G: <math>m\bar{3} T_h</math></b>								
$A_u$	$x_1^-$	23	$T$	1	$\varepsilon; g_1 = g_2 = g_3; d_{14} = d_{25} = d_{36}$	2	1	0
$E_g$ (La)	$(x_3^+, y_3^+)$	$m_x m_y m_z$	$D_{2h}$	1	$[u_3 - a(u_1 + u_2), b(u_1 - u_2)]$ $\delta u_1 + \delta u_2 + \delta u_3 = 0$	3	3	0
$E_u$	$(x_3^-, y_3^-)$	$2_x 2_y 2_z$	$D_2$	1	$[g_3 - a(g_1 + g_2), b(g_1 - g_2)]$ $\delta g_1 + \delta g_2 + \delta g_3 = 0$ $[b(d_{14} - d_{25}), a(d_{14} + d_{25}) - d_{36}]$ $\delta d_{14} + \delta d_{25} + \delta d_{36} = 0$	6	3	0
$T_g$ (La)	$(0, 0, z_1^+)$ $(x_1^+, x_1^+, x_1^+)$ $(x_1^+, y_1^+, z_1^+)$	$2_z/m_z$ $\bar{3}_p$ 1	$C_{2hz}$ $C_{3p}$ $C_i$	3 4 1	$u_6$ $u_4 = u_5 = u_6$ $(u_4, u_5, u_6)$	6 4 12	6 4 12	0 0 0
$T_u$	$(0, 0, z_1^-)$ $(x_1^-, x_1^-, x_1^-)$ $(x_1^-, y_1^-, z_1^-)$	$m_x m_y 2_z$ $3_p$ 1	$C_{2vz}$ $C_{3p}$ $C_1$	3 4 1	$P_3$ $P_1 = P_2 = P_3$ $(P_1, P_2, P_3)$	6 8 24	3 4 12	6 8 24
<b>Parent symmetry G: 432 O</b>								
$A_2$	$x_2$	23	$T$	1	$d_{14} = d_{25} = d_{36}$	2	1	0
$E$ (La)	$(x_3, 0)$ $(x_3, y_3)$	$4_z 2_x 2_{xy}$ $2_x 2_y 2_z$	$D_{4z}$ $D_2$	3 1	$\delta u_1 = \delta u_2 = -\frac{1}{2}\delta u_3$ $[u_3 - a(u_1 + u_2), b(u_1 - u_2)]$ $\delta u_1 + \delta u_2 + \delta u_3 = 0$	3 6	3 6	0 0
$T_1$ (Li)	$(0, 0, z_1)$ $(x_1, x_1, 0)$ $(x_1, x_1, x_1)$ $(x_1, y_1, z_1)$	$4_z$ $2_{xy}$ $3_p$ 1	$C_{4z}$ $C_{2xy}$ $C_{3p}$ $C_1$	3 6 4 1	$P_3$ $P_1 = P_2$ $P_1 = P_2 = P_3$ $(P_1, P_2, P_3)$	6 12 8 24	3 12 4 24	6 12 8 24
$T_2$ (La, Li)	$(0, 0, z_2)$ $(x_2, -x_2, z_2)$ $(x_2, x_2, x_2)$ $(x_2, y_2, z_2)$	$2_{x\bar{y}} 2_{xy} 2_z$ $2_{xy}$ $3_p 2_{x\bar{y}}$ 1	$\hat{D}_{2z}$ $C_{2xy}$ $D_{3p}$ $C_1$	3 6 4 1	$u_6$ $u_4 = -u_5, u_6$ $u_4 = u_5 = u_6$ $(u_4, u_5, u_6)$	6 12 4 24	6 12 4 24	0 12 0 24

continuously. A table of all crystallographic point groups  $G$  along with all their physically irreducible representations, corresponding ferroic point groups  $F$  and related data has been compiled by Janovec *et al.* (1975). These data are presented in an improved form in Table 3.1.3.1 together with corresponding principal tensor parameters and numbers of ferroic, ferroelectric and ferroelastic domain states. This table facilitates solving of the following typical problems:

(1) *Inverse Landau problem* (Ascher & Kobayashi, 1977) of equitranslational phase transitions: For a given equitranslational symmetry descent  $\mathcal{G} \Downarrow^t \mathcal{F}$  (determined for example from diffraction experiments), find the representation  $\Gamma_\eta$  of  $\mathcal{G}$  that specifies the transformation properties of the primary order parameter. Solution: In Table 3.1.3.1, one finds a physically irre-

ducible representation  $\Gamma_\eta$  of the point group  $G$  of  $\mathcal{G}$  with epikernel  $F$  (point group of  $\mathcal{F}$ ). For some symmetry descents from cubic point groups  $G = 432, 43m$  and  $m\bar{3}m$ , the inverse Landau problem has two solutions, which are given in Table 3.1.3.2.

If for a given symmetry descent  $\mathcal{G} \Downarrow^t \mathcal{F}$  no appropriate  $R$ -irep exists in Table 3.1.3.1, then the primary order parameter  $\eta$  transforms according to a reducible representation of  $G$ . These transitions are always discontinuous and can be accomplished with several reducible representations. Some symmetry descents can be associated with an irreducible representation and with several reducible representations. All these transitions are treated in the software *GI★KobO-1* and in Kopský (2001). All point-group symmetry descents are listed in Table 3.4.2.7 and can be traced in lattices of subgroups (see Figs. 3.1.3.1 and 3.1.3.2).

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Table 3.1.3.1 (cont.)

$R$ -irep $\Gamma_\eta$	Standard variables	Ferroic symmetry			Principal tensor parameters	Domain states		
		$F_1$		$n_F$		$n_f$	$n_a$	$n_e$
<b>Parent symmetry <math>G</math>: <math>\bar{4}3m</math> <math>T_d</math></b>								
$A_2$	$x_2$	23	$T$	1	$\varepsilon; g_1 = g_2 = g_3$ $A_{14} = A_{25} = A_{36}; \pi_{23}^a = \pi_{31}^a = \pi_{12}^a$	2	1	0
$E$ (La)	$(x_3, 0)$	$\bar{4}_z 2_x m_{xy}$	$D_{2dz}$	3	$\delta u_1 = \delta u_2 = -\frac{1}{2} \delta u_3$ $[u_3 - a(u_1 + u_2), b(u_1 - u_2)]$ $\delta u_1 + \delta u_2 + \delta u_3 = 0$	3	3	0
	$(x_3, y_3)$	$2_x 2_y 2_z$	$D_2$	1				
$T_1$	$(0, 0, z_1)$	$\bar{4}_z$	$S_{4z}$	3	$g_6; d_{32} = -d_{31}, d_{24} = -d_{15}$ $g_4 = g_5$ $d_{13} = -d_{23}, d_{12} = -d_{21}$ $d_{35} = -d_{34}, d_{26} = -d_{16}$ $g_4 = g_5 = g_6$ $d_{13} = d_{21} = d_{32}, d_{12} = d_{23} = d_{31}$ $d_{35} = d_{16} = d_{24}, d_{26} = d_{34} = d_{15}$ $(g_4, g_5, g_6)$ $(d_{13} - d_{12}, d_{21} - d_{23}, d_{32} - d_{31})$ $(d_{35} - d_{26}, d_{16} - d_{34}, d_{24} - d_{15})$	6	3	0
	$(x_1, x_1, 0)$	$m_{xy}$	$C_{sxy}$	6				
	$(x_1, x_1, x_1)$	$3_p$	$C_{3p}$	4				
$T_2$ (La)	$(0, 0, z_2)$	$m_{xy} m_{xy} 2_z$	$\hat{C}_{2vz}$	3	$P_3; u_6$ $P_1 = -P_2, P_3; u_4 = -u_5, u_6$ $P_1 = P_2 = P_3; u_4 = u_5 = u_6$ $(P_1, P_2, P_3); (u_4, u_5, u_6)$	6	6	6
	$(x_2, -x_2, z_2)$	$m_{xy}$	$C_{sxy}$	6				
	$(x_2, x_2, x_2)$	$3_x m_{xy}$	$C_{3xp}$	4				
	$(x_2, y_2, z_2)$	1	$C_1$	1				
<b>Parent symmetry <math>G</math>: <math>m\bar{3}m</math> <math>O_h</math></b>								
$A_{2g}$	$x_2^+$	$m\bar{3}$	$T_h$	1	$A_{14} = A_{25} = A_{36}; \pi_{23}^a = \pi_{31}^a = \pi_{12}^a$	2	1	0
$A_{1u}$	$x_1^-$	432	$O$	1	$\varepsilon; g_1 = g_2 = g_3;$	2	1	0
$A_{2u}$	$x_2^-$	$\bar{4}3m$	$T_d$	1	$d_{14} = d_{25} = d_{36}$	2	1	0
$E_g$ (La)	$(x_3^+, 0)$	$4_z/m_z m_x m_{xy}$	$D_{4hz}$	3	$\delta u_3$ $[\delta u_3 - a(\delta u_1 + \delta u_2), b(\delta u_1 - \delta u_2)]$	3	3	0
	$(x_3^+, y_3^+)$	$m_x m_y m_z$	$D_{2h}$	1				
$E_u$	$(x_3^-, 0)$	$4_z 2_x 2_{xy}$	$D_{4z}$	3	$g_1 = g_2, g_3; d_{14} = -d_{25}$ $g_1 = -g_2; d_{14} = d_{25} = d_{36}$ $[g_3 - a(g_1 + g_2), b(g_1 - g_2)]$ $[b(d_{14} - d_{25}), a(d_{14} + d_{25}) - d_{36}]$	12	3	0
	$(0, y_3^-)$	$\bar{4}_z 2_x m_{xy}$	$D_{2dz}$	3				
	$(x_3^-, y_3^-)$	$2_x 2_y 2_z$	$D_2$	1				
$T_{1g}$	$(0, 0, z_1^+)$	$4_z/m_z$	$\hat{C}_{4hz}$	3	$A_{33}, A_{32} = A_{31}, A_{24} = A_{15}, A_{14} = -A_{25}$ $A_{11} = A_{22},$ $A_{13} = A_{23}, A_{12} = A_{21}$ $A_{35} = A_{34}, A_{26} = A_{16}$ $A_{11} = A_{22} = A_{33}$ $A_{13} = A_{21} = A_{32}, A_{12} = A_{32} = A_{31}$ $A_{35} = A_{16} = A_{24}, A_{26} = A_{34} = A_{15}$ $(A_{11}, A_{22}, A_{33})$ $(A_{13} + A_{12}, A_{21} + A_{23}, A_{32} + A_{31})$ $(A_{35} + A_{26}, A_{16} + A_{34}, A_{24} + A_{15})$	6	3	0
	$(x_1^+, x_1^+, 0)$	$2_{xy}/m_{xy}$	$C_{2hxy}$	6				
	$(x_1^+, x_1^+, x_1^+)$	$\bar{3}_p$	$C_{3ip}$	4				
$T_{2g}$ (La)	$(0, 0, z_2^+)$	$m_{xy} m_{xy} m_z$	$\hat{D}_{2hz}$	3	$u_6$ $u_4 = -u_5, u_6$ $u_4 = u_5 = u_6$ $(u_4, u_5, u_6)$	6	6	0
	$(x_2^+, -x_2^+, z_2^+)$	$2_{xy}/m_{xy}$	$C_{2hxy}$	6				
	$(x_2^+, x_2^+, x_2^+)$	$\bar{3}_p m_{xy}$	$D_{3dp}$	4				
	$(x_2^+, y_2^+, z_2^+)$	1	$C_i$	1				
$T_{1u}$	$(0, 0, z_1^-)$	$4_z m_x m_{xy}$	$C_{4vz}$	3	$P_3$ $P_1, P_2$ $P_1 = P_2$ $P_1 = -P_2, P_3$ $P_1 = P_2 = P_3$ $(P_1, P_2, P_3)$	6	3	6
	$(x_1^-, y_1^-, 0)$	$m_z$	$C_{3z}$	3				
	$(x_1^-, x_1^-, 0)$	$m_{xy} 2_{xy} m_z$	$\hat{C}_{2vxy}$	6				
	$(x_1^-, -x_1^-, z_1^-)$	$m_{xy}$	$C_{sxy}$	6				
	$(x_1^-, x_1^-, x_1^-)$	$3_p m_{xy}$	$C_{3vp}$	4				
	$(x_1^-, y_1^-, z_1^-)$	1	$C_1$	1				
$T_{2u}$	$(0, 0, z_2^-)$	$\bar{4}_z m_x 2_{xy}$	$\hat{D}_{2dz}$	3	$g_6; d_{32} = -d_{31}, d_{24} = -d_{15}$ $g_4, g_5; d_{13}, d_{12}, d_{21}, d_{23}$ $d_{35}, d_{26}, d_{16}, d_{34}$ $g_4 = -g_5; d_{13} = d_{23}, d_{21} = d_{21}$ $d_{35} = d_{34}, d_{16} = d_{26}$ $g_4 = -g_5, g_6; d_{13} = d_{23}, d_{21} = d_{21}$ $d_{35} = d_{34}, d_{16} = d_{26}$ $d_{32} = -d_{31}, d_{24} = -d_{15}$ $g_4 = g_5 = g_6;$ $d_{13} = -d_{12} = d_{21} = -d_{23} = d_{32} - d_{31}$ $d_{35} = -d_{26} = d_{16} = -d_{34} = d_{24} = -d_{15}$ $(g_4, g_5, g_6)$ $(d_{13} - d_{12}, d_{21} - d_{23}, d_{32} - d_{31})$ $(d_{35} - d_{26}, d_{16} - d_{34}, d_{24} - d_{15})$	6	3	0
	$(x_2^-, y_2^-, 0)$	$m_z$	$C_{sz}$	3				
	$(x_2^-, -x_2^-, 0)$	$m_{xy} 2_{xy} m_z$	$\hat{C}_{2vxy}$	6				
	$(x_2^-, -x_2^-, z_2^-)$	$2_{xy}$	$C_{2xy}$	6				
	$(x_2^-, x_2^-, x_2^-)$	$3_p 2_{xy}$	$D_{3p}$	4				
	$(x_2^-, y_2^-, z_2^-)$	1	$C_1$	1				

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The solution of the inverse Landau problem – *i.e.* the identification of the representation  $\Gamma_\eta$  relevant to symmetry descent  $G \Downarrow F$  – enables one to determine the corresponding normal mode (so-called soft mode) of the transition (see *e.g.* Rousseau *et al.*, 1981). We note that this step requires additional knowledge of the crystal structure, whereas other conclusions of the analysis hold for *any* crystal structure with a given symmetry descent  $G \Downarrow F$ . Normal-mode determination reveals the dynamic microscopic nature of the instability of the crystal lattice which leads to the phase transition (for more details and examples, see Section 3.1.5).

The representation  $\Gamma_\eta$  further determines the principal tensor parameters associated with the primary order parameter  $\eta$ . If one of them is a vector (polarization) the soft mode is infrared-active in the parent phase; if it is a symmetric second-rank tensor (spontaneous strain), the soft mode is Raman active in this phase. Furthermore, the *R*-irep  $\Gamma_\eta$  determines the polynomial in components of  $\eta$  in the Landau free energy (basic invariant polynomials, called *integrity bases*, are available in the software *GI★KoBo-1* and in Kopský, 2001) and allows one to decide whether the necessary conditions of continuity of the transition (so-called Landau and Lifshitz conditions) are fulfilled.

(2) *Direct Landau problem of equitranslational phase transitions:* For a given space group  $\mathcal{G}$  of the parent phase and the *R*-irep  $\Gamma_\eta$  (specifying the transformation properties of the primary order parameter  $\eta$ ), find the corresponding equitranslational space group  $\mathcal{F}$  of the ferroic phase. To solve this task, one first finds in Table 3.1.3.1 the point group  $F$  that corresponds to point group  $G$  of space group  $\mathcal{G}$  and to the given *R*-irep  $\Gamma_\eta$ . The point-group symmetry descent  $G \Downarrow F$  thus obtained specifies uniquely the equitranslational subgroup  $\mathcal{F}$  of  $\mathcal{G}$  that can be found in the lattices of equitranslational subgroups of space groups available in the software *GI★KoBo-1* (see Section 3.1.6).

(3) *Secondary tensor parameters of an equitranslational phase transition  $\mathcal{G} \Downarrow \mathcal{F}$ :* These parameters are specified by the representation  $\Gamma_\lambda$  of  $G$  associated with a symmetry descent  $\Gamma \Downarrow L$ , where  $L$  is an intermediate group [see equation (3.1.3.1)]. In other words, the secondary tensor parameters of the transition  $G \Downarrow F$  are identical with principal tensor parameters of the transition  $G \Downarrow L$ . To each intermediate group  $L$  there corresponds a set of secondary tensor parameters. All intermediate subgroups of a symmetry descent  $G \Downarrow F$  can be deduced from lattices of subgroups in Figs. 3.1.3.1 and 3.1.3.2.

The representation  $\Gamma_\lambda$  specifies transformation properties of the secondary tensor parameter  $\lambda$  and thus determines *e.g.* its

infrared and Raman activity in the parent phase and enables one to make a mode analysis. Representation  $\Gamma_\lambda$  together with  $\Gamma_\eta$  determine the coupling between secondary and primary tensor parameters. The explicit form of these faint interactions (Aizu, 1973; Kopský, 1979*d*) can be found in the software *GI★KoBo-1* and in Kopský (2001).

(4) *Changes of property tensors at a ferroic phase transition.* These changes are described by tensor parameters that depend only on the point-group-symmetry descent  $G \Downarrow F$ . This means that *the same principal tensor parameters and secondary tensor parameters appear in all equitranslational and in all non-equitranslational transitions with the same  $G \Downarrow F$ .* The only difference is that in non-equitranslational ferroic phase transitions a principal tensor parameter corresponds to a secondary ferroic order parameter. It still plays a leading role in tensor distinction of domains, since it exhibits different values in any two ferroic domain states (see Section 3.4.2.3). Changes of property tensors at ferroic phase transitions are treated in detail in the software *GI★KoBo-1* and in Kopský (2001).

We note that Table 3.1.3.1 covers only those point-group symmetry descents  $G \Downarrow F$  that are ‘driven’ by *R*-ireps of  $G$ . All possible point-group symmetry descents  $G \Downarrow F$  are listed in Table 3.4.2.7. Principal and secondary tensor parameters of symmetry descents associated with reducible representations are combinations of tensor parameters appearing in Table 3.1.3.1 (for a detailed explanation, see the manual of the software *GI★KoBo-1* and Kopský, 2000). Necessary data for treating these cases are available in the software *GI★KoBo-1* and Kopský (2001).

#### 3.1.3.3.1. Explanation of Table 3.1.3.1

*Parent symmetry  $G$ :* the short international (Hermann–Mauguin) and the Schoenflies symbol of the point group  $G$  of the parent phase are given. Subscripts specify the orientation of symmetry elements (generators) in the Cartesian crystallophysical coordinate system of the group  $G$  (see Figs. 3.4.2.3 and 3.4.2.4, and Tables 3.4.2.5 and 3.4.2.6).

*R-irep  $\Gamma_\eta$ :* physically irreducible representation  $\Gamma_\eta$  of the group  $G$  in the spectroscopic notation. This representation defines transformation properties of the primary order parameter  $\eta$  and of the principal tensor parameters. Each complex irreducible representation is combined with its complex conjugate and thus a real physically irreducible representation *R*-irep is formed. Matrices  $D^{(\alpha)}$  of *R*-ireps are given explicitly in the the software *GI★KoBo-1*.

Table 3.1.3.2. Symmetry descents  $G \Downarrow F_1$  associated with two irreducible representations

$G$	$\Gamma_\eta$	$F_1$	Proper or improper		Domain states			Full or partial	
			Ferroelectric	Ferroelastic	$n_f$	$n_e$	$n_a$	Ferroelectric	Ferroelastic
432	$T_1$	$2_{xy}$	proper	improper	12	12	12	full	full
	$T_2$		improper	proper					
	$T_1$	1	improper	improper	24	24	24	full	full
	$T_2$		proper	proper					
$\bar{4}3m$	$T_1$	$m_{xy}$	improper	improper	12	12	12	full	full
	$T_2$		proper	proper					
	$T_1$	1	improper	improper	24	24	24	full	full
	$T_2$		proper	proper					
$m\bar{3}m$	$T_{1g}$	$2_{xy}/m_{xy}$	non	improper	12	0	12	non	full
	$T_{2g}$		non	proper					
	$T_{1g}$	$\bar{1}$	non	improper	24	0	24	non	full
	$T_{2g}$		non	proper					
	$T_{1u}$	$m_{xy}2_{xy}m_z$	proper	improper	12	12	6	full	partial
	$T_{2u}$		improper	improper					
	$T_{1u}$	$m_z$	proper	improper	24	24	12	full	partial
	$T_{2u}$		improper	improper					
	$T_{1u}$	1	proper	improper	48	48	24	full	partial
	$T_{2u}$		improper	improper					



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provided by the *conversion equations* that express Cartesian tensor components as linear combinations of principal and secondary covariant components (for more details on tensorial covariants and conversion equations see Appendix E of the manual for *GI★KoBo-1* and Kopský, 2001).

We illustrate the situation on a transition with symmetry descent  $4_z 2_x 2_{xy} \Downarrow 2_x 2_y 2_z$ . In Table 3.1.3.1, we find that the principal tensor parameter transforms according to irreducible representation  $B_1$  with standard variable  $x_3$ . The corresponding covariant  $u_3 = u_1 - u_2$  can be found in Appendix E of the manual of *GI★KoBo-1* (or in Kopský, 2001), where one also finds an invariant containing  $u_1$  and  $u_2$ :  $u_{1,1} = u_1 + u_2$ . The corresponding conversion equations are:  $u_1 = \frac{1}{2}(u_{1,1} + u_3)$ ,  $u_2 = \frac{1}{2}(u_{1,1} - u_3)$ . In the parent phase  $u_3 = u_1^{(p)} - u_2^{(p)} = 0$ , hence  $u_1^{(p)} = u_2^{(p)} = \frac{1}{2}u_{1,1}$ , whereas in the ferroic phase  $u_1^{(f)} = \frac{1}{2}(u_{1,1} + u_3) = u_1^{(p)} + \frac{1}{2}u_{1,1} = u_1^{(p)} + \delta u_1$ ,  $u_2^{(f)} = u_2^{(p)} - \frac{1}{2}u_{1,1} = u_2^{(p)} + \delta u_2 = u_1^{(p)} - \delta u_1$ . The symmetry-breaking increments  $\delta u_1 = -\delta u_2$  describe thus the changes of the Cartesian components that correspond to the nonzero principal tensor component  $u_1 - u_2$ .

An analogous situation occurs frequently in trigonal and hexagonal parent groups, where  $u_1 - u_2$  (or  $g_1 - g_2$ ) transforms like the first or second component of the principal tensor parameter. In these cases, the corresponding symmetry-breaking increments of Cartesian components are again related:  $\delta u_1 = -\delta u_2$  (or  $\delta g_1 = -\delta g_2$ ).

We note that relations like  $A_{11} = -A_{12} = -A_{26}$  do not imply that these components transform as the standard variable. Though these components are proportional to the principal tensor parameter in the first domain state, they cannot be transformed to corresponding components in other domain states as easily as covariant tensor components of the principal tensor parameter.

In general, it is useful to consider a tensor parameter as a vector in the carrier space of the respective representation. Then the Cartesian components are projections of this vector on the Cartesian basis of the tensor space.

The presentation of the principal tensor parameters in the column *Principal tensor parameters* of this table is a compromise: whenever conversion equations lead to simple relations between morphic Cartesian components and/or symmetry-breaking increments, we present these relations, in some cases together with corresponding covariants. In the more complicated cases, only the covariants are given. The corresponding conversion equations and labelling of covariants are given at the beginning of that part of the table which covers hexagonal and cubic parent groups  $G$ . In the main tables of the software *GI★KoBo-1*, the principal tensor parameters and the secondary tensor parameters up to rank 4 are given consistently in covariant form. Labelling of covariant components and conversion equations are given in Appendix E of the manual.

The principal tensor parameters presented in Table 3.1.3.1 represent a particular choice of property tensors for standard variables given in the second column. To save space, property tensors are selected in the following way: polarization  $\mathbf{P}$  and strain  $u$  are always listed; if none of their components transform according to  $D^{(n)}$ , then components of one axial and one polar tensor (if available) appearing in Table 3.1.3.3 are given. Principal parameters of two different property tensors are separated by a semicolon. If two different components of the same property tensor transform in the same way, they are separated by a comma.

As tensor indices we use integers 1, 2, 3 instead of vector components  $x, y, z$  and contracted indices 1, 2, 3, 4, 5, 6 in matrix notation for pairs  $xx, yy, zz, yz \approx zy, zx \approx xz, xy \approx yx$ , respectively

*Important note:* To make Table 3.1.3.1 compatible with the software *GI★KoBo-1* and with Kopský (2001), coefficients of property tensors in matrix notation with contracted indices 4, 5, 6 do not contain the numerical factors 2 and 4 which are usually

Table 3.1.3.3. *Important property tensors*

$i = 1, 2, 3; \mu, \nu = 1, 2, \dots, 6$ .

Tensor components	Property	Tensor components	Property
$\varepsilon$	enantiomorphism		chirality
$P_i$	polarization	$P_i$	pyroelectricity
$u_\mu$	strain	$\varepsilon_{ij}$	dielectric permittivity
$g_\mu$	optical activity		
$d_{i\mu}$	piezoelectricity	$r_{i\mu}$	electro-optics
$A_{i\mu}$	electrogyration		
$\pi_{\mu\nu}$	piezo-optics	$Q_{\mu\nu}$	electrostriction

introduced to preserve a compact form (without these factors) of linear constitutive relations [see Chapter 1.1, Nye (1985) and especially Appendices E and F of Sirotin & Shaskolskaya (1982)]. This explains the differences in matrix coefficients appearing in Table 3.1.3.1 and those presented in Chapter 1.1 or in Nye (1985) and in Sirotin & Shaskolskaya (1982). Thus *e.g.* for the symmetry descent  $6_z 2_x 2_y \Downarrow 3_z 2_x$ , we find in Table 3.1.3.1 the principal tensor parameters  $d_{11} = -d_{12} = -d_{26}$ , whereas according to Chapter 1.1 or *e.g.* to Nye (1985) or Sirotin & Shaskolskaya (1982) these coefficients for  $F_1 = 3_z 2_x$  are related by equations  $d_{11} = -d_{12} = -2d_{26}$ .

Property tensors and symbols of their components that can be found in Table 3.1.3.1 are given in the left-hand half of Table 3.1.3.3. The right-hand half presents other tensors that transform in the same way as those on the left and form, therefore, covariant tensor components of the same form as those given in the column *Principal tensor parameters*. Principal and secondary tensor parameters for all property tensors that appear in Table 3.1.3.3 are available in the software *GI★KoBo-1*.

$n_f$ : number of ferroic single domain states that differ in the primary order parameter  $\eta$  and in the principal tensor parameters.

$n_a$ : number of ferroelastic single domain states. If  $n_a = n_f$ ,  $n_a < n_f$  or  $n_a = 1$ , the ferroic phase is, respectively, a full, partial or non-ferroelastic one.

$n_e$ : number of ferroelectric single domain states. If  $n_e = n_f$ ,  $n_e < n_f$  or  $n_e = 0, 1$ , the ferroic phase is, respectively, a full, partial or non-ferroelectric one ( $n = 0$  or  $n = 1$  correspond to a non-polar or to a polar parent phase, respectively) (see Section 3.4.2).

#### 3.1.3.4. Examples

*Example 3.1.3.4.1. Phase transition in triglycine sulfate (TGS).* Assume that the space groups of both parent (high-symmetry) and ferroic (low-symmetry) phases are known:  $\mathcal{G} = P2_1/c (C_{2h}^5)$ ,  $\mathcal{F}_1 = P2_1 (C_2^2)$ . The same number of formula units in the primitive unit cell in both phases suggests that the transition is an equitranslational one. This conclusion can be checked in the lattice of equitranslational subgroups of the software *GI★KoBo-1*. There we find for the low-symmetry space group the symbol  $P112_1(\mathbf{b}/4)$ , where the vector in parentheses expresses the shift of the origin with respect to the conventional origin given in *IT A* (2002).

In Table 3.1.3.1, one finds that the corresponding point-group-symmetry descent  $2_z/m_z \Downarrow 2_z$  is associated with irreducible representation  $\Gamma_\eta = A_u$ . The corresponding principal tensor parameters of lowest rank are the pseudoscalar  $\varepsilon$  (enantiomorphism or chirality) and the vector of spontaneous polarization with one nonzero morphic component  $P_3$  – the transition is a proper ferroelectric one. The non-ferroelastic ( $n_a = 1$ ) full ferroelectric phase has two ferroelectric domain states ( $n_f = n_e = 2$ ). Other principal tensor parameters (morphic tensor components that transform according to  $\Gamma_\eta$ ) are available in the software *GI★KoBo-1*:  $g_1, g_2, g_3, g_6; d_{31}, d_{32}, d_{33}, d_{36}, d_{14}, d_{15}, d_{24}, d_{25}$ . Property tensors with these components are listed in Table 3.1.3.3. As shown in Section 3.4.2, all these components