

the elastic moduli at zero field and pressure are shown in Fig. 2. Data published by Rosen and Klimker [9] show similar, but significantly different, effects. The differences between Fig. 2 and Rosen's data are primarily in the variation of c_{12} ; consequently, the variations of the compressibilities β_{\parallel} and β_{\perp} with turn angle and structure, that are calculated from the Fig. 2 data, are considerably different than given by Rosen. In the present context, however, there appears no indication that magnetic ordering has any influence on the second order elastic moduli at 298°K.

The magnetic structures in Er are even more complex. The spontaneous ferromagnetic phase, with $T_c = 20^\circ\text{K}$, contains a spiral component. The ferromagnetic component transforms to an antiferromagnetic arrangement, between 20 and 53°K, and a modulated moment arrangement between 53 and 80°K.

The effects of the ordering on the elastic moduli are shown in Fig. 3. These effects are clearly not evident at 298°K.

The values of the second order elastic stiffness moduli at 298°K are listed in Table 2. We include here the data for Ho as given by Palmer and Lee [8]. Included in Table 2 are the values of the density [18] that are used for computing the moduli. The calculated adiabatic and isothermal bulk moduli, K_s and K_T , and also given and the last column gives the parameters $(\beta_{\parallel} - \beta_{\perp})/\beta_V$ calculated from isothermal values of the compressibilities. The latter parameter is that used for computing equation (3). The variations of the data with increasing atomic number (Gd \rightarrow Er) are noteworthy. The c/a ratio decreases whereas the density increases because of the so-called lanthanide contraction associated with the addition of electrons to the 4f shell. There is

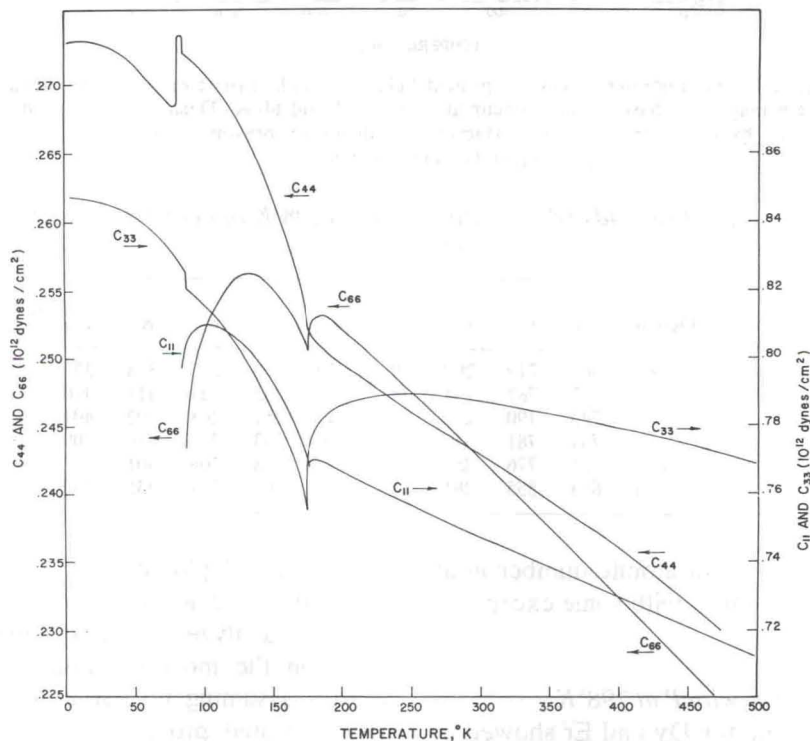


Fig. 2. Temperature dependence of the principal elastic moduli of Dy at zero magnetic field. $T_N = 179^\circ\text{K}$ and $T_c = 87^\circ\text{K}$. Absence of data for c_{11} and c_{66} at $T < 85^\circ\text{K}$ is result of spontaneous macroscopic distortion by anisotropic magnetostriction.

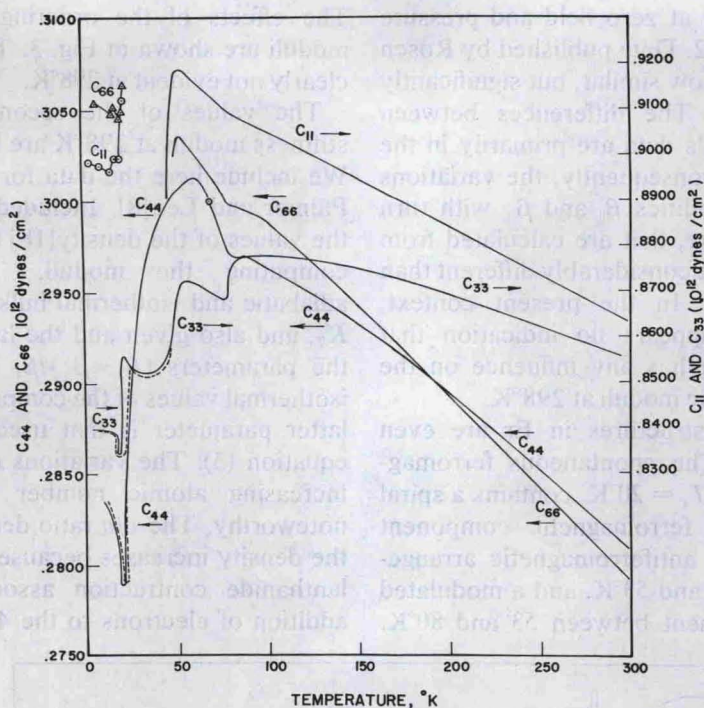


Fig. 3. Temperature dependence of principal elastic moduli in Er at zero magnetic field, where magnetic phase changes occur at ~ 20 , ~ 53 and 80°K . Dashed lines indicate thermal hysteresis in $T_c \sim 20^\circ\text{K}$. Macroscopic distortion prevented measurements of c_{11} and c_{66} in spiral phase.

Table 2. Values of second order elastic moduli at 298°K of heavy rare earth metals (Kbars)

| | c/a | Density | c_{11} | c_{33} | c_{44} | c_{66} | C_H | c_{12} | c_{13} | K_s | K_T | $\left(\frac{\beta_{\parallel} - \beta_{\perp}}{\beta_V}\right)_T$ |
|-------|-------|---------|----------|----------|----------|----------|-------|----------|----------|-------|-------|--|
| Gd | 1.590 | 7.888 | 667 | 719 | 207 | 208 | 250 | 250 | 213 | 378 | 373 | +0.007 |
| Dy(F) | 1.573 | 8.560 | 747 | 787 | 243 | 243 | 278 | 262 | 223 | 411 | 410 | -0.005 |
| Dy(R) | | 8.545 | 743 | 790 | 255 | 246 | 290 | 251 | 208 | 402 | 401 | -0.009 |
| Dy(P) | | 8.560 | 731 | 781 | 240 | 239 | 276 | 253 | 223 | 410 | 409 | -0.015 |
| Ho | 1.570 | 8.800 | 761 | 776 | 257 | 256 | 290 | 248 | 206 | 401 | | +0.019 |
| Er | 1.569 | 9.064 | 863 | 855 | 281 | 279 | 328 | 305 | 227 | 455 | 450 | +0.043 |

a general increase with atomic number in all of the c_{ij} and K values, with some exceptions for the Ho data.

(b) Variations of c_{ij} with P at 298°K

Plots of the data for Dy and Er showed no significant or systematic departure from linear relations between the c_{ij} and pressure. The Gd data showed small but significant curvature

for the c_{33} vs P plot at $P > 2$ Kbar and for c_{44} at $P = 0$ and 2 Kbar. All of the data were, however, analyzed by least square statistics to obtain the most probable values for the dc_{ij}/dP , assuming no curvature. The results with indicated probable errors are given in Table 3. The similarities and differences in the effects of pressure on the c_{ij} of the three metals are evident in the pressure derivatives