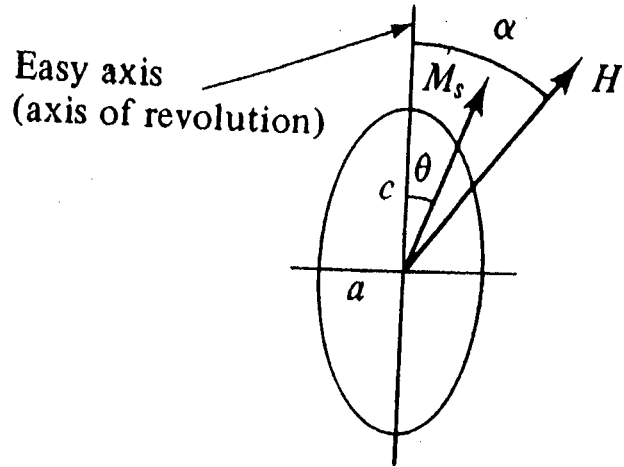


HYSTERESIS LOOP

$$10^{-3} < H_c < 2 \times 10^4$$

(a)



(b)

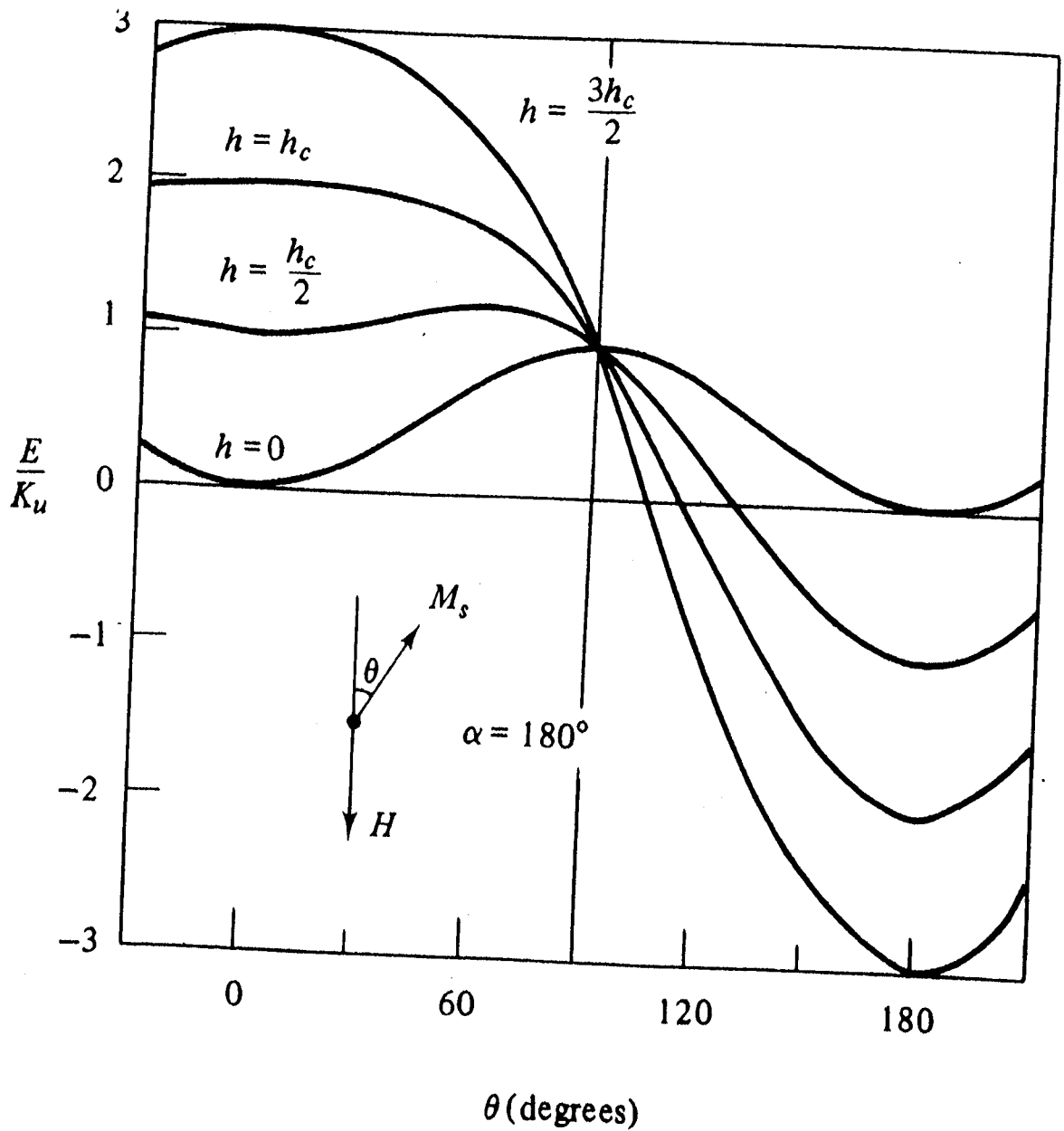


Fig. 9.37 Rotation of magnetization in an ellipsoid. $h =$ reduced field.

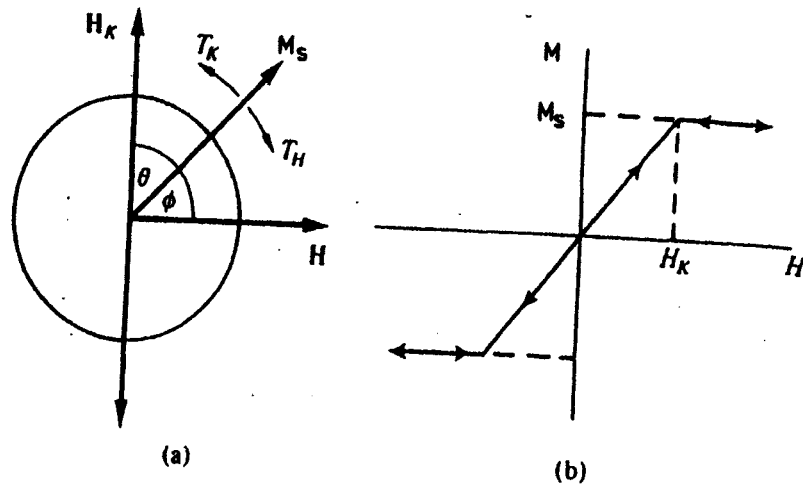


Fig. 13.5. (a) A spherical single-domain particle with anisotropy field H_K and with the magnetic field H perpendicular to the easy axis; (b) magnetization curve obtained for the situation depicted in (a).

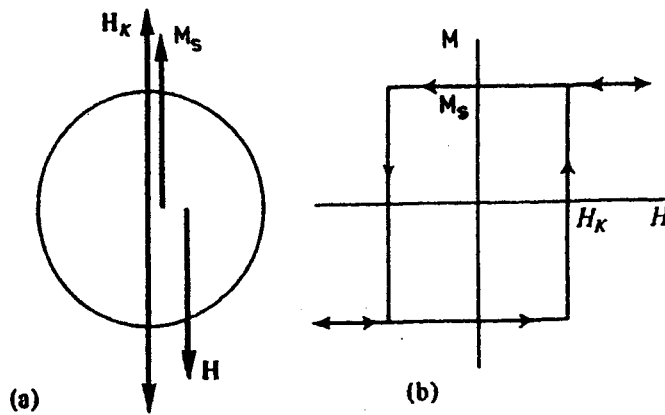


Fig. 13.6 (a) A spherical single-domain particle with anisotropy field H_K and with the magnetic field H parallel to the easy axis; (b) magnetization curve obtained for the situation depicted in (a).

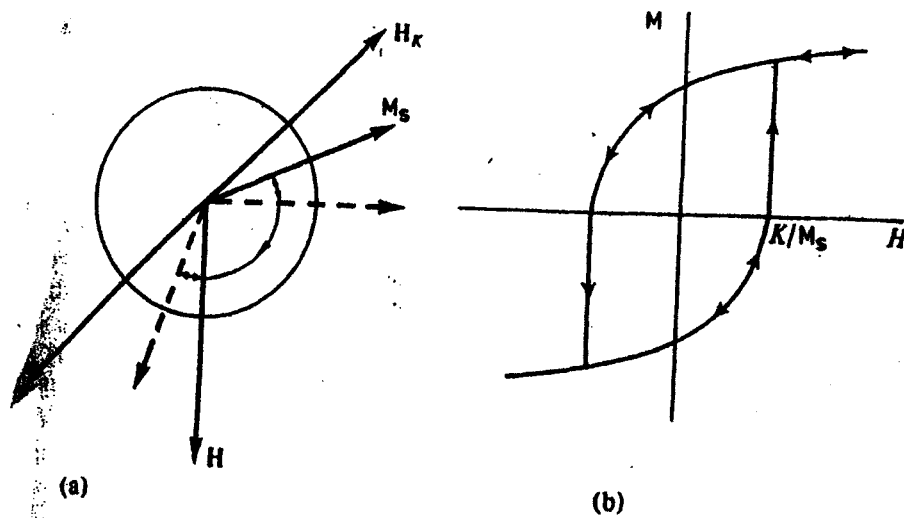
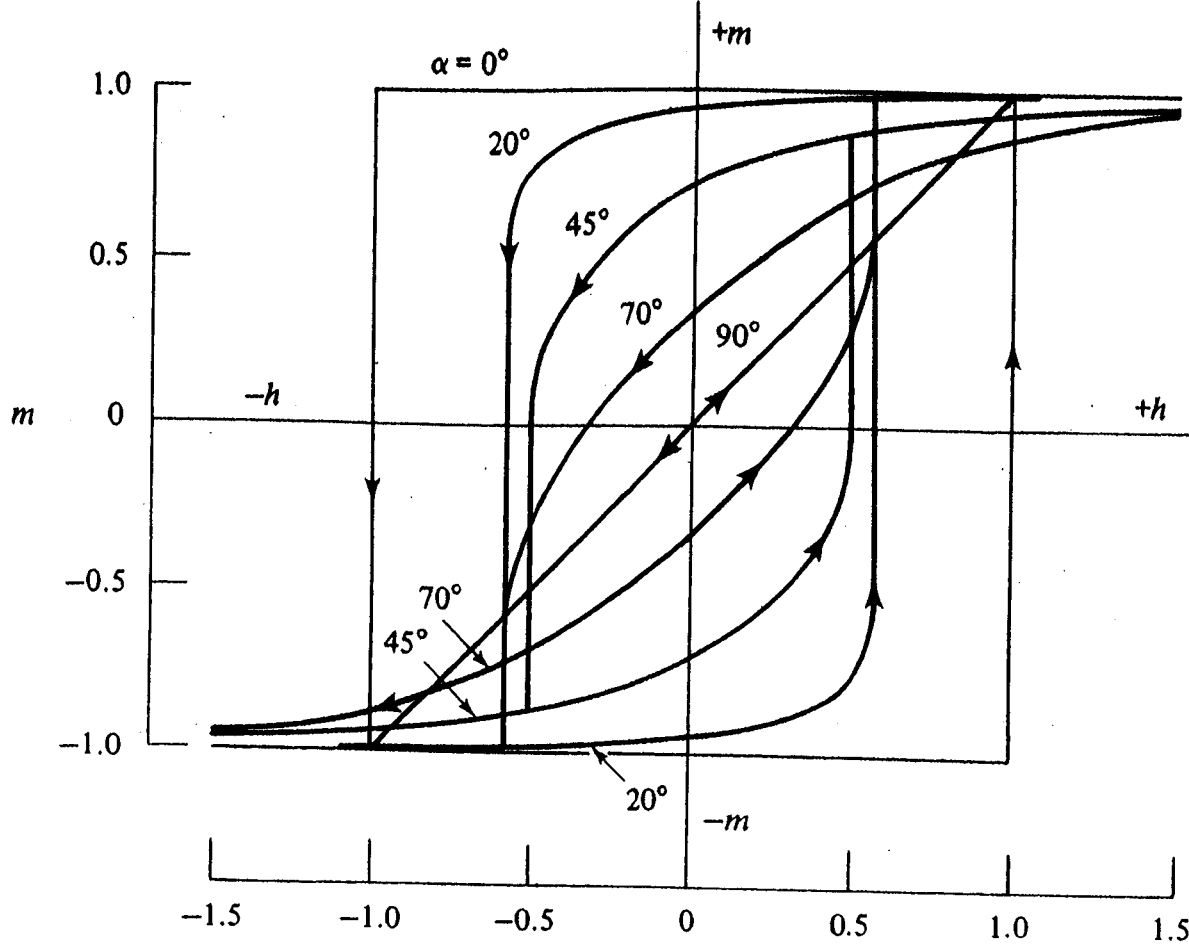
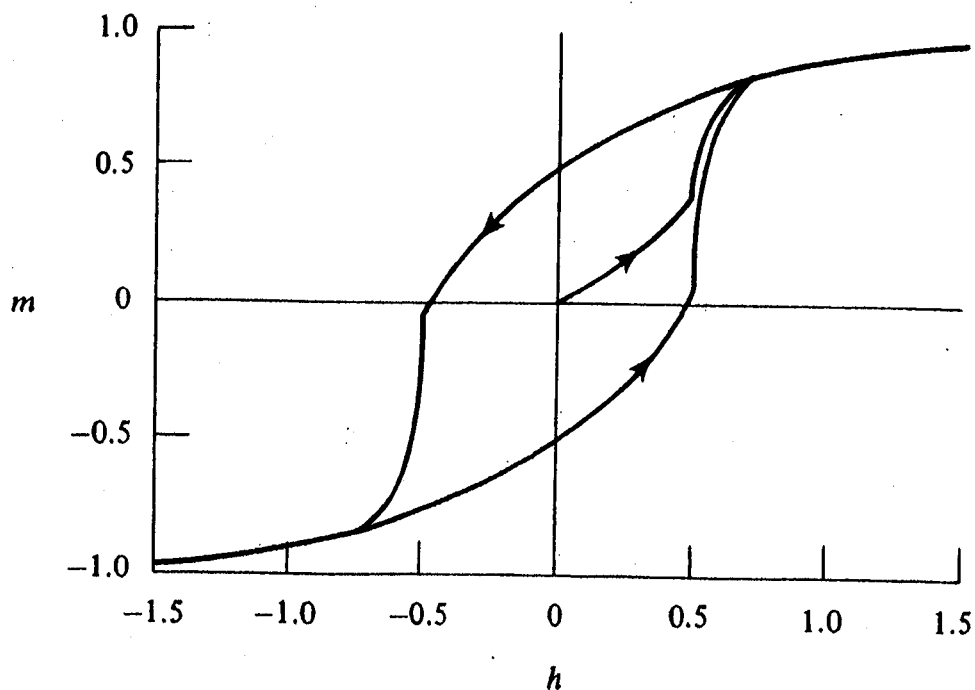


Fig. 13.7 (a) A spherical single-domain particle with anisotropy field H_K and with the magnetic field H at an arbitrary angle to the easy axis; (b) magnetization curve obtained for the situation depicted in (a).

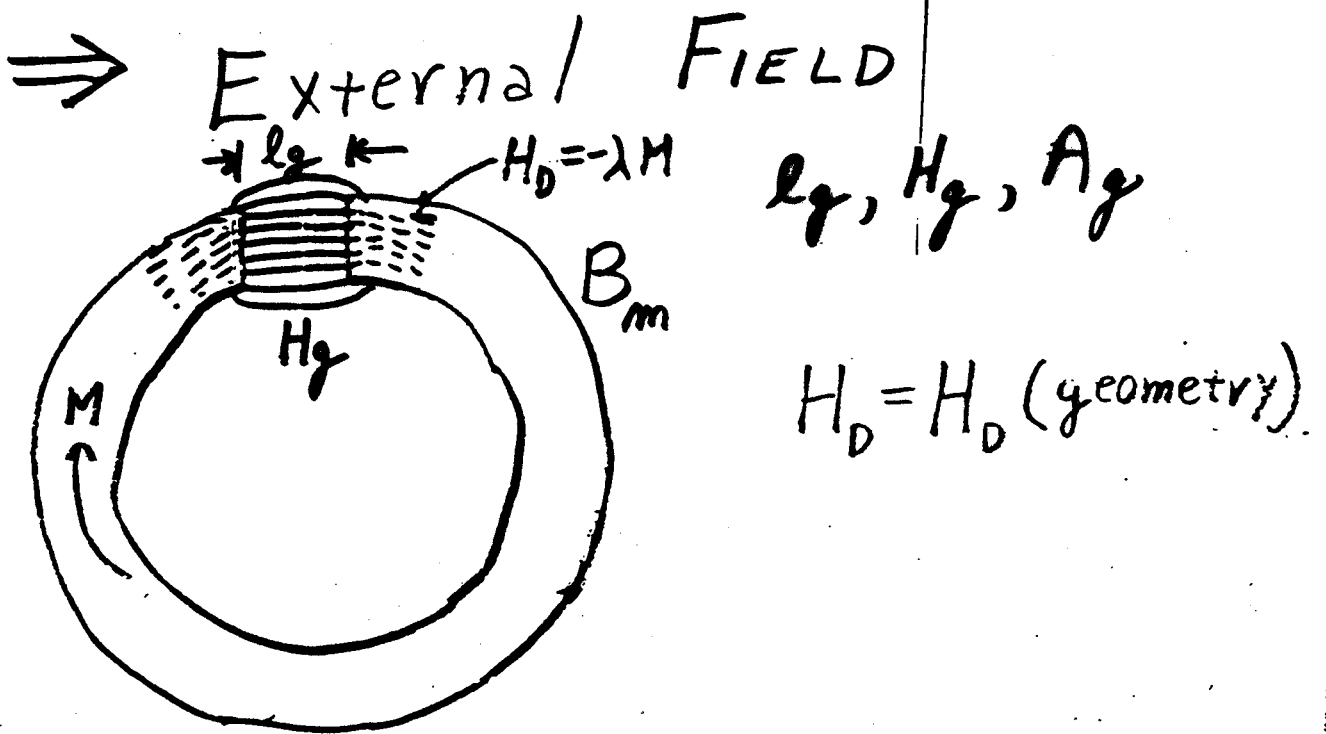


Hysteresis loops for uniaxial anisotropy. ($\alpha =$ angle between field and easy axis.)



Hysteresis loop of an assembly of single-domain, uniaxial particles having their easy axes randomly oriented. Stoner and Wohlfarth [7.23], Rhodes [9.31].

PERMANENT MAGNET

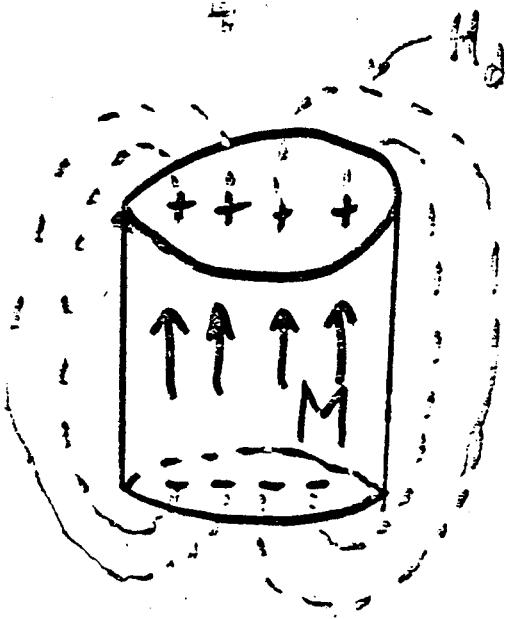


OPEN CIRCUIT ⇒

FREE Poles

DEMAGNETIZATION FIELDS

$$\underline{H_d = -N_x M_x - N_y M_y - N_z M_z = -N \cdot M}$$



$$B = H + 4\pi M$$

$$\underline{\nabla \cdot B = 0}$$

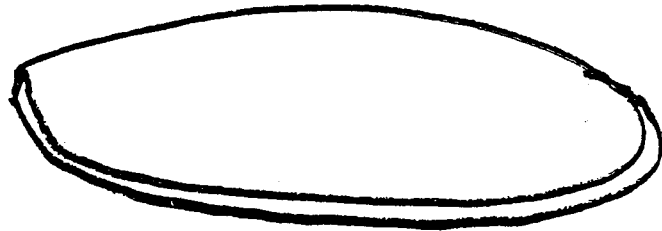
$$\underline{N_x + N_y + N_z = 4\pi}$$

N = DEMAGNETIZING FACTOR
= f(geometry only)

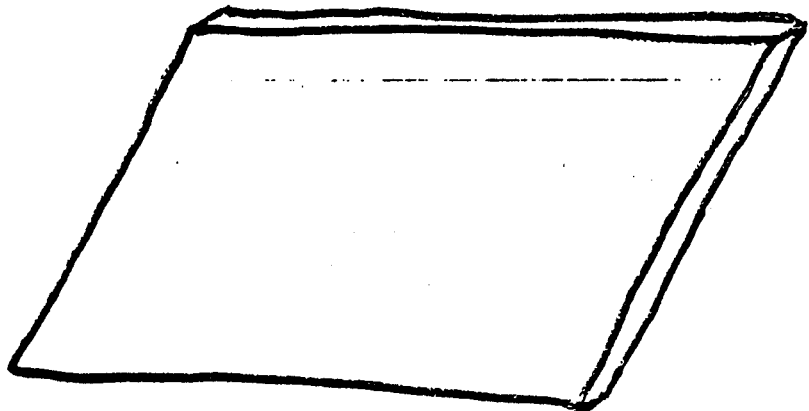


$$N_z = 0$$

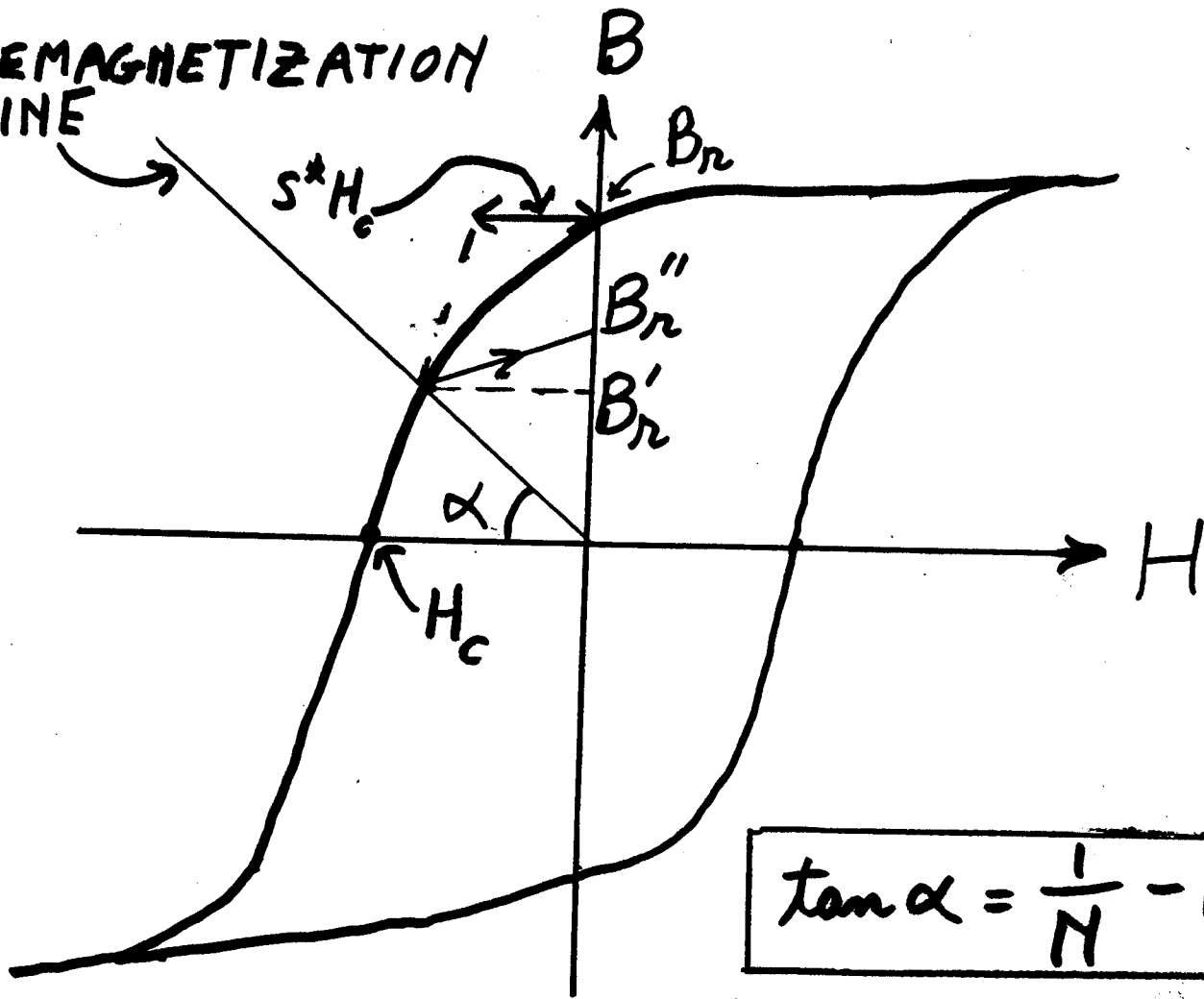
$$N_x = N_y = 2\pi$$



$$N_z = 4\pi, N_x = N_y = 0$$



DEMAGNETIZATION
LINE

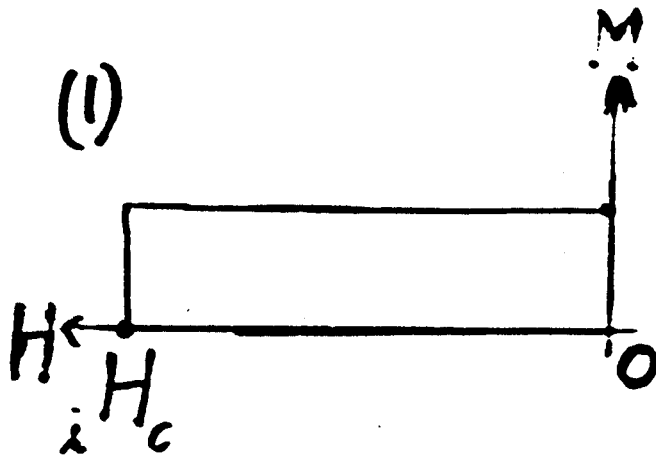


$$\tan \alpha = \frac{1}{N} - 1$$

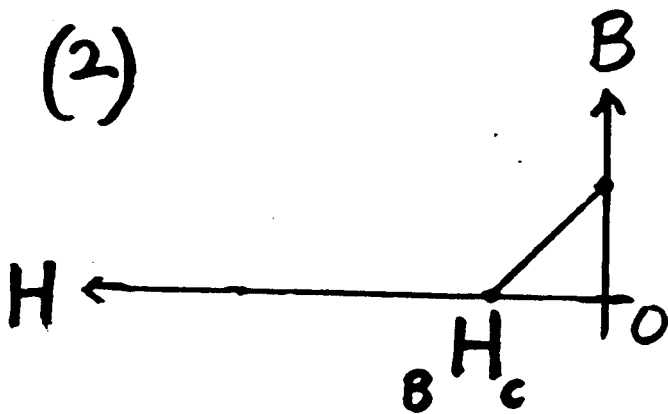
REDUCED REMNANCE, B_r' , due to
DEMAGNETIZATION N .

B_r'' = PARTIAL RECOVERY when
magnetic circuit closed by playback
head core.

PERMANENT MAGNETS



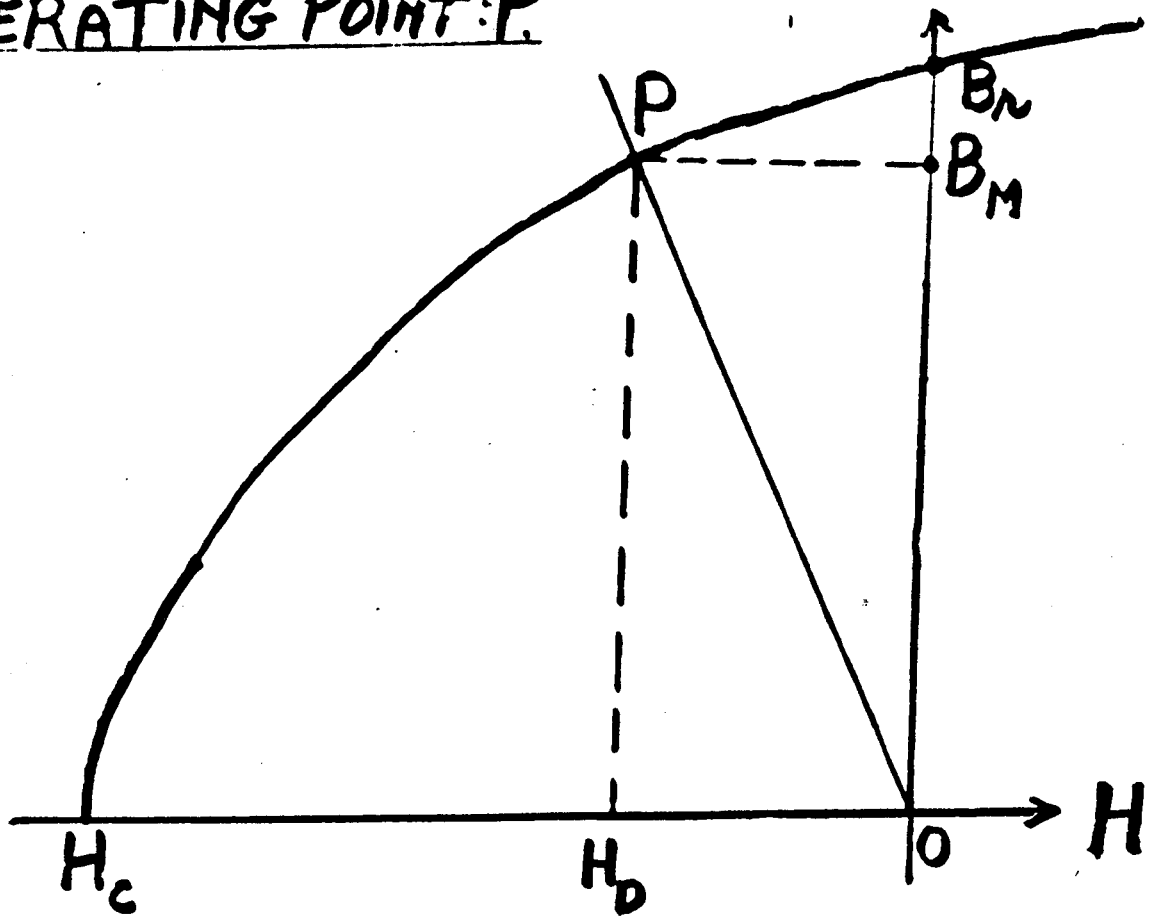
$$B = H + 4\pi M$$
$$\Rightarrow B H_c = -4\pi M$$



ENERGY Storage $\propto (BH)_{max}$ of (2).

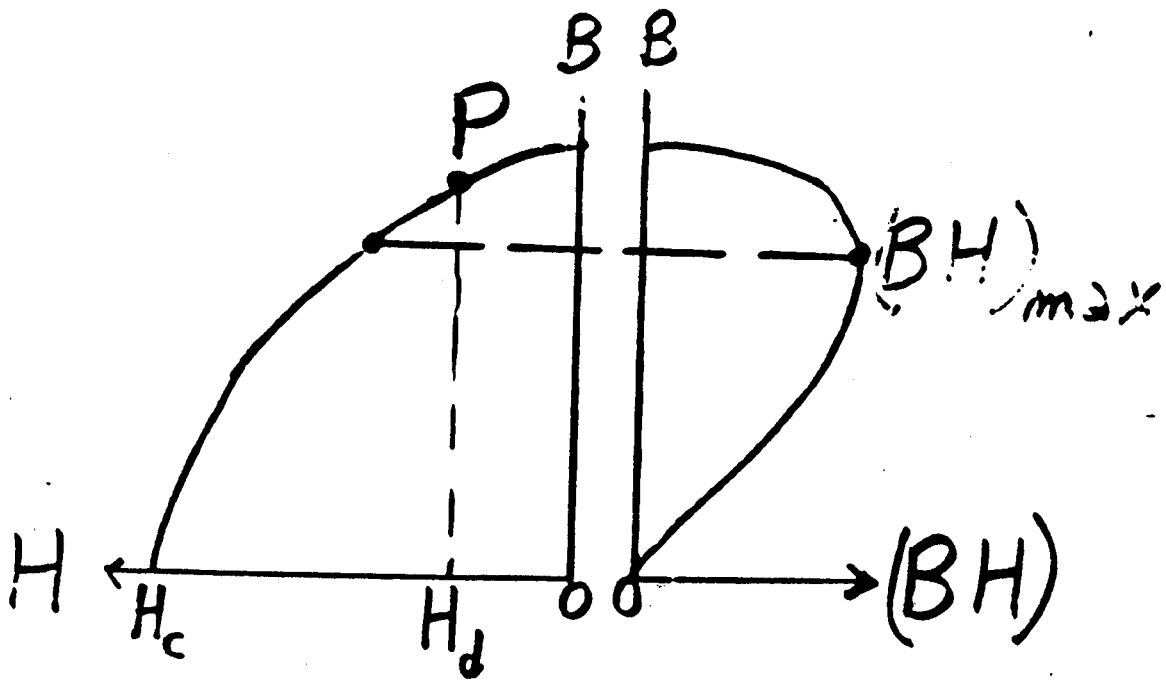
Stability re: H_0 and LINEARITY = $f(I H_c)$.

OPERATING POINT: P.



$$B = H_{\text{app.}} + H_0 + 4\pi M$$

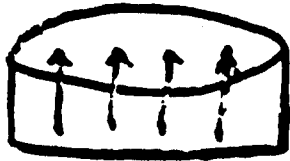
H_0 Shifts B_n to B_M .



i.e. WANT H_d LARGER \Rightarrow

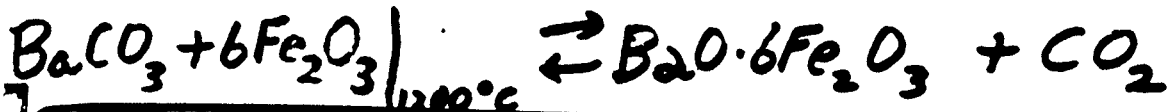
MAGNET THICKER OR Shorter

[NOTE:

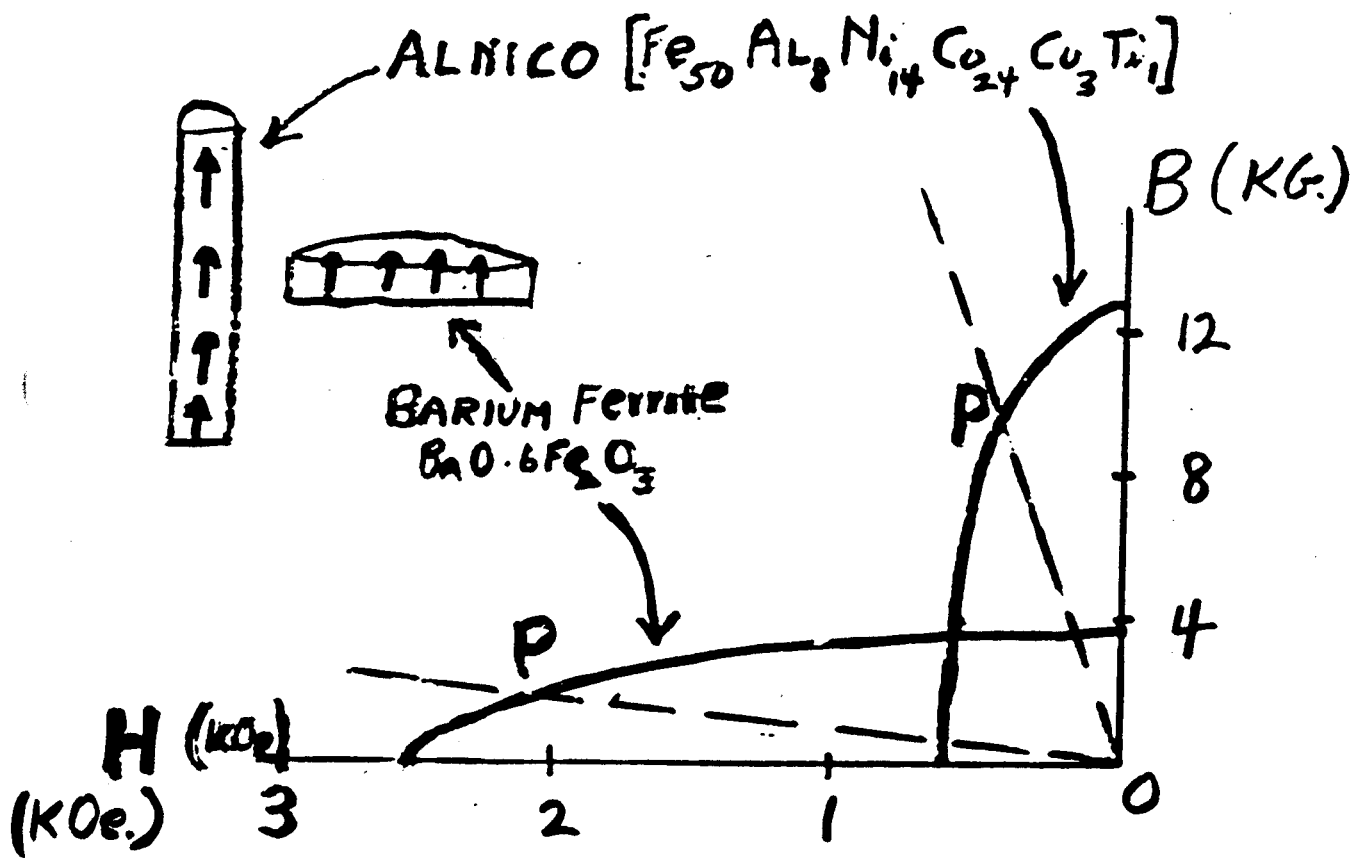


requires large H_c

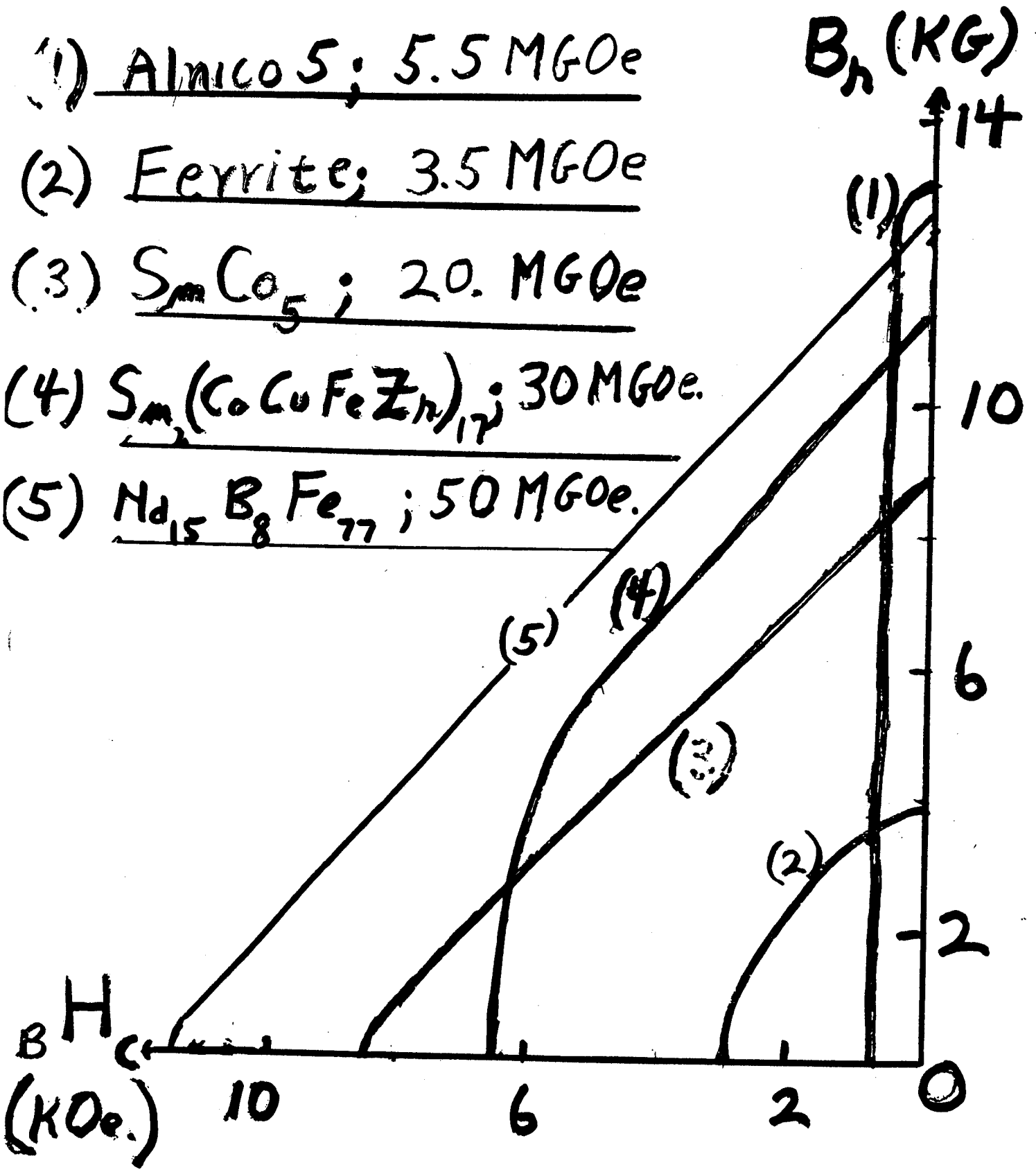
COMMON-USE-PERMANENT MAGNETS



	B_r	BH_c	$(BH)_{\text{max.}}$	iH_c
Barium Ferrite	3,950	2400	3.5	2470
Alnico - 5	13,000	640	5.5	SINTERED - CAST.



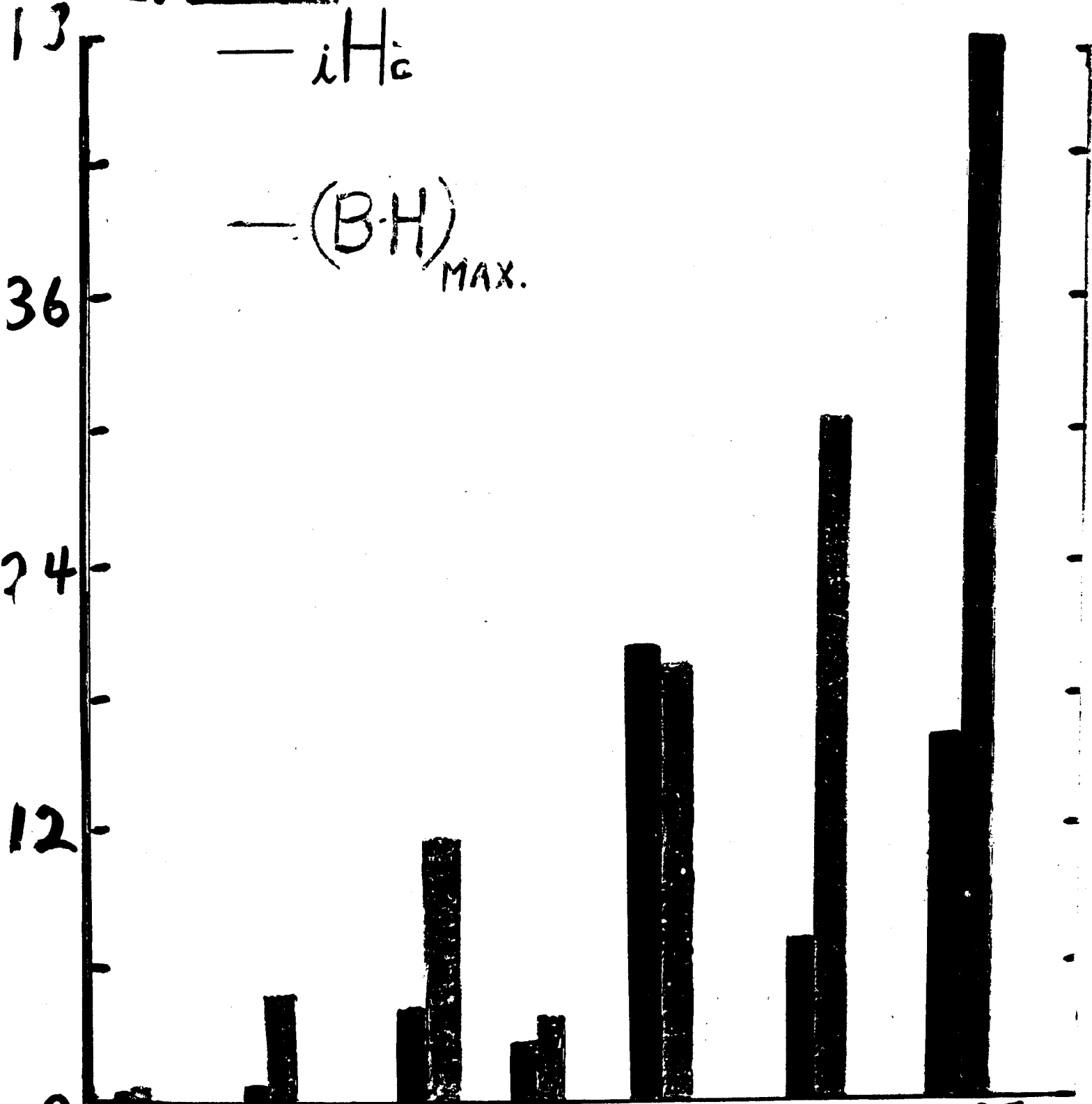
- (1) Alnico 5; 5.5 MG0e
- (2) Ferrite; 3.5 MG0e
- (3) S_mCo_5 ; 20. MG0e
- (4) $S_m(CoCuFeZn)_{17}$; 30 MG0e.
- (5) $Nd_{15}B_8Fe_{77}$; 50 MG0e.



KOe., MG Oe.

— iH_c

— (B·H)
MAX.



0 Steel 1885 Alnico 1938 CoPt 1952 BaFe 1954 SmCo₅ 1975 SmCo_{2,17} 1980 NdBFe 1986

Sm Co₅ - 1973

RARE EARTH (Sm) ⇒

CRYSTALLOGRAPHY - hcp

large ANISOTROPY - 10^8 ergs/cm³

Co₅ ⇒ MAGNETIZATION, $B_r = 9000$ G.

FINE-GRAINED (1-10 μm) MATERIAL }
WALL NUCLEATION CONTROLLED }

⇒ HIGH COERCIVE FORCE ; $H_c = 15$ kOe.

$(BH)_{MAX} \sim 20$ MG.Oe.

Curie Temp $\sim 750^\circ\text{C}$

SmCo₅ -

Powder Metallurgy

- 1) Melt Alloy
 - 2) Crush and Ball Mill \Rightarrow Powder
 - 3) Field Aligned and Compacted.
 - 4) Sintered at 1100°C in Argon ($\frac{1}{2}$ hr.)
-

Difficulties

- a) Reactivity of Samarium
 - b) Crystal Structure damage (2). $H_c \downarrow$
 - c) Low Packing Density of Aligned Compact
-

Plastic Bonded Magnets

Aligned SmCo₅ particles embedded
in flexible plastic MATERIAL

$\text{Sm}_2\text{Co}_{17}$ - 1978 - 1983

$\text{Sm}_2\text{Co}_{10}\text{Cu}_{1.48}\text{Fe}_{3.16}\text{Zr}_{.194}$ (TDK JAPAN)

$B_r = 11,200 \text{ G}$, $iH_c = 6,740 \text{ oe}$.

$(BH)_{\text{max}} = 30 \text{ MGoe}$, $T_c = 900^\circ\text{C}$

METALLURGY - SAME AS SmCo_5 except

STEP ANNEAL -

{ $850^\circ\text{C} \Rightarrow 500\text{\AA}$ CELLULAR STRUCTURE
 $600^\circ\text{C} \Rightarrow 2^{\text{nd}}$ PHASE (COPPER) IN CELL WALLS
(DOMAIN WALL PINNING)

SAMARIUM, COBALT - EXPENSIVE.

MATERIALS - CAST

$\text{Fe}_{77} \text{Nd}_{15} \text{B}_8$ Permanent Magnet (SUMITOMO)

Mat'ls melted, ground, and sintered.

3 μm non-magnetic neodymium rich precip in iron rich host matrix -

coercive force mechanism - Wall nucleation

	Fe Nd B	Sm Co
$B H_c$	12 kOe	6.5 kOe
$I H_c$	15 kOe.	15 kOe.
B_r	12.4 kG.	10 kG.
$(BH)_{max}$	50 MG Oe	30 MG Oe
Temp Coeff. of remnance	-0.13%/K	+0.02%/K

MATERIALS - VERY INEXPENSIVE

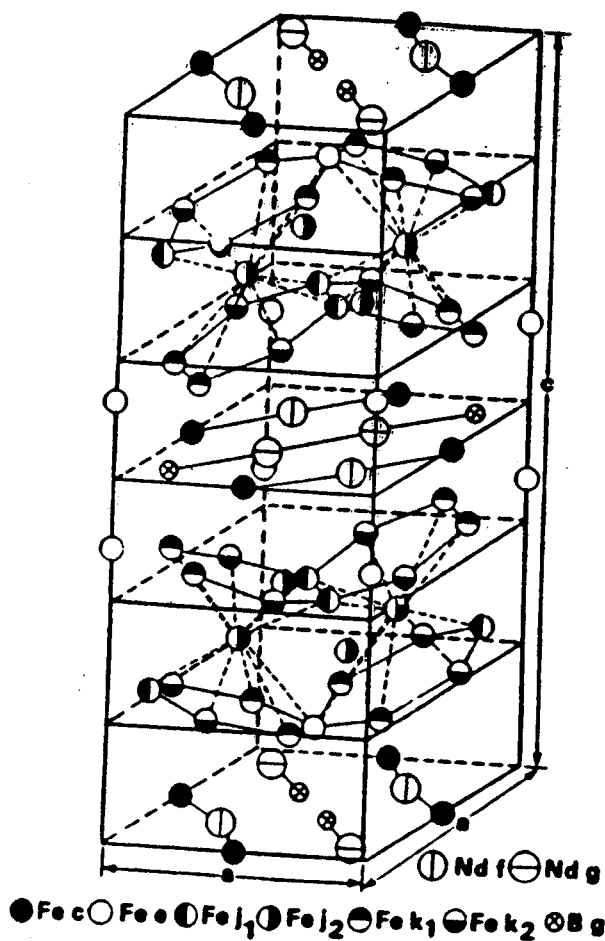


FIG. 1. Unit cell of $\text{Nd}_2\text{Fe}_{14}\text{B}$ ($P4_2/mnm$ space group). The c/a ratio in the figure is exaggerated to emphasize the puckering of the hexagonal iron nets.

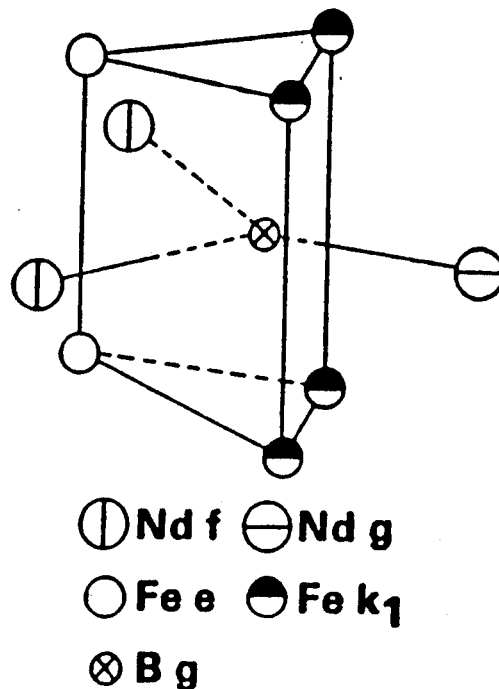
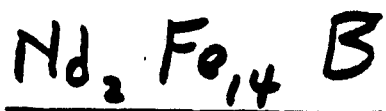


FIG. 2. Trigonal prism containing a boron atom in the $\text{Nd}_2\text{Fe}_{14}\text{B}$ structure. Distances between the $\text{B}(g)$ atom and its nearest neighbors are $\text{B}-\text{Fe}(k_1) = 2.09 \text{ \AA}$, $\text{B}-\text{Fe}(e) = 2.14 \text{ \AA}$, $\text{B}-\text{Nd}(g) = 2.86 \text{ \AA}$, and $\text{B}-\text{Nd}(f) = 3.34 \text{ \AA}$. The vertical edge lengths are $\text{Fe}(e)-\text{Fe}(e) = 2.75 \text{ \AA}$, $\text{Fe}(k_1)-\text{Fe}(k_1) = 3.12 \text{ \AA}$. Distances in the triangular faces are $\text{Fe}(e)-\text{Fe}(k_1) = 2.51 \text{ \AA}$, $\text{Fe}(k_1)-\text{Fe}(k_1) = 2.59 \text{ \AA}$.

TABLE I. Atomic sites, occupancies, and coordinates x , y , and z (in units of the lattice constants $a = 8.80 \text{ \AA}$, $c = 12.19 \text{ \AA}$) for $\text{Nd}_2\text{Fe}_{14}\text{B}$ obtained from analysis of room-temperature neutron-diffraction data. The space group is $P4_2/mnm$. The quality-of-fit index for the nuclear structure is $R(\text{nuc}) = 4.2\%$. These results imply a density of 7.60 g/cm^3 which agrees well with the measured value of 7.55 g/cm^3 .

Atom	Site	Occupancy	x	y	z
Nd	f	4	0.266	0.266	0.0
Nd	g	4	0.139	-0.139	0.0
Fe	k_1	16	0.224	0.568	0.128
Fe	k_2	16	0.039	0.359	0.176
Fe	j_1	8	0.097	0.097	0.205
Fe	j_2	8	0.318	0.318	0.247
Fe	e	4	0.5	0.5	0.113
Fe	c	4	0.0	0.5	0.0
B	g	4	0.368	-0.368	0.0



TETRAGONAL PRISMS

1) CENTER - BORON

2) BASAL PLANES - 3 IRON ATOMS
[Stretched vertically from hexagonal nets]

3) LINK IRON PLANES ABOVE + BELOW

4) COORDINATE ALL RARE-EARTH + BORON
{ more complete than $\text{Nd}_2\text{Fe}_{17}$ } $\left\{ \begin{array}{l} T_c = 310^\circ\text{C} \\ T_c = 155^\circ\text{C} \end{array} \right.$

5) ALSO IN TRANSITION METAL-METALLOID
{ Fe-B, Co-B, Ni-B, Fe-C, Pd-Si }
{ CRYSTALLINE AND AMORPHOUS }

MAGNETIC STRUCTURE

1) All Nd. + Fe MOMENTS FERROMAGNETIC

2) UNIAXIAL ANISOTROPY (|| c axis)

a) (CRYSTAL FIELD SPLITTING OF $4f$ Nd. level)

b) ($\sim 10\%$ from Fe sublattice itself -)

$$\text{e.g., } \left\{ \text{Gd}_2\text{Fe}_{14}\text{B} \Rightarrow H_c \sim 2\frac{1}{2} \text{ kOe.} \right\}$$

$$\left\{ \text{Nd}_2\text{Fe}_{14}\text{B} \Rightarrow H_c \sim 14 \text{ kOe.} \right\}$$