

X. Critical Phenomena: Ferromagnetism

A. Key Features in Ferromagnetism

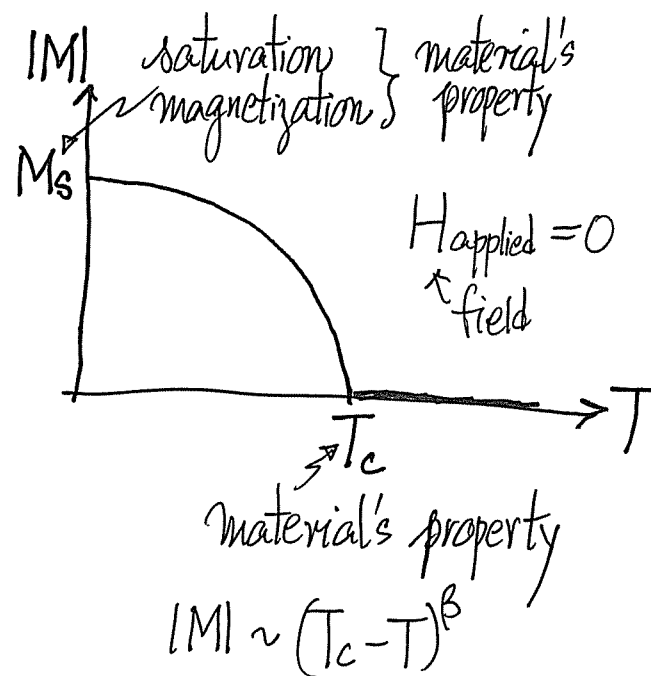
Spontaneous Magnetization (自發)

- No external applied magnetic field

$M \neq 0$ for $T < T_c$ ← Curie temperature
 "spontaneous" (∵ no applied field to magnetize sample)
 $M = 0$ for $T > T_c$

Material	M_s (10^6 A/m)	T_c (K)
Iron	1.75	1043
Cobalt	1.45	1404
Nickel	0.512	631
Gadolinium	2.00	289
Terbium	1.44	230
Dysprosium	2.01	85
Holmium	2.55	20

Sources: American Institute of Physics Handbook (D. W. Gray, Ed.) (New York: McGraw-Hill, 1963).



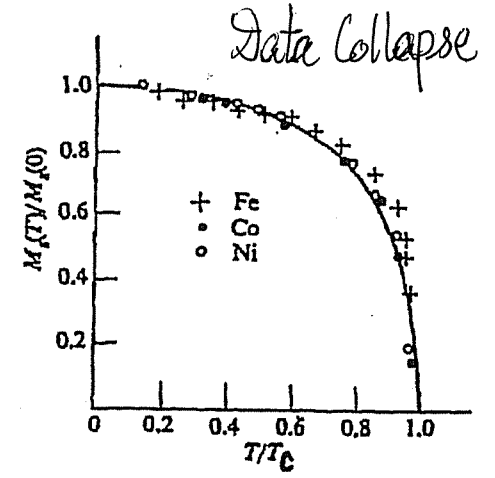
Critical point:
 $(T = T_c, H_{\text{applied}} = 0)$

- "Universal behavior" when properties of different Ferromagnetic (FM) materials are viewed in reduced quantities

FM material A: Measure $M(T)$
 $\Rightarrow M_S^{(A)}, T_c^{(A)}$

FM material B: Measure $M(T)$
 $\Rightarrow M_S^{(B)}, T_c^{(B)}$

Look at $\frac{M(T)}{M_S}$ vs $\frac{T}{T_c}$
 Almost the same curve for different materials



- For $T \lesssim T_c$, $M \sim (T_c - T)^\beta$ and same β for many different materials [c.f. vapor-liquid case][†]
 - M is the order parameter of FM transition
 - M changes continuously at $T_c \Rightarrow$ "Continuous phase transition" (transitions that are NOT first order)
- $T < T_c, M \neq 0$ Ferromagnetic phase (more ordered)
 $T > T_c, M = 0$ Paramagnetic phase (disordered)

[†] Recall the law of corresponding state. Here is another example.

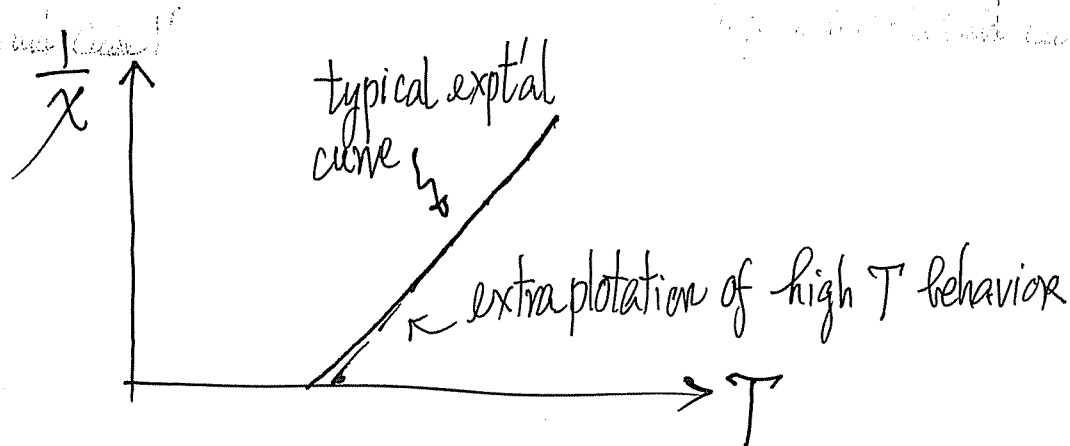
Paramagnetic behavior at $T > T_c$

$T > T_c$, $H_{\text{applied}} = 0$, $M = 0$ (no spontaneous magnetization)

With an applied field, $M \neq 0$ when $H_{\text{applied}} \neq 0$

Response is described by $M = \chi H_{\text{applied}}$

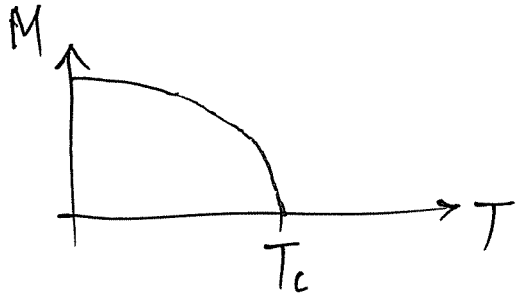
χ follows $\chi = \frac{C}{T - T_c}$ similar to Curie's law [recall $\chi \sim \frac{1}{T}$ for paramagnetic behavior]



It says, χ gets bigger and bigger as $T \rightarrow T_c$ from above.

As $T \rightarrow T_c$, χ diverges. OK! $M = \chi \underbrace{H_{\text{applied}}}_{=0} \Rightarrow \chi$ diverges when sample becomes FM!

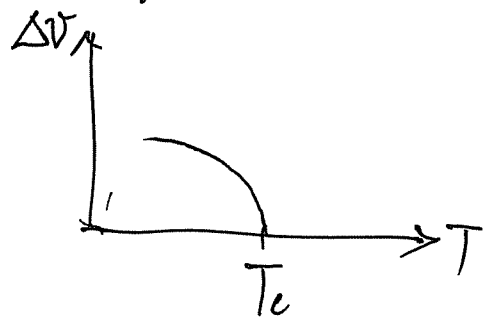
Ferromagnetic-Paramagnetic Transition



Ordered \leftrightarrow Disordered

$$M \sim (T_c - T)^\beta$$

Liquid-vapor Transition



Ordered - Disordered

$$\Delta v \sim (T_c - T)^{\beta'}$$

It turns out that very different physical scenarios may carry the same value of critical exponent!

This is what physicists meant by universal behavior.

[Magnetic system and liquid-vapor system behave the same way near the critical point] (See Appendix A for Percolation Problem)

B. Hints from theory of paramagnetism

Think like a physicist!

atoms \rightarrow tiny magnets (magnetic moments)

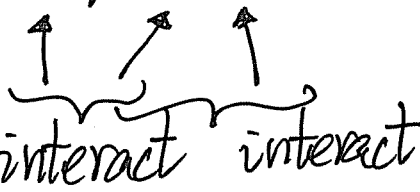
$M \neq 0 \Rightarrow$ moments have tendency to align (even $H_{\text{applied}} = 0$)

Paramagnetism: No magnetic moment - magnetic moment interaction
 & $\vec{\mu}$ only responds to \vec{H}_{applied}

Ferromagnetism: $\vec{\mu}$'s aligned! $\vec{\mu}$ must "feel" some local or internal

magnetic field!

How come?



Interaction between neighboring magnetic moments

a moment wants neighboring moments to point to its direction (ferromagnetic interaction)

Interaction: tend to align
 kT : tend to randomize

low T or strong interaction \Rightarrow alignment
 $\Rightarrow M \neq 0$

How strong is the interaction?

Data: $T_c(\text{iron}) = 1043 \text{ K} \sim 10^3 \text{ K}$
 $T_c(\text{nickel}) = 631 \text{ K}$ } Meaning: kT_c is not sufficient
 to randomize alignment

2 to 3 times of room temperature

$kT_c \sim 0.05$ to 0.1 eV

strong interaction! [Much stronger than EM theory predicts]

The points are:

- Data indicate strong moment-moment interaction
 \Rightarrow alignment gains much energy
- Interaction has quantum origin
 \Rightarrow exchange interaction (due to anti-symmetric electron-wavefunction)
 electrons are fermions (a quantum idea)

1-page review on Paramagnetism

$J = 1/2$ ($S = 1/2$)
 $- +\mu_B B$ (anti-align with applied B)
 (two-level system)

 $- -\mu_B B$ (align with applied B)

$$Z = z^N, \quad z = e^{\beta\mu_B B} + e^{-\beta\mu_B B} = 2 \cosh\left(\frac{\mu_B B}{kT}\right)$$

$\langle \mu_z \rangle = \mu_B \tanh\left(\frac{\mu_B B}{kT}\right)$
 one moment

Whole system: $N \langle \mu_z \rangle = N \mu_B \tanh\left(\frac{\mu_B B}{kT}\right)$

Per unit volume $M = \frac{N}{V} \langle \mu_z \rangle = \frac{N}{V} \mu_B \tanh\left(\frac{\mu_B B}{kT}\right) = M_s \tanh\left(\frac{\mu_B B}{kT}\right)$

saturation magnetization
 ↓
 B-field acting on each independent (non-interacting) magnetic moment
 same as $B_J(x)$ with $J = 1/2$

The underlying Hamiltonian is:

$$H_{\text{para}} = \sum_i -\vec{\mu}_i \cdot \vec{B}_{\text{applied}} \quad \text{with } \mu_{iz} = \begin{cases} +\mu_B & \text{(aligned with } \vec{B}_{\text{applied}}) \\ -\mu_B & \text{(anti-aligned with } \vec{B}_{\text{applied}}) \end{cases}$$

(cover all magnetic moments)

More on Paramagnetic Hamiltonian

The last equation of H_{para} may as well be written as:

both forms are used in textbooks and journal articles

$$H_{para} = -(\mu_B B_{applied}) \sum_i \sigma_i \quad \text{with } \sigma_i = \begin{cases} +1 & \text{(align with } B_{applied}) \\ -1 & \text{(anti-align with } B_{applied}) \end{cases}$$

(cover all magnetic moments)

$$= -(\mu_B B_{applied}) \sum_i S_i$$

(cover all magnetic moments)

an energy indicating how strong $B_{applied}$ is

Sometimes (e.g. in Ising model literature), this term is written as

$$H_{para} = -H \sum_i S_i \quad \text{OR} \quad -H \sum_i \sigma_i \quad \text{OR} \quad -B \sum_i \sigma_i \quad \text{OR} \quad -B \sum_i S_i$$

an energy representing $\mu_B B_{applied}$ an energy representing $\mu_B B_{applied}$

[unfortunate symbol, as H means many other things!]
(but must get used to it!)

C. Modelling: An art and a science

- Good model:
 - { include the essential physics in simplest form
 - { fewer symbols / simplified notations
 - { non-trivial results (e.g. ferro to para transition)

Motivating a model:

- Essential physics = $\vec{\mu}_i$ and $\vec{\mu}_j$ interaction $\ll kT$

FM interaction: $\vec{\mu}_i$ and $\vec{\mu}_j$ align (μ_{iz} and μ_{jz} align)

\Rightarrow lower energy

$\vec{\mu}_i$ and $\vec{\mu}_j$ anti-align (μ_{iz} and μ_{jz} anti-align)

\Rightarrow higher energy

In general, $\vec{\mu}_i$ ($\vec{\mu}_j$) could have many \hat{x} -components (J quantum number).

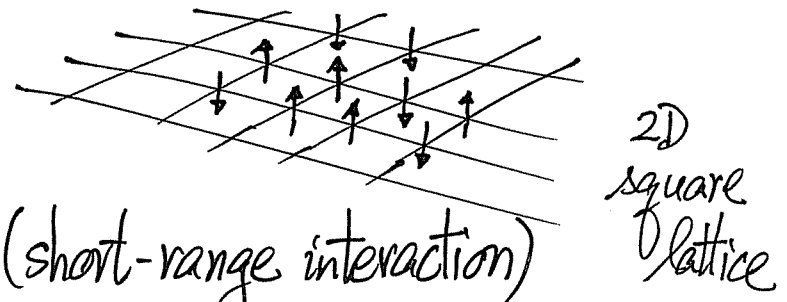
Why bother! Grasp the essential physics $\left\{ \begin{array}{l} +\mu_B \\ -\mu_B \end{array} \right.$ OR $J = 1/2$ OR $S = 1/2$ only

Simplest model : μ_z takes on only $\left\{ \begin{matrix} \mu_B \text{ ("up")} \\ -\mu_B \text{ ("down")} \end{matrix} \right\}$ OR $\left\{ J_z = \pm \frac{1}{2} \right\}$ OR $\left\{ S_z = \pm \frac{1}{2} \right\}$

[Including more components will only add details]

Nearest-neighbor interaction only

- Interact only when two magnetic moments are nearest neighbors (short-range interaction)



[Including longer-range interactions will only add details]

Big physics = short-range interaction can lead to long-range order

\hat{x} -component aligned : $\uparrow \uparrow$ and $\downarrow \downarrow \Rightarrow E_{\text{lower}}$

\hat{z} -component anti-aligned : $\uparrow \downarrow$ and $\downarrow \uparrow \Rightarrow E_{\text{higher}}$

$\begin{matrix} i & j \\ \uparrow & \uparrow \end{matrix}$

 $\begin{matrix} i & j \\ \uparrow & \uparrow \end{matrix}$

nearest neighbors nearest neighbors

differ by
 $E_{\text{higher}} - E_{\text{lower}} = "2J"$

In physical units, the interaction energy is of the form

$$-J_{ij} \vec{\mu}_i \cdot \vec{\mu}_j$$

to make sure the expression is an energy
↑
"± μ_B²" for { aligned
anti-aligned

Simplify notations

- ↑ ↑ as +1 +1 or + + | ↓ ↓ as -1 -1 or - -
up up up up down down down down
- ↑ ↓ as +1 -1 or + - | ↓ ↑ as -1 +1 or - +
up down up down down up down up

Recall: Pauli matrices (spin-half)

$$\sigma_z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix} \Rightarrow \text{eigenvalues} = \begin{matrix} +1 \\ \uparrow \\ \text{up} \end{matrix}, \begin{matrix} -1 \\ \downarrow \\ \text{down} \end{matrix}$$

Write interaction energy as

$$S_i \quad S_j \quad \boxed{-J_{ij} S_i S_j}$$

with $\begin{cases} S_i \text{ taking on } +1 \text{ or } -1 \\ S_j \text{ taking on } +1 \text{ or } -1 \end{cases}$

aligned: $(+1, +1)$ and $(-1, -1)$ give $S_i S_j = 1$ and energy $-J_{ij}$

anti-aligned: $(+1, -1)$ and $(-1, +1)$ give $S_i S_j = -1$ and energy $+J_{ij}$ } differ by $2J_{ij}$

• J_{ij} is an energy characterizing the strength of interaction

• $J_{ij} > 0$, S_i and S_j tend to align \Rightarrow FM interaction

[$J_{ij} < 0$, S_i and S_j tend to anti-align \Rightarrow Anti-ferromagnetic interaction]

Assuming $J_{ij} = J$ (same) for all nearest-neighbors (ij) , the interaction energy (Hamiltonian) is

$$E(\{S_i\}) = -J \sum_{\langle ij \rangle} S_i S_j$$

sum over all distinct nearest-neighbor pairs

Ising model
with no external
applied magnetic
field

In general, there may also be an applied (external) \vec{B}_{applied} .

An additional energy due to $(-\vec{\mu}_i \cdot \vec{B}_{\text{applied}})$ and sum up all i .

$$\vec{B}_{\text{applied}} = B_{\text{applied}} \hat{z} \quad (\text{can always call the direction } \hat{z})$$

$$\text{Zeeman energy} = \underbrace{-\mu_B B_{\text{applied}}}_{\substack{\uparrow \\ \text{energy} \\ \text{scale}}} \sum_i \underbrace{S_i}_{\substack{\uparrow \\ \text{add up} \\ \text{all magnetic moments}}}$$

\uparrow contributing $-\mu_B B_{\text{applied}}$ (aligned)
 \uparrow contributing $+\mu_B B_{\text{applied}}$ (anti-aligned)
 $(S_i = +1 \text{ or } -1)$

Simplify notations

$$\text{Zeeman energy} = E_B(\{S_i\}) = \underbrace{-B}_{\substack{\uparrow \\ \text{an energy } (\mu_B B_{\text{applied}}) \text{ characterizing} \\ \text{how strong the applied field is.}}} \sum_i S_i$$

D. Ising Model

$$E(\{S_i\}) = -J \sum_{\langle ij \rangle} S_i S_j - B \sum_i S_i \quad (1)$$

interaction
between magnetic moments
(or people called them spins)

energy representing
external applied field

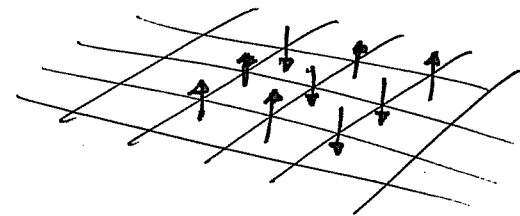
- this competes with kT
- S_i takes on $+1$ or -1
- When $B=0$, can spins align even at finite temperature?
- Can study Ising model on 1D chain (exactly solvable),
2D lattices (some exactly solvable), 3D lattices (no exact solution),
4D lattices, ...
- Lenz (1920) constructed the model for Ising (1925) to study in his thesis

Reminder:

- $J=0$, $E(\{S_i\}) = -B \sum_{i=1}^N S_i$ (N spins or N dipole moments)
is the paramagnetic problem studied earlier
- $\langle S_i \rangle$ is an average number between -1 to +1
- Previously in paramagnetism, $\langle \mu_z \rangle = \mu_B \langle S_i \rangle$
- Now, with $J \neq 0$ in the Ising Model, we also want to find $\langle S_i \rangle$, which is proportional to the Magnetization.

E. Ising Model: What can be done, formally?

$$E(\{S_i\}) = -J \sum_{\langle ij \rangle} S_i S_j - B \sum_i S_i \quad (1)$$



Consider: N moments (spins) on a 2D square lattice
 In Stat. Mech., want to evaluate $Z = \sum_{\text{all } N\text{-spin states } \{S_i\}} e^{-\frac{E(\{S_i\})}{kT}}$

- What to sum over?
 - 2^N strings of the form $\{S_1, S_2, \dots, S_N\}$ with $S_i = \begin{cases} +1 \\ -1 \end{cases}$
- What goes into $e^{-\frac{E(\{S_i\})}{kT}}$?
 - For every string $\{S_i\}$, calculate $E(\{S_i\})$ from Eq. (1) and evaluate one term in Z
 - Repeat for each of 2^N strings (2^N is a huge number for, say, $N=100^2$)
 - Done! In principle.

Can this be done?

- Analytically?
 - 1D chain: Yes (see Problem Set 3)
 - 2D square lattice: Yes (but not so easy)
 - 3D simple cubic or other lattice: No!
- Numerically?
 - Any dimension? (Write a program to evaluate $Z(T, N, B)$ exactly?)
[huge # of $\{S_i\}$'s]
 - Any dimension? (An algorithm to carry out the importance sampling implied by the canonical ensemble)
[Monte-Carlo simulation]
- Approximately?
 - Mean field theories
 - $1 + \epsilon$ ($\epsilon \ll 1$) dimension; $4 - \epsilon$ ($\epsilon \ll 1$) dimension
 - Renormalization methods