

Project for the exam: AMS (DSE)

Paola Serra and Marzio De Corato

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# Theoretical Framework

# Statistical Mechanics

"Ludwig Boltzmann, who spent much of his life studying statistical mechanics, died in 1906, by his own hand. Paul Ehrenfest, carrying on the work, died similarly in 1933. Now it is our turn to study statistical mechanics." States of Matter (1975), by David L. Goodstein

# Concepts of statistical mechanics: aim [5]

**Aim** To predict the macroscopic properties of systems on the basis of their microscopic structure. This connection between these two scales is performed with statistical methods

# Concepts of statistical mechanics: entropy [5]

**Central problem of thermodynamics:** characterize the actual state of equilibrium among all virtual states

**Entropy postulate**: there exist a function S of the extensive variables  $(X_0, X_1...X_r)$  called entropy, that assumes the maximum value for a state of equilibrium among all vritual states and that possesses the following properties:

- Extensivity  $S^{(1\cup 2)} = S^1 + S^2$
- Convexity  $S((1-\alpha)X^1 + \alpha X^2) \ge (1-\alpha)S(x^1) + \alpha S(X^2)$
- Monotonicity  $\frac{\partial S}{\partial F}|_{X_1...X_r} = \frac{1}{T} > 0$

The equilibrium state corresponds to the maximum entropy compatible with the constrains

# Concepts of statistical mechanics: entropy [5]

• Fundamental postulate of statistical mechanics  $S = k_b \ln |\Gamma|$ Where S is the thermodynamic entropy,  $k_b$  is Boltzmann constant and  $|\Gamma|$  the volume in the phase space

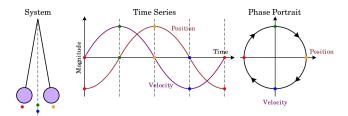


Image taken from [1]

# Concepts of statistical mechanics: ensemble [5]

- Statistical ensemble: a large number of virtual copies of a system; each of them is a possible state of the real system (epistemic probability). It is the formalization of a repeated experiment proposed by Gibbs (empirical probability)
- Microcanical ensamble: p = 1/W W is the number of microstates
- Canonical ensamble:  $p = \frac{1}{Z} \exp\left(-\frac{E}{kT}\right)$  where  $Z = \sum_{i} \exp\left(\frac{-E_{i}}{k_{b}T}\right)$

$$\langle A(x) \rangle = \frac{1}{Z} \int dx_s A(x_s) exp \left[ -\frac{H^{(S)}(x_s)}{k_b T} \right]$$
 (1)

$$Z = \int dx_s \exp\left[-\frac{H^S(x_s)}{k_b T}\right]$$
 (2)

$$\langle H(x)^2 \rangle - \langle H(x) \rangle^2 = k_b T^2 C$$
 (3)

# Ising Model

# Ising model [4, 3]

- An array of atoms that can take states  $\pm 1$ . The energy of the system is given by  $E(\mathbf{x},J,H) = -\left[\frac{1}{2}\sum_{m,n}J_{mn}x_mx_n + \sum_nHx_n\right]$  where J is the coupling constant between two neighbour sites, and H is an external field.
- The probability of the system to be in the state x is given by  $p(x|\beta, J, H) = \frac{1}{Z(\beta, J, H)} \exp[-\beta E(x, J, H)]$  (canonical ensemble) where  $\beta = 1/k_b T$   $Z(\beta, J, H) = \sum_x exp[-\beta E(x, J, H)]$
- It is useful to characterize the order level of a lattice (macroscopic) with the (spatial) correlation functions (whose input are microscopic quantities). In particular, for the Ising model, these are given by the following expression (with H=0)

$$g(m) = \frac{\langle \sigma_i \sigma_{i+m} \rangle - \langle \sigma_i \rangle \langle \sigma_{i+m} \rangle}{1 - \langle \sigma_i \rangle \langle \sigma_{i+m} \rangle} = \langle \sigma_i \sigma_{i+m} \rangle$$

# Numerical simulations

"Never make a calculation until you know the answer. Make an estimate before every calculation, try a simple physical argument (symmetry! invariance! conservation!) before every derivation, guess the answer to every paradox and puzzle. Courage: No one else needs to know what the guess is. Therefore make it quickly, by instinct. A right guess reinforces this instinct. A wrong guess brings the refreshment of surprise. In either case life as a spacetime expert, however long, is more fun!" John Archibald Wheeler

# Numerical simulation [5]

- Molecular dynamics: the equation of motion are solved numerically PROS: information of both the dynamical and static properties of the system are explored
- Monte Carlo: a fictitious evolution process of the system is solved in order to get the equilibrium distribution PROS 1) also the systems whose dynamics is not defined can be explored 2) a fictitious dynamics can be considered in order to reach the equilibrium faster

# Monte Carlo method [5]

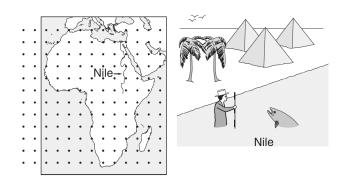


Image taken from [2]

# Monte Carlo method [5]

- We would calculate an integral of type  $\langle A \rangle = \int_0^1 dx A(x) \rho(x)$  where  $\rho$  is the probability distribution.
- Evaluate the integrand in N+1 points uniformly arranged between 0 and 1  $\langle A \rangle \approx \frac{1}{N+1} \sum_{i=0}^{N} A(x_i) \rho(x_i)$
- A better convergence is reached if the  $x_i$  density is proportional to  $\rho(x)$ . In this case we have  $\langle A \rangle = 1/N \sum_{i=1}^N A(\mathbf{x}_i)$ .

# Monte Carlo method [2]

- Lets consider the generic integral  $I = \int_a^b dx \ f(x)$
- This can be recast in the following form  $I = \int_0^1 dx \ w(x) \frac{f(x)}{w(x)}$
- If w(x) is the derivative of u(x) (non-decreasing,non negative) we have  $I=\int_0^1 du \; \frac{f[x(u)]}{w[x(u)]}$
- If one considers L random values of u uniformely distributed in the interval [0,1] we have  $I \approx \frac{1}{L} \sum_{i=1}^{L} \frac{f[x(u)]}{w[x(x)]}$
- ullet The choice of w is crucial since  $\sigma=rac{1}{L}\left[\left\langle \left(rac{f}{w}
  ight)^2 
  ight
  angle \left\langle rac{f}{w} 
  ight
  angle^2
  ight]$
- Brute Force  $f=10^{-260}$  and  $\sigma=\frac{1}{Lf}$ ...is not a good idea
- PROBLEM: we do not know the form of the denominator (if we know it we do not need the Monte Carlo method)

# Monte Carlo method: Metropolis idea [5]

- The quantity  $\langle A \rangle$  is evaluated over a distribution  $\rho(x)$  where  $x_i$  are random, independent and distributed with a probability density  $\rho(x)$
- The independence conditions can be relaxed with the condition that the correlations between  $x_i$  and  $x_{i+1}$  go to zero fairly rapidly as I grows
- The true dynamics is replaced with a fictitious stochastic dynamics. The state at t+1 depends only from the state at t.  $\rightarrow$  Markov Chain

# Monte Carlo method: Metropolis idea [5]

- The evolution of probability is described by the **master equation**  $\Delta p_a(t) = \sum_{b\neq a}' [W_{ab}p_b(t) W_{ba}p_a(t)]$
- stationary  $o \sum_{b 
  eq a}' [W_{ab} p_b(t) W_{ba} p_a(t) = 0 \quad orall a$
- ullet Detailed balance property  $W_{ab}W_{bc}W_{ca}=W_{ac}W_{cb}W_{ba}$
- $W_{ab}p_b(t) W_{ba}p_a(t) = 0 \quad \forall a, b$

# Monte Carlo method: Metropolis idea [5]

- We would sample  $p_a^{eq}$
- This can be performed as long as the  $W_{ab}$  is ergodic and the detailed balance property holds
- The transition between any two arbitrary states can take place as long as one waits for a sufficient amount of time

# Application [5]

- $H(\sigma) = -\sum_{\langle ij \rangle} J\sigma_i \sigma_j \sum_i h\sigma_i$
- $P_{\sigma} = \frac{e^{-H(\sigma)/k_bT}}{Z}$   $Z = \sum_{\sigma} e^{-H(\sigma)/k_bT}$
- The observable are calculated as

$$E = \langle H \rangle = \sum_{sigma} H(\sigma) P^{B}_{\sigma'} \quad M = \sum_{\sigma} (\sigma_{i}\sigma) P^{B}_{\sigma}$$

- The markov chain states should be the microstates  $\sigma$  and  $P_{\sigma}$  the stationary distribution
- $W_{\sigma\sigma'} = W_{\sigma'\sigma} \frac{P_{\sigma}}{P_{\sigma'}} = W_{\sigma'\sigma} \exp{-\frac{H(\sigma) H(\sigma')}{k_b T}}$
- The Z term is no more present !!!

# Application [5]

• 
$$W_{\sigma'\sigma} = \begin{cases} \kappa H(\sigma') < H(\sigma) \\ \kappa \exp\left\{-\left[H(\sigma') - H(\sigma)\right]/k_bT\right\} \end{cases} \quad \{A\} \neq [A]_T = \frac{1}{T}$$

- $A(\sigma) = \langle A \rangle \left[ 1 + O(N^{-1/2}) \right] \forall \sigma \in \Gamma$
- $S = 0.5Nk_b ln2$   $\frac{|\Gamma|}{2^N} \approx 2^{-0.5N}$  N = 100  $10^{-15}$
- $\approx [A]_T = \frac{1}{T} \sum_{t=T_0}^{T\langle A \rangle_0 + T} A_{\sigma(t)}$
- $\langle \Delta A_T^2 \rangle \approx \frac{1}{T} \left( A_{\sigma(t)} [A]_T \right)^2$
- $\sigma_t$  and  $\sigma_{t'}$  are indipendent if |t-t'| is larger that a characteristic time  $au_0$
- $\left\langle \Delta A_T^2 \right\rangle \approx \frac{\tau_0}{T} \left( A_{\sigma(t)} [A]_T \right)^2$

# Goals and methods

#### Goals and methods

- Reproduce the main result for a 2D anti ferromagnetic lattice (J=-1) with no external magnetic field (H=0) with the montecarlo-metropolis
- Once checked that the script provide the correct results apply it to a lattice (J=+1). In this case the spins represent an opinion and the sites people. The goal is to find the stationary states (at T=0 and  $T\neq 0$ )
- Introduce in the lattice some blocks that never change their status. These islands represent groups that never change mind and only diffuse their ideas. (at T=0 and  $T\neq 0$ )

# Results

#### Simulation features

- 10x10 lattice
- ullet Periodic boundary conditions o the topology of a torus (genus equal to 1)
- 6000 steps

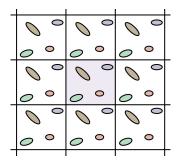


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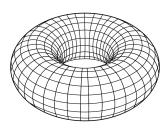
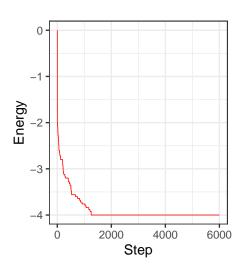
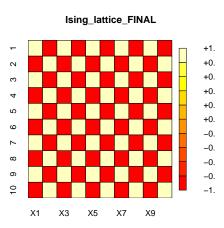
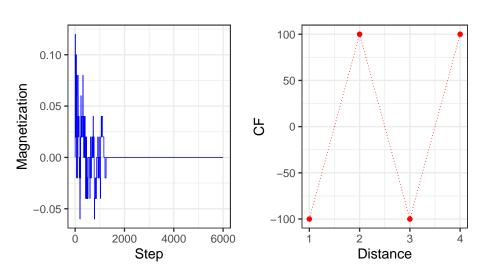
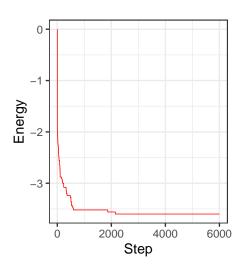


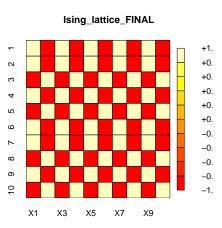
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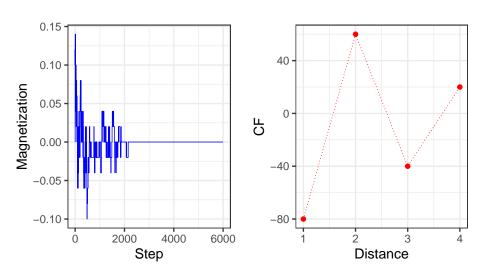


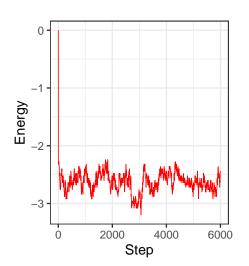


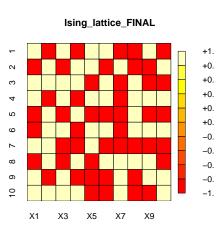


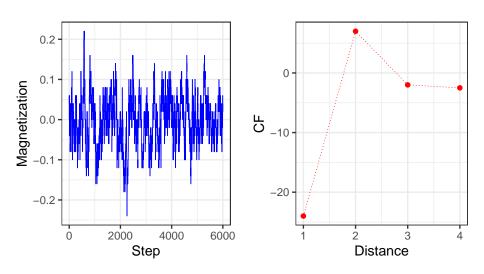


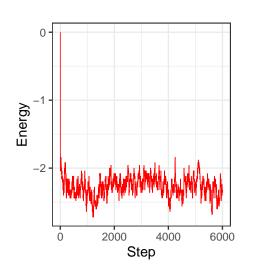


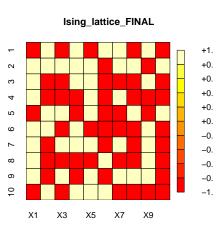


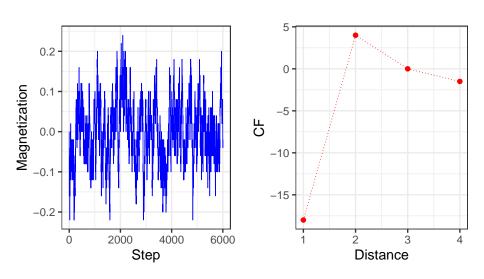




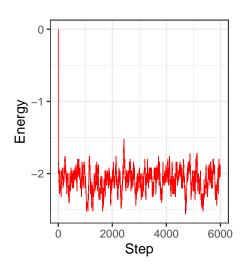


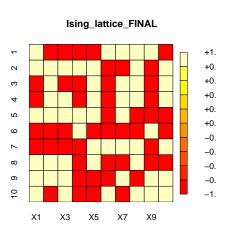




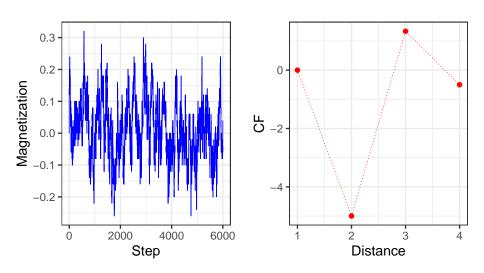


### Ferromagnetic J=-1, T=30

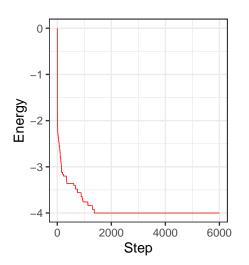


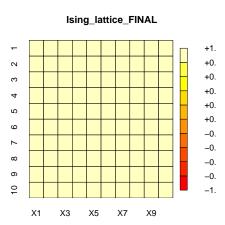


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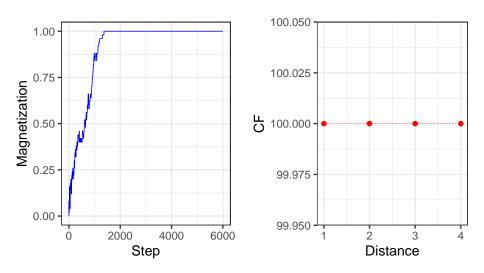


### Ferromagnetic J=+1, T=0

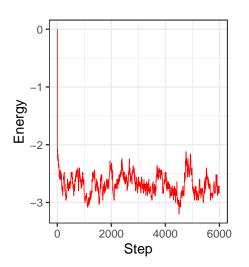


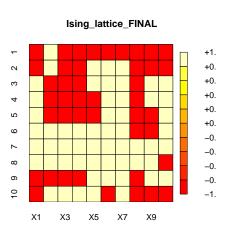


# Ferromagnetic J=+1, T=0

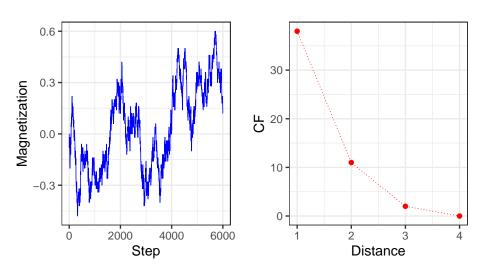


# Ferromagnetic J=+1, T=8

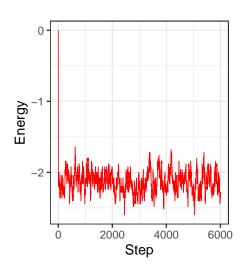


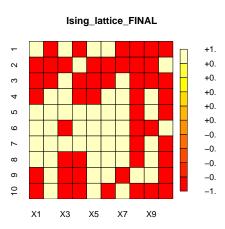


# Ferromagnetic J=+1, T=8

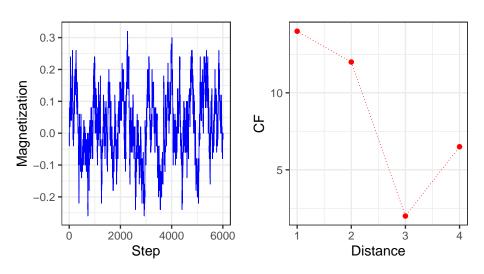


# Ferromagnetic J=+1, T=16

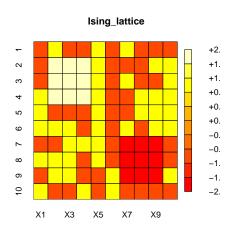




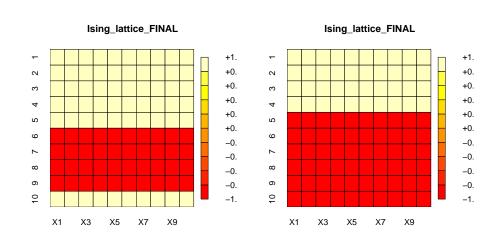
# Ferromagnetic J=+1, T=16



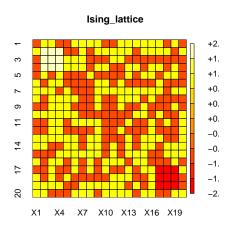
#### Ferromagnetic with two centers T=0



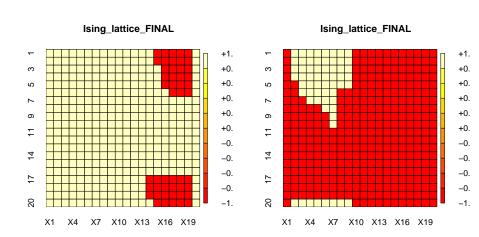
#### Ferromagnetic with two centers T=0



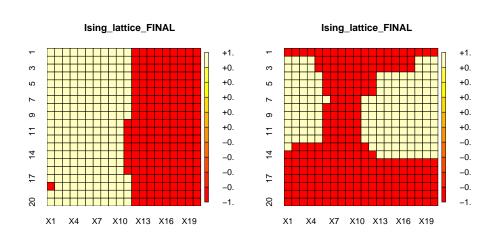
#### Ferromagnetic with two centers T=0, $20\times20$



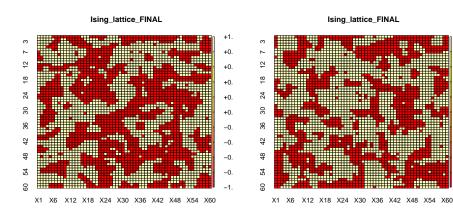
## Ferromagnetic with two centers T=0, $20\times20$



## Ferromagnetic with two centers T=0.25, $20\times20$



#### Ferromagnetic with two centers T=0, $60\times60$



+1.

+0. +0. +0.

+0.

+0.

-0. -0.

-0.

## Bibliography I

- [1] https://commons.wikimedia.org/wiki.
- [2] Daan Frenkel and Berend Smit. *Understanding molecular simulation:* from algorithms to applications. Vol. 1. Elsevier, 2001.
- [3] Jeffrey C. Grossman and Elif Ertekin, NSE C242 Phys C203, Spring 2008, U.C. Berkeley, howpublished = https://nanohub.org/resources/4579/download/05-01-08-forweb.pdf,
- [4] David JC MacKay and David JC Mac Kay. *Information theory, inference and learning algorithms*. Cambridge university press, 2003.
- [5] Luca Peliti. *Statistical mechanics in a nutshell*. Princeton University Press, 2011.

# Concepts of statistical mechanics: entropy [5]

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**Entropy postulate**: there exist a function S of the extensive variables  $(X_0, X_1...X_r)$  called entropy, that assumes the maximum value for a state of equilibrium among all vritual states and that possesses the following properties:

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The equilibrium state corresponds to the maximum entropy compatible with the constrains

# Concepts of statistical mechanics: entropy [5]

- Fundamental postulate of statistical mechanics  $S=k_b \ln |\Gamma|$
- Where S is the thermodynamic entropy,  $k_b$  is Boltzmann constant and  $|\Gamma|$  the volume in the phase space

$$S(X_{0},...,X_{r}) = k_{b} \ln \int_{\Gamma} dx = k_{b} \int dx \prod_{i=0}^{r} \left[\theta(X_{i}(x) - (X_{i} - \Delta X_{i}))\theta(X_{i} - X_{i}(x))\right]$$
(4)

- $(X_0, ..., X_r)$  are the extensive variables
- The  $\theta$  functions assures that the integrand is not null only in the interval  $X_i \Delta X_i \leq X_i(x) \leq X_i$

# Concepts of statistical mechanics: micro-canonical ensamble

Lets focus on a particular observable A (extensive)

$$S(X; a) = k_b \ln \int_{\Gamma} dx \delta(A(x) - a)$$
 (5)

$$S(X) = S(X; a*) \ge S(X; a) \tag{6}$$

$$\frac{|\Gamma(a)|}{|\Gamma|} = \frac{1}{|\Gamma|} \int_{\Gamma} dx \delta(A(x) - a))$$

$$= \exp\left\{\frac{1}{k_b} \left[S(X; a) - S(X; a^*)\right]\right\}$$

$$\simeq \exp\left\{\frac{1}{k_b} \left[\frac{\partial^2 S}{\partial A^2}|_{a^*} (a - a^*)^2\right]\right\}$$

$$a^* = \langle A(x) \rangle = \frac{1}{|\Gamma|} \int_{\Gamma} dx A(x)$$
(8)

#### Concepts of statistical mechanics: canonical ensemble

$$a^* = \frac{1}{|\Gamma|} \int_{\Gamma} dx_s dx_R A(x_s) \tag{9}$$

$$\langle A(x) \rangle = \int dx_s dx_r A(x_S) \delta(H^{(s)})$$
 (10)

$$\langle A(x) \rangle = \frac{1}{|\Gamma|} \int dx_s dx_r A(x_s) \delta(H^S(x_s) + H^R(x_R) - H^S(x_S)) \tag{11}$$

$$\langle A(x) \rangle = \frac{1}{|\Gamma|} \int dx_s A(x_s) \times \int dx_r \delta(H^R(x_R) - (E - H^{(S)}(x_s)))$$
 (12)

$$\int dx_r \delta(H^R(x_R) - (E - H^{(S)}(x_S))) \simeq \exp\left\{\frac{1}{k_b} S^R(E - H^S)\right\}$$
(13)

## Concepts of statistical mechanics: canonical ensemble

$$\exp\left\{\frac{1}{k_b}S^R(E-H^S)\right\} \simeq \exp\left[\frac{1}{k_b}S^R(E)\right] \exp\left[-\frac{1}{k_b}\frac{\partial S^{(R)}}{\partial E}|_E H^{(S)}(x_S)\right]$$
(14)

$$\langle A(x) \rangle = \frac{1}{Z} \int dx_s A(x_s) exp \left[ -\frac{H^{(S)}(x_s)}{k_b T} \right]$$
 (15)

$$Z = \int dx_s \exp\left[-\frac{H^S(x_s)}{k_b T}\right]$$
 (16)

$$\langle A(x) \rangle = \frac{1}{Z} \int dE \int dx \delta(H(x) - E) A(x) exp(-\frac{E}{k_b T})$$
 (17)

$$\langle A(x)\rangle = \frac{1}{Z} \int dE' a^*(E') exp \left[ -\frac{E' - TS(E')}{k_b T} \right]$$
 (18)

# Concepts of statistical mechanics: canonical ensemble[5]

$$Z \simeq \exp\left[-\frac{E^* - TS(E^*)}{k_b T}\right] = \exp\left(-\frac{F}{k_b T}\right)$$
 (19)

$$\frac{\partial \ln Z(\beta)}{\partial \beta} = -\frac{1}{Z} \int dx H(x) \exp\left[-\frac{H}{k_b T}\right] = -\langle H(x) \rangle = -E$$
 (20)

$$\frac{\partial^2 \ln Z(\beta)}{\partial \beta^2} = \left\langle H(x)^2 \right\rangle - \left\langle H(x) \right\rangle^2 \tag{21}$$

$$\langle H(x)^2 \rangle - \langle H(x) \rangle^2 = -\frac{\partial E}{\partial (1/k_b T)} = k_b T^2 \frac{\partial E}{\partial T} = k_b T^2 C$$
 (22)

In this way a statistical quantity the variance has been connected to a thermodynamic quantity: the temperature.