

Long Range Correlations in Driven Systems (I)

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Non-equilibrium systems

Systems with currents
(driven by electric field, T gradients etc.)

In many cases these systems reach a steady state
(but non-equilibrium steady state).

What is the nature of these steady states?

In general, the probability distribution to be in a microstate C evolves by the master equation

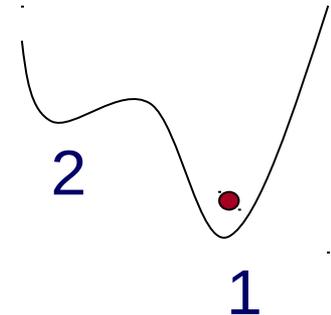
$$\frac{\partial P(C)}{\partial t} = \sum_{C'} W(C' \rightarrow C)P(C') - \sum_C W(C \rightarrow C')P(C)$$

in steady state $\frac{\partial P(C)}{\partial t} = 0$

equilibrium (non-driven) reach a steady state which satisfies the detailed balance condition for every microstates C and C' :

$$W(C' \rightarrow C)P(C') = W(C \rightarrow C')P(C)$$

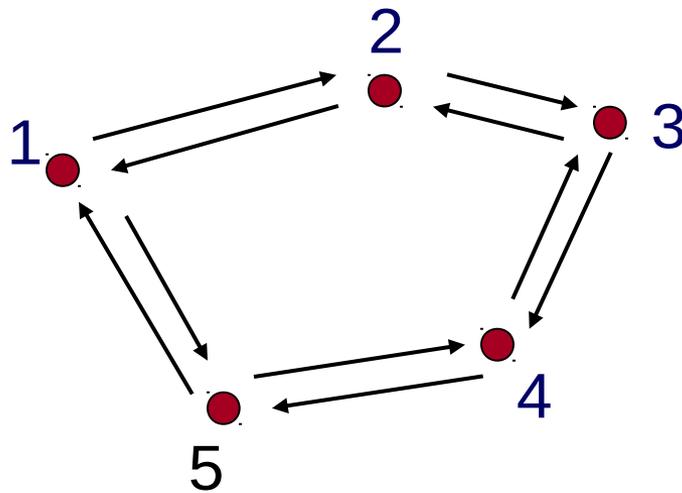
(no net probability current between two states)



Given transition rates $W(C \rightarrow C')$

a necessary and sufficient condition for detailed balance: for any set of microstates C_1, \dots, C_k

$$W(1 \rightarrow 2)W(2 \rightarrow 3)\dots W(k \rightarrow 1) = W(k \rightarrow k-1)\dots W(2 \rightarrow 1)W(1 \rightarrow k)$$

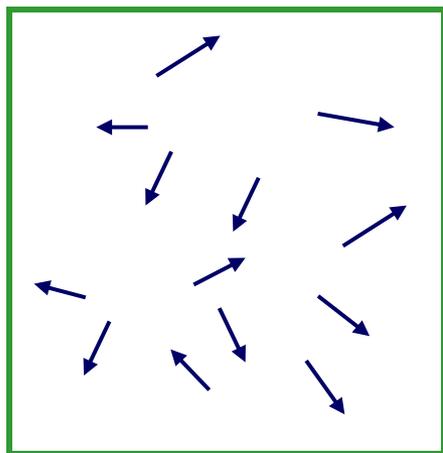


Equilibrium states (detailed balance, no currents)

- Collective phenomena
- Phase transitions (first or second order)
- Long range order
- Spontaneous symmetry breaking
- Phase separation
- Critical behavior
- Fluctuations in the equilibrium state (spatial or temporal)
- Relaxation processes to equilibrium states
- Effect of disorder
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Rules governing **equilibrium** collective phenomena

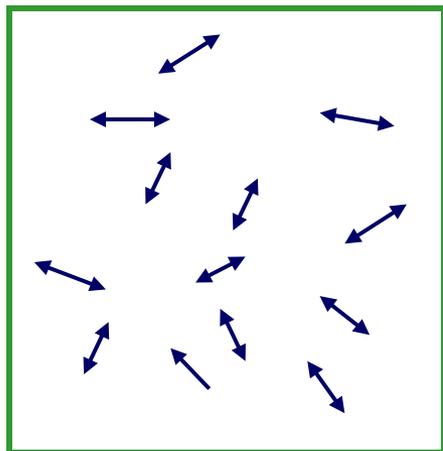
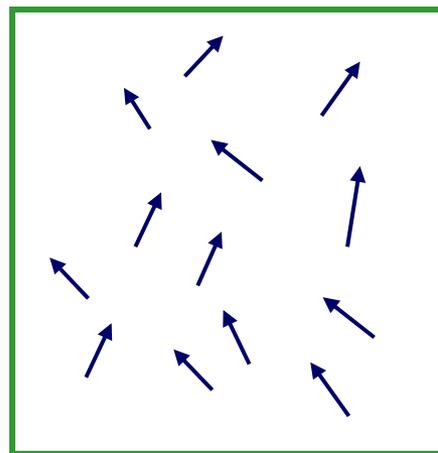
- Landau's symmetry rules for the order of the transition (ferromagnets - second order ; nematic transition - first order)
- Symmetry classification into universality classes
- No long range order in low dimensional systems
- Renormalization group criteria for the order of the transition
- Gibbs phase rule (dimension of the coexistence manifold)
 $D=2+c-n$
- 180° rule
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magnetic transition



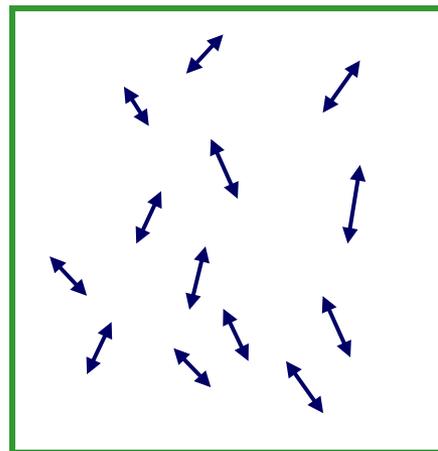
2nd order



nematic liquid crystals

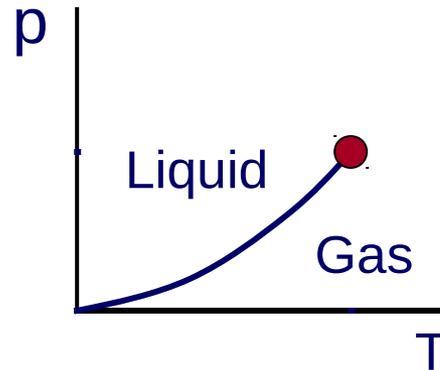


1st order



Gibbs phase rule

(dimension of manifold of n coexisting phases in c - components fluid mixtures)



$$c=1$$
$$n=2$$

Gibbs phase rule $D=2+c-n$

(for the dimension of manifold of n coexisting phases
In fluid mixtures)

c - number of components in fluid mixtures

n - number of coexisting phases

D - dimension of the manifold of n coexisting phases

180° rule

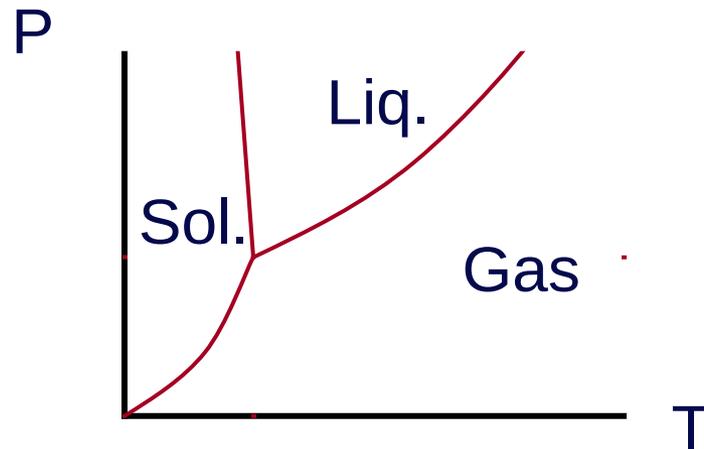
(for coexistence lines in phase diagrams)



yes



no



Do similar rules exist for **non-equilibrium** (driven) systems)?
(for which “free energy” cannot be defined)

In fact most of these rules **do not apply** in non-equilibrium systems.

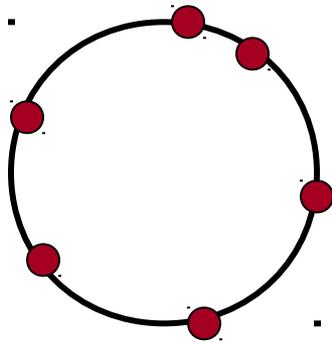
Phase separation in 1d

In thermal equilibrium:

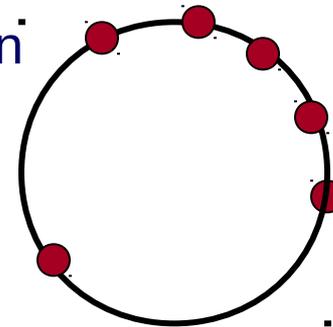
short range interactions
 $T > 0$



Density is macroscopically
homogeneous



No liquid-gas transition



Landau, Peierls 1930's: no phase separation, long range order, spontaneous symmetry breaking, phase transitions in 1d.

A simple physical argument for **no long-range order in 1d**

Ising model:
$$H = -J \sum_n s_n s_{n+1} \quad s = \pm 1, \quad J > 0$$

Ground state:

+++++

Consider the evolution of this state: since $T > 0$
a “wrong” droplet will be created in time

+++++ - - - +++++

Once created, the droplet may increase (or decrease)
without energy cost.

In one-dimension “wrong” droplets are not eliminated

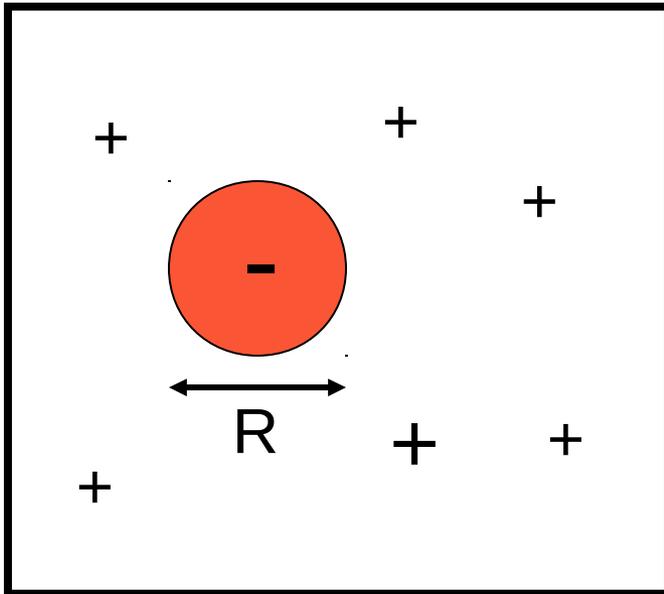
+++++ - - - ++++++ - - - ++++++ - - - ++++++

The energy of a droplet does not depend on its length (the energy cost of each droplet is $4J$).

The length of droplets will fluctuate in time, droplets will merge and long range order will be destroyed in time.

Robust argument: the only ingredients are $T > 0$ and **short range** interactions.

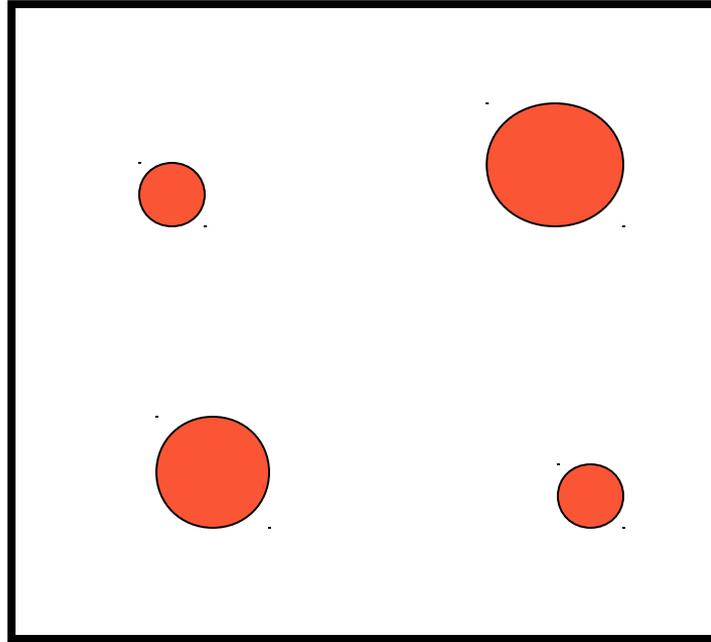
Maintaining long range order in higher dimensions:



Larger droplet cost more (surface) energy.

$$\frac{dR}{dt} \approx -\frac{\sigma}{R}$$

σ - surface tension



Wrong droplets are generated by fluctuations but are eliminated by surface tension.

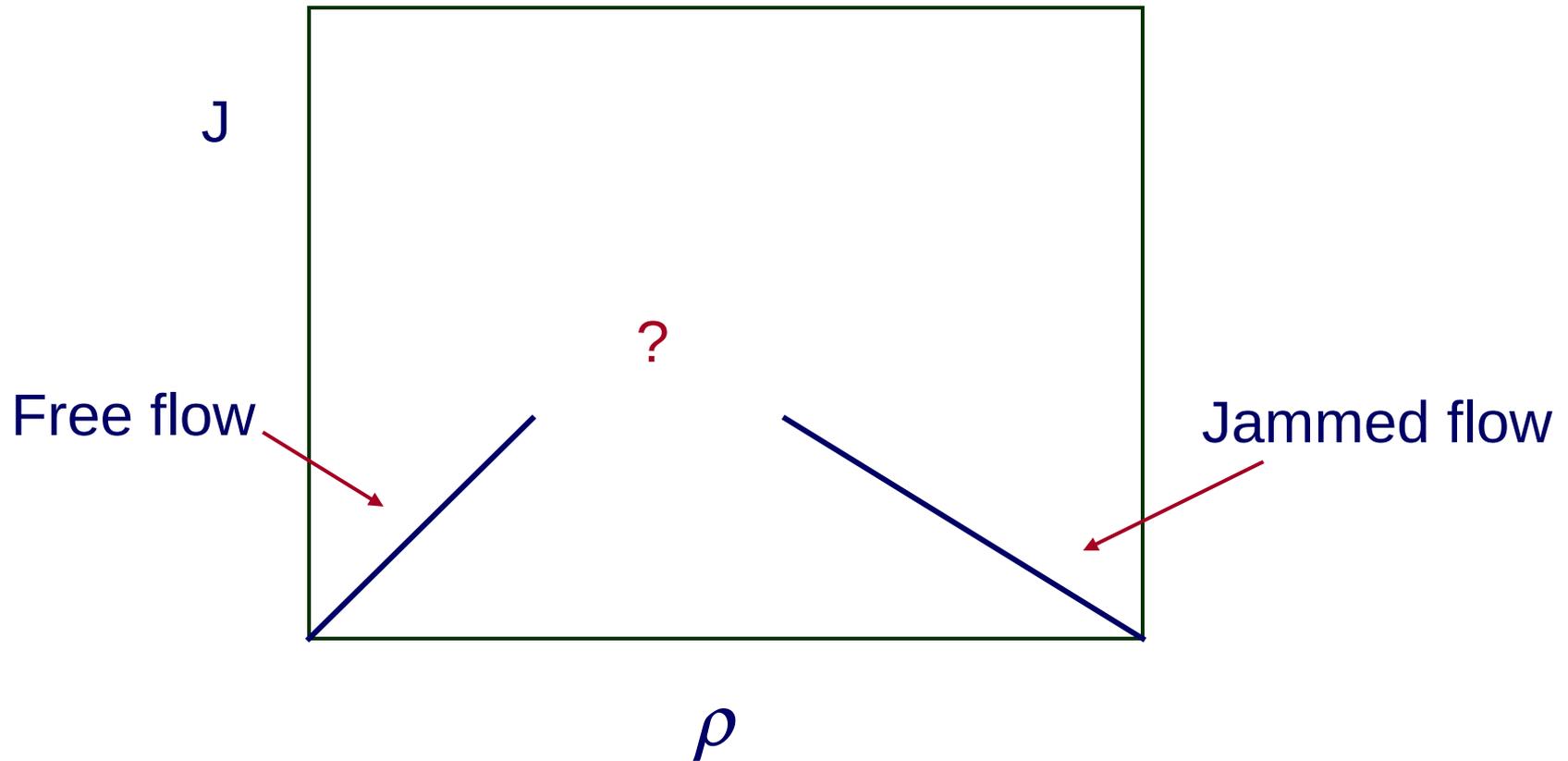
At sufficiently low T no large droplets are formed and long range order is maintained.

Can one have phase separation in 1d driven systems (?)

local, noisy dynamics
homogeneous, ring geometry
no detailed balance

A criterion for phase separation in such systems (?)

Traffic flow - fundamental diagram

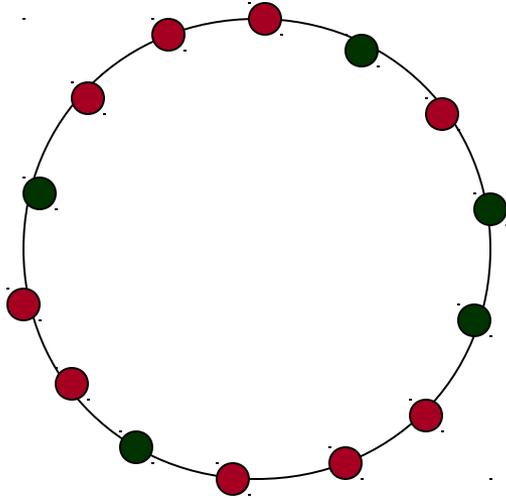


Is there a jamming phase transition?
or is it a broad crossover?

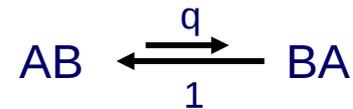
Main points

- Phase transitions do exist in one dimensional driven systems.
- In many traffic models studied in recent years Jamming is a crossover phenomenon. Usually it does not take place via a genuine phase transition.

Asymmetric Simple Exclusion Process (ASEP)



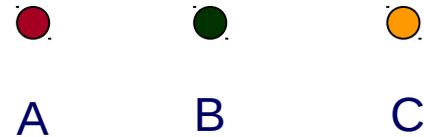
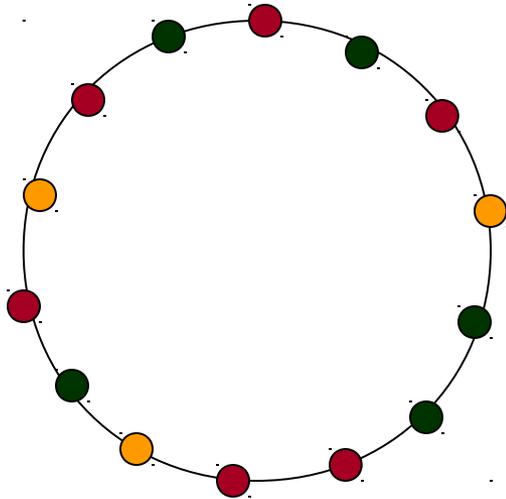
dynamics



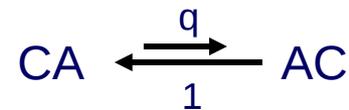
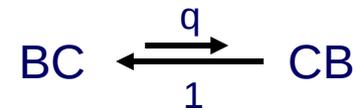
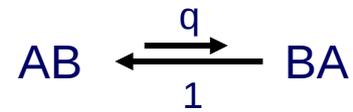
Steady State:

- ★ $q=1$ corresponds to an Ising model at $T = \infty$
- ★ All microscopic states are equally probable.
- ★ Density is macroscopically homogeneous.
No liquid-gas transition (for any density and q).

ABC Model



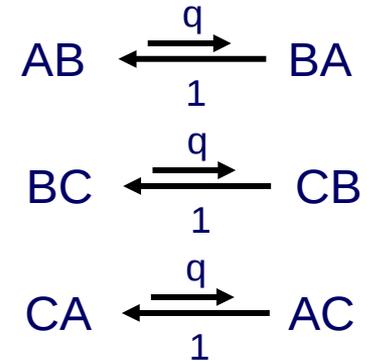
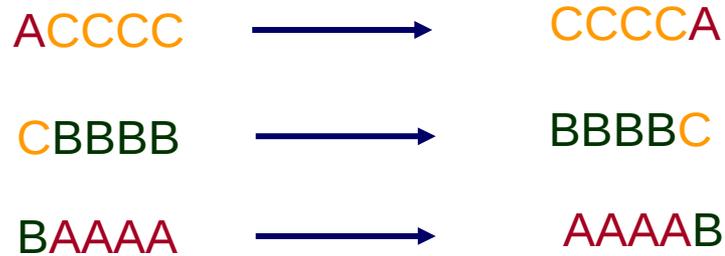
dynamics



Evans, Kafri, Koduvely, Mukamel PRL 80, 425 (1998)

A model with similar features was discussed by Lahiri, Ramaswamy PRL 79, 1150 (1997)

Simple argument:



...AACBBBCCAAACBBBCCC...

...AABBBCCCAABBBCCCC...

...AAAAABBBBBCCCCCAA...



fast rearrangement



slow coarsening

- logarithmically slow coarsening

...AAAAABBBBBBCCCCC...AA...

$$t \propto q^{-l} \quad l \propto \ln t$$

- needs $n > 2$ species to have phase separation
- Phase separation takes place for any q (except $q=1$)
- Phase separation takes place for any density N_A, N_B, N_C
- strong phase separation: no fluctuation in the bulk; only at the boundaries.

...AAAAAAAAAABBBBBBBBBBBBBBCCCCCCCCCCCC...

Special case $N_A = N_B = N_C$

The argument presented before is general, independent of densities.

For the equal densities case the model has **detailed balance** for **arbitrary q** .

We will demonstrate that for any microscopic configuration $\{X\}$

One can define “energy” $E(\{X\})$ such that the steady state

Distribution is

$$P(\{X\}) \propto q^{E(\{X\})}$$

AAAAAABBBBBBCCCCC

E=0



With this weight one has:

$$W(AB \rightarrow BA)P(\dots AB \dots) = W(BA \rightarrow AB)P(\dots BA \dots)$$

$=q$ $=1$

$$P(\dots BA \dots) / P(\dots AB \dots) = q$$

This definition of “energy” is possible only for $N_A = N_B = N_C$

AAAAABBBBBBCCCCC \longrightarrow AAAABBBBBBCCCCCA

E \longrightarrow E + N_B - N_C

$$N_B = N_C$$

Thus such “energy” can be defined only for $N_A = N_B = N_C$

AABB BBCC CAAAAA BBBB CCCC

The rates with which an A particle makes a full circle clockwise
And counterclockwise are equal

$$q^{N_B} = q^{N_C}$$

Hence no currents for any N.

For $N_B \neq N_C$ the current of A particles satisfies $J_A \propto q^{N_B} - q^{N_C}$

The current is **non-vanishing** for finite N. It vanishes only in the
limit $N \rightarrow \infty$. Thus no detailed balance in this case.

The model exhibits strong phase separation

...AAAAAABBBABBBBBBCCCCCCCCAA...

The probability of a particle to be at a distance l on the wrong side of the boundary is q^l

The width of the boundary layer is $-1/\ln q$

$$N_A = N_B = N_C$$

The “energy” E may be written as

$$P(\{x\}) = q^{E(\{x\})}$$

$$E(\{x\}) = \sum_{i=1}^N \sum_{k=1}^{N-1} \left(1 - \frac{k}{N}\right) (C_i B_{i+k} + A_i C_{i+k} + B_i A_{i+k})$$

summation over $(i+k)$ modulo N

Local dynamics



• long range

Partition sum

Excitations near a single interface: **AAAAAAAABBBBBB**

$$Z_1(q) = \sum p(n)q^n$$

$P(n)$ = degeneracy of the excitation with energy n

$$P(0)=1$$

$$P(1)=1$$

$$P(2)=2 \text{ (2, 1+1)}$$

$$P(3)=3 \text{ (3, 2+1, 1+1+1)}$$

$$P(4)=5 \text{ (4, 3+1, 2+2, 2+1+1, 1+1+1+1)}$$

$P(n)$ = no. of partitions of an integer n

$$Z_1(q) = \frac{1}{(1-q)(1-q^2)\dots}$$

Partition sum: $Z(q) = N \left[\frac{1}{(1-q)(1-q^2)\dots} \right]^3$

Correlation function: $\langle A_1 A_r \rangle \approx 1/3$

with $\langle A_1 \rangle \langle A_r \rangle = 1/9$

for $-1/\ln q < r < N/3$

Summary of ABC model

- The model exhibits phase separation for any $q \neq 1$
- Needs $n > 2$ species for phase separation.
- Strong phase separation (probability to find a particle in the bulk of the “wrong” is exponentially small.
- Phase separation is a result of effective **long range Interactions** generated by the **local** dynamics.
- Logarithmically slow coarsening process.