

# Multi-dimensional quantum paths to optimization problems

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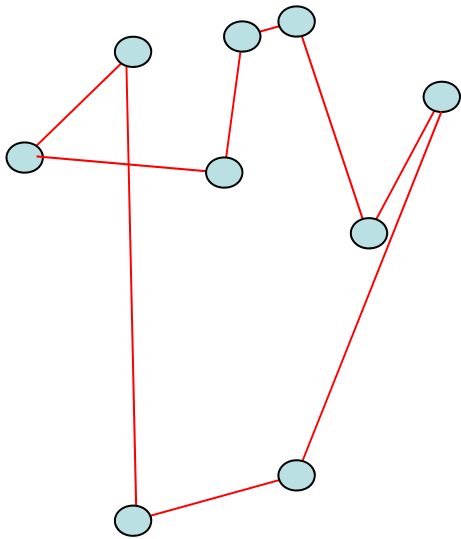
*Tokyo Institute of Technology*



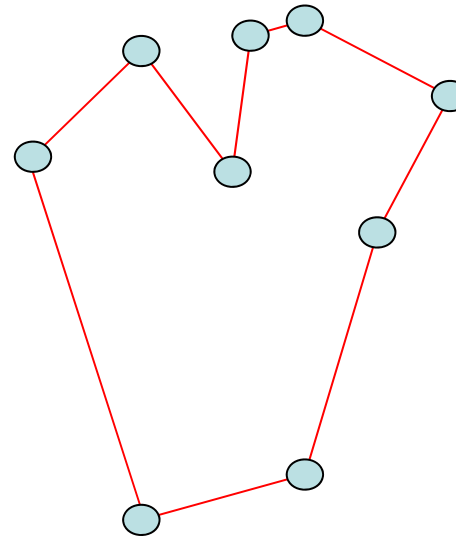
# Quantum annealing

# Combinatorial optimization

- Travelling Salesman Problem (TSP)



Configuration 1



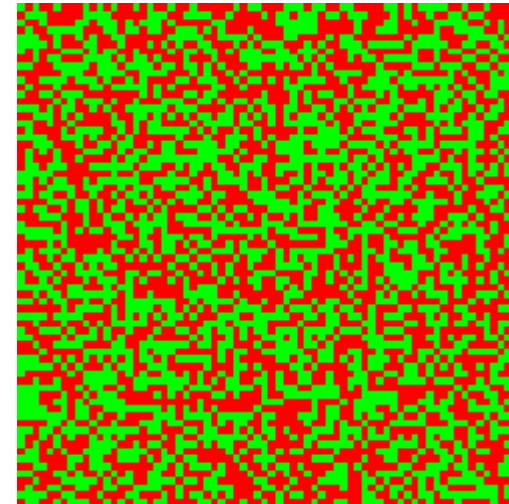
Configuration 2

Minimize the cost function (=tour length)

# Problem in physics

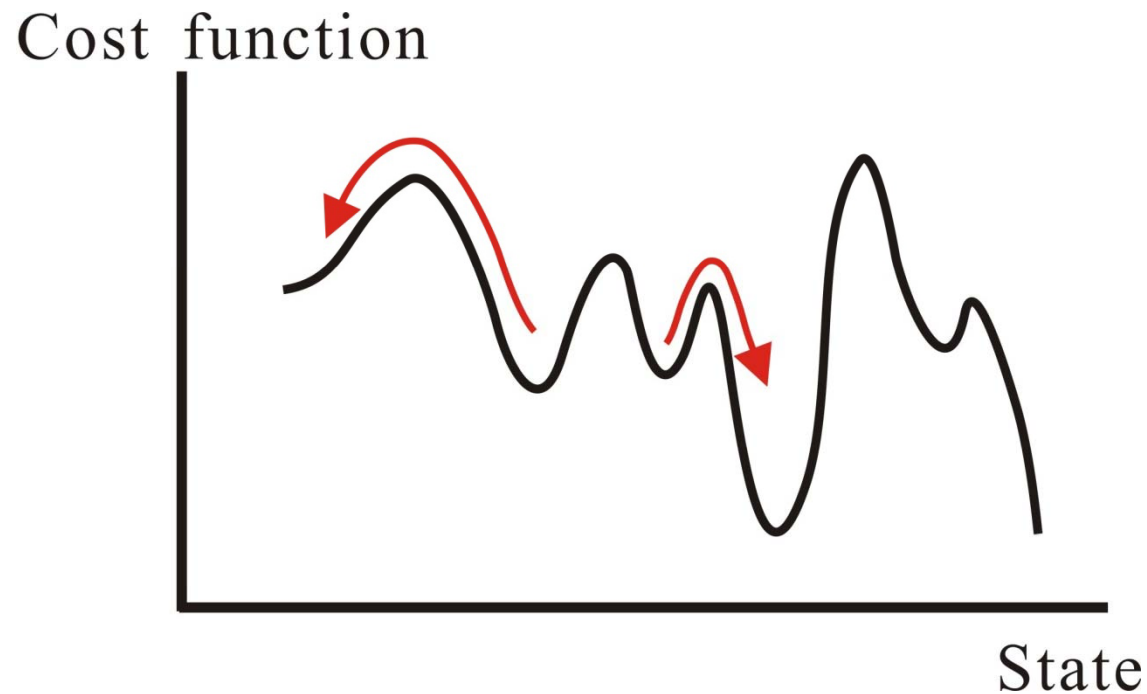
Ground state of Ising SG

$$H = -\sum J_{ij} S_i S_j$$



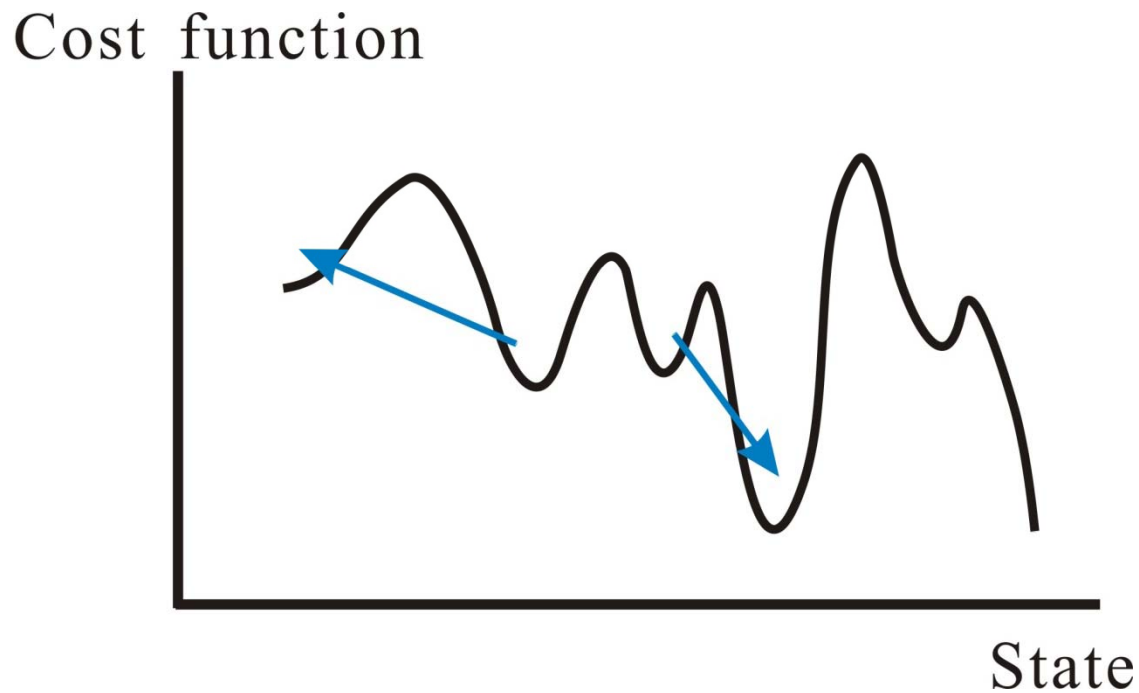
# Simulated Annealing (SA)

- *Generic, approximate* algorithm
- Phase-space search by **thermal** fluctuations



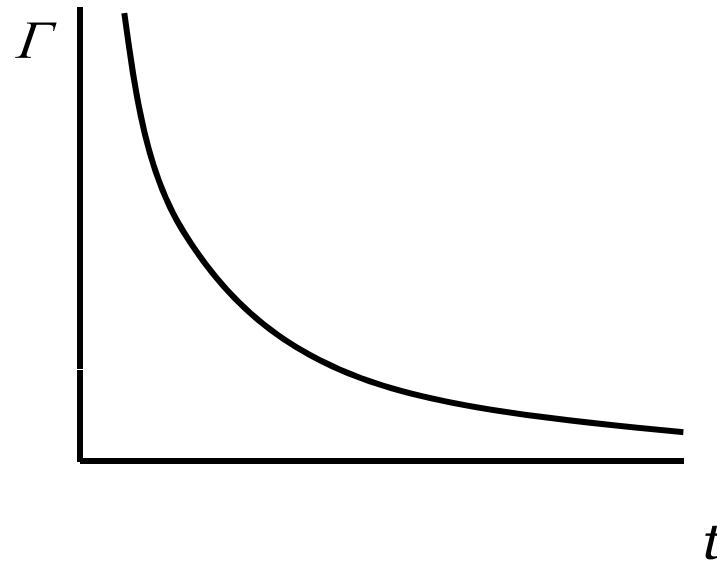
# Quantum Annealing (QA)

- *Generic, approximate* algorithm
- Phase-space search by **quantum** fluctuations



# Implementation

$$H(t) = H_{\text{classical}} + H_{\text{quantum}} = -\sum J_{ij} \sigma_i^z \sigma_j^z - \Gamma(t) \sum \sigma_i^x$$



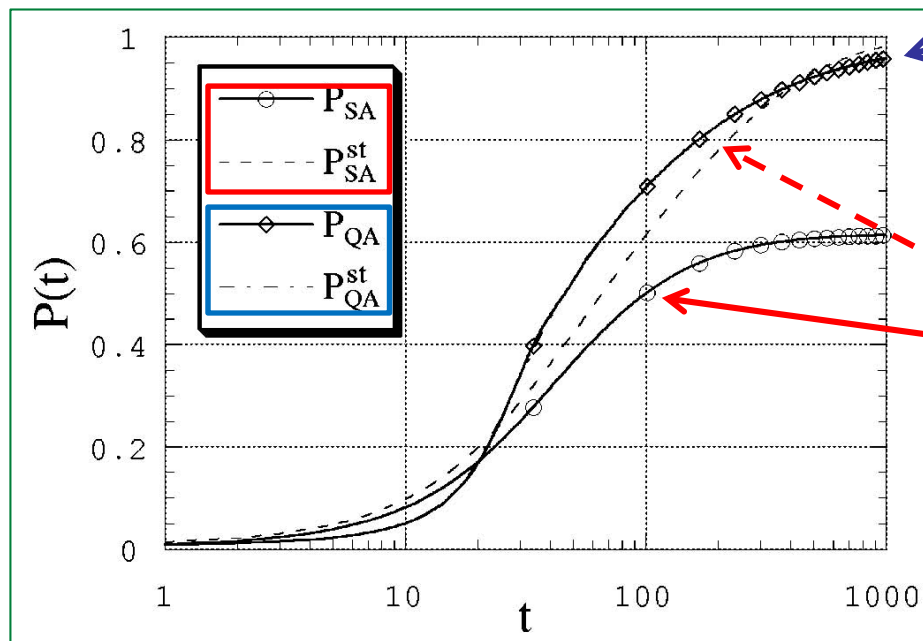
# Numerical evidence

$$H = -\sum J_{ij} \sigma_i^z \sigma_j^z - \Gamma(t) \sum_i \sigma_i^x$$



# Master eqn vs. Schrödinger eqn

Spin glass (SK model) with 8 spins



$$\Gamma(t) = \frac{3}{\sqrt{t}}$$

Schrödinger

$$T(t) = \frac{3}{\sqrt{t}}$$

Master /Equilibrium

# Monte Carlo for TSP (1002 cities)

$$H(t) = \frac{t}{\tau} H_{\text{classical}} + \left(1 - \frac{t}{\tau}\right) H_{\text{quantum}}$$

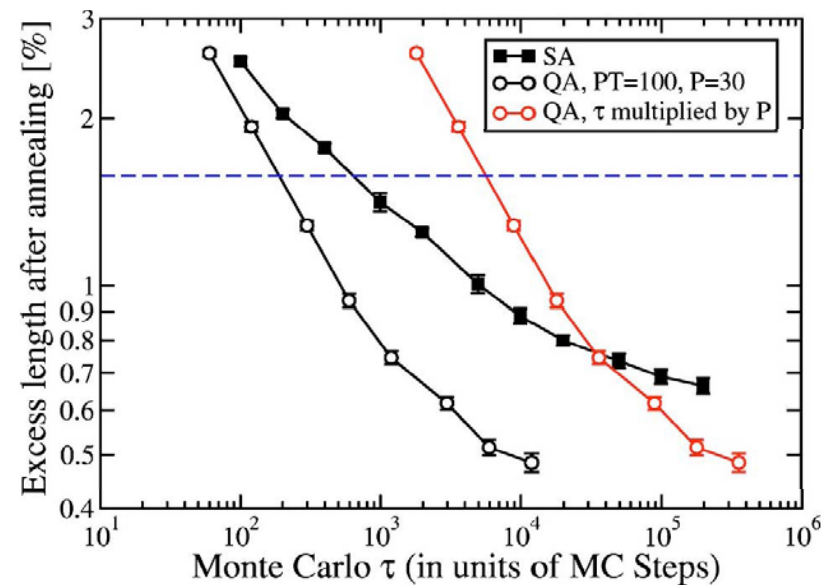
$$H(0) = H_{\text{quantum}} \Rightarrow H(\tau) = H_{\text{classical}}$$

cf: classical SA

$$T(0) = \text{large} \Rightarrow T(\tau) = 0$$

Residual energy:  $H(\tau) - E_{\text{true}}$

*Martonak, Santoro & Tosatti (2004)*



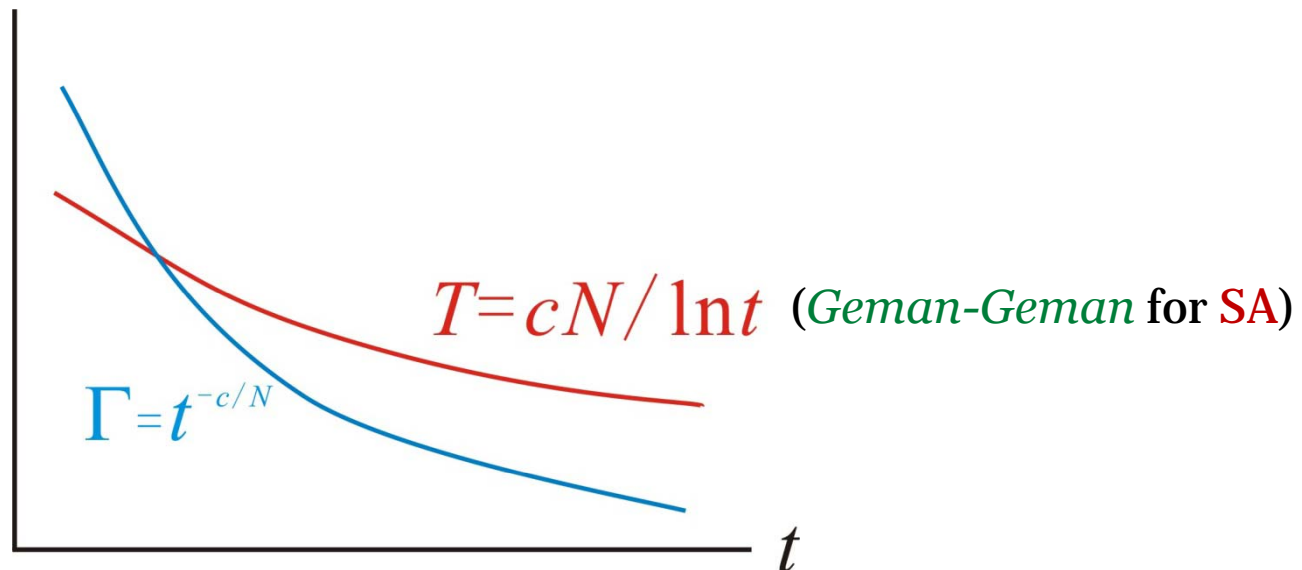
# Theoretical background

$$H = -\sum J_{ij} \sigma_i^z \sigma_j^z - \Gamma \sum_i \sigma_i^x$$

# Convergence theorem

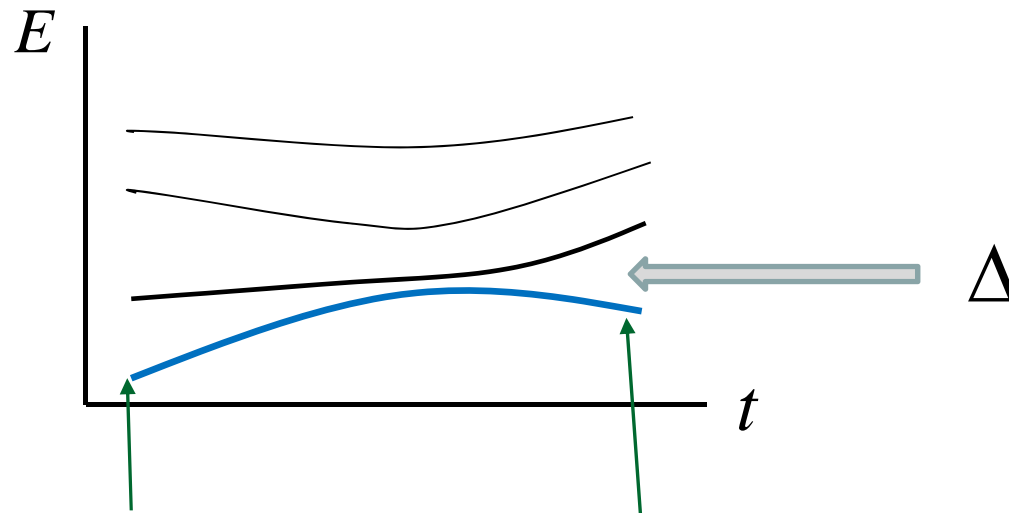
$$H = H_{\text{classical}} + H_{\text{quantum}} = -\sum J_{ij} \sigma_i^z \sigma_j^z - \Gamma(t) \sum \sigma_i^x$$

Convergence condition  $\Gamma(t) = t^{-c/N}$  *Morita & Nishimori*



*p*-spin ferromagnet  
- Recent results -

# Adiabatic evolution



Trivial initial state

Non-trivial final state

$$H(t) = -\left(1 - \frac{t}{\tau}\right) \sum \sigma_i^x - \frac{t}{\tau} \sum J_{ij} \sigma_i^z \sigma_j^z$$

# Computational complexity

## Finite-size analysis

Adiabatic theorem

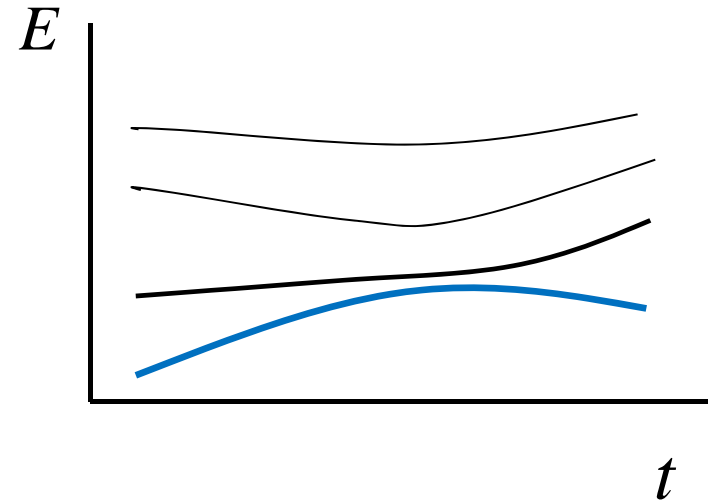
$$\tau \propto \Delta^{-2}$$

Gap scaling

$$\Delta \propto \begin{cases} e^{-aN} \\ N^{-b} \end{cases}$$

Complexity

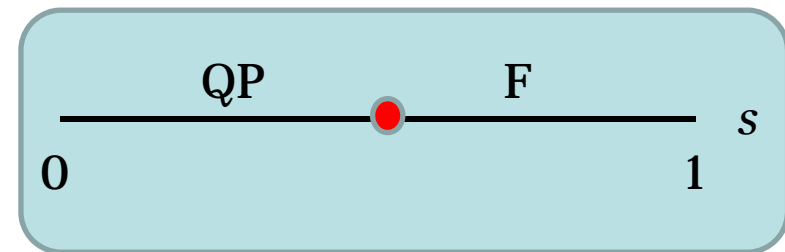
$$\tau \propto \begin{cases} e^{2aN} & \text{(hard)} \\ N^{2b} & \text{(easy)} \end{cases}$$



# $p$ -spin ferromagnet

$$H(s) = -sN \left( \frac{1}{N} \sum_{i=1}^N \sigma_i^z \right)^p - (1-s) \sum_{i=1}^N \sigma_i^x \quad (s = t / \tau)$$

- 1<sup>st</sup> order transition at finite  $s$



- Exponentially small energy gap.

$$\Delta \propto e^{-aN}$$

- Exponentially large time for adiabatic computation.

$$\tau \propto e^{bN}$$

Jörg *et al* (2010) “The problem that quantum annealing **cannot** solve”

Seki and Nishimori (2012) “The problem that quantum annealing **CAN** solve”



# An additional quantum term

$$H(\lambda, s) = s \left( H_0 \right) + (1-s) H_{\text{TF}}$$

$$H_0 = -N \left( \frac{1}{N} \sum_{i=1}^N \sigma_i^z \right)^p, \quad H_{\text{TF}} = -\sum_{i=1}^N \sigma_i^x$$

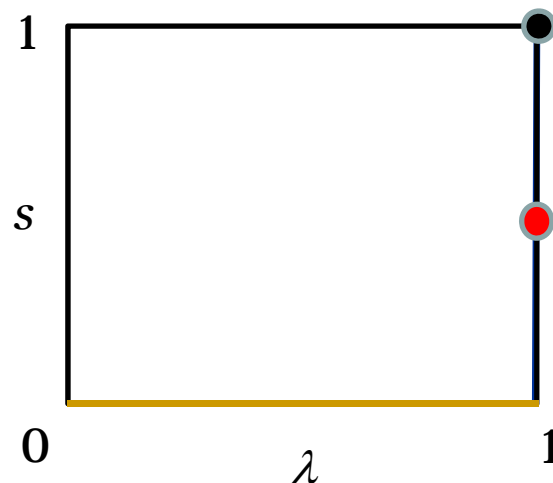
$$H(\lambda, s) = s \left( \lambda H_0 + (1-\lambda) H_{\text{AFF}} \right) + (1-s) H_{\text{TF}}$$

$$H_{\text{AFF}} = N \left( \frac{1}{N} \sum_{i=1}^N \sigma_i^x \right)^2$$

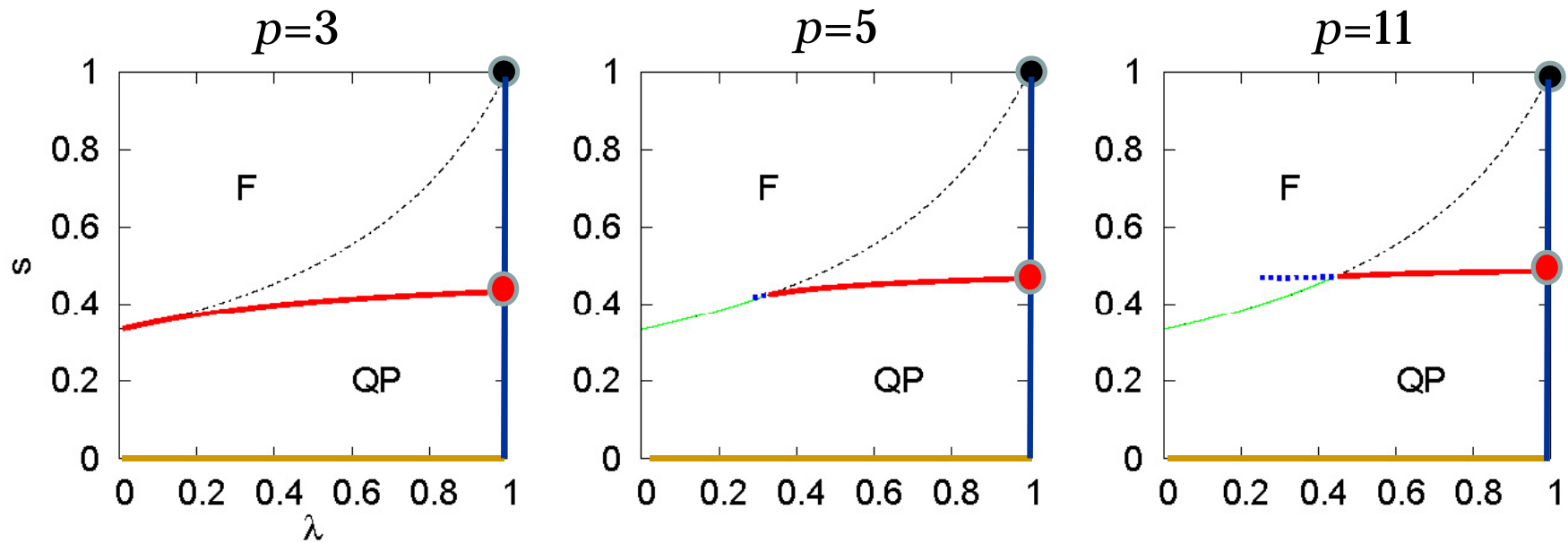
Conventional case:  $\lambda=1$

Start:  $s=0, \lambda=\text{any}$

Goal:  $s=1, \lambda=1$  ●



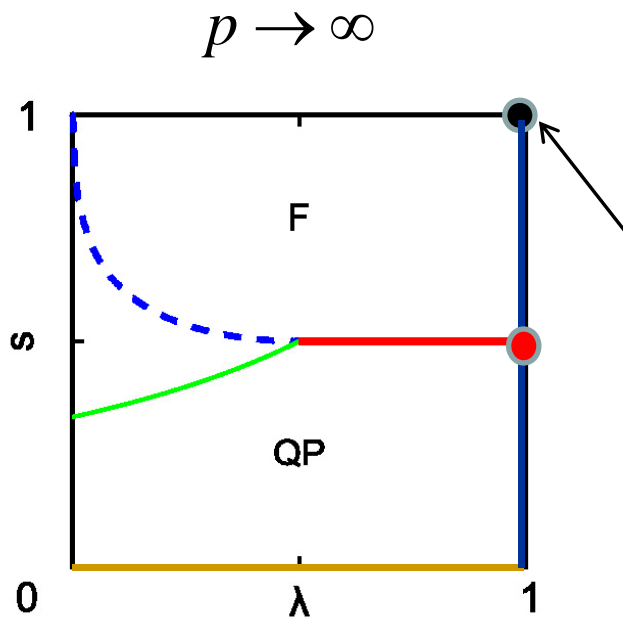
# Result



— 1st     
 - - - 1st     
 — 2nd

$$H(\lambda, s) = s(\lambda H_0 + (1 - \lambda)H_{\text{AFF}}) + (1 - s)H_{\text{TF}}$$

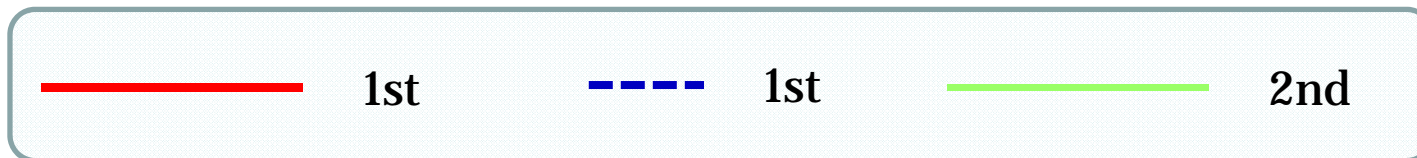
# Result (cont'd)



Grover problem

$$\frac{H_0}{N} = -\left(\frac{1}{N} \sum_{i=1}^N \sigma_i^z\right)^p = -1 \quad (\sigma_i^z = 1, \forall i)$$

$$\frac{H_0}{N} = 0 \quad \text{otherwise}$$



**Is**  $H_{\text{AFF}} = N \left( \frac{1}{N} \sum_{i=1}^N \sigma_i^x \right)^2$  **essential?**

Seoane and Nishimori (2012)

$$H(\lambda, s) = s \left( \lambda H_0 + (1 - \lambda) H_{\text{AFF}}^{(k)} \right) + (1 - s) H_{\text{TF}}$$

$$H_{\text{AFF}}^{(k)} = N \left( \frac{1}{N} \sum_{i=1}^N \sigma_i^x \right)^k$$

**Even**  $k$  ( $=2, 4, \dots$ ) is OK .

**Odd**  $k$  ( $=3, 5, \dots$ ) is OK, *partially*.

$$\left| \langle \Psi_{\text{ferro}} | \Psi_{\text{AFF}}^{(k)} \rangle \right|^2 \propto 1 / \sqrt{N}$$

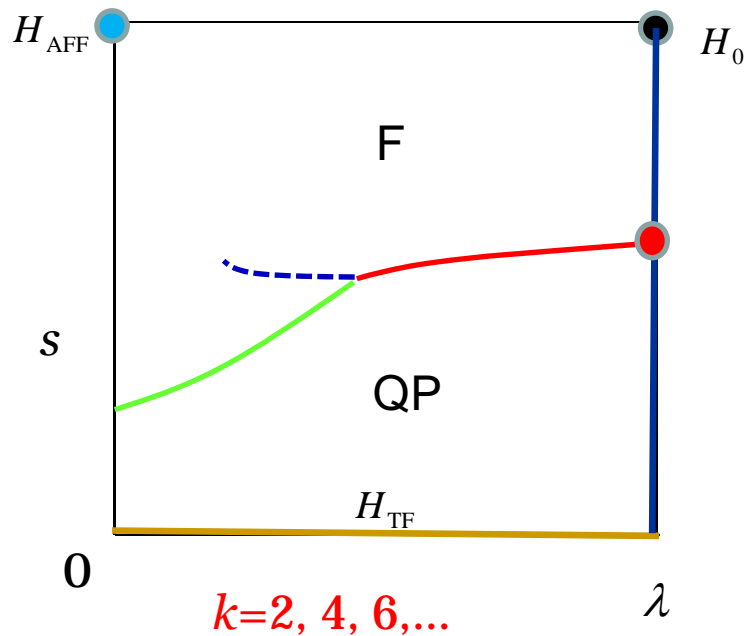
$$\left| \langle \Psi_{\text{ferro}} | \Psi_{\text{AFF}}^{(k)} \rangle \right|^2 \propto e^{-aN}$$

$k=2$ : Bapst and Semerjian (2012)

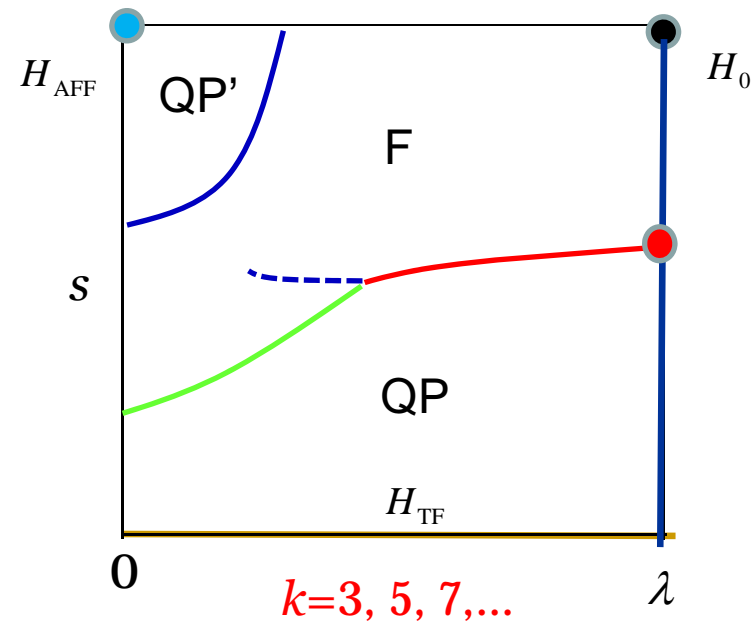
# Even $k$ and odd $k$

$$H(\lambda, s) = s(\lambda H_0 + (1 - \lambda)H_{\text{AFF}}^{(k)}) + (1 - s)H_{\text{TF}}$$

$$H_{\text{AFF}}^{(k)} = N \left( \frac{1}{N} \sum_{i=1}^N \sigma_i^x \right)^k$$



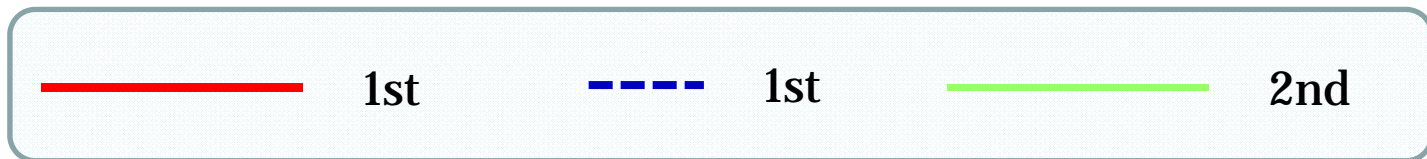
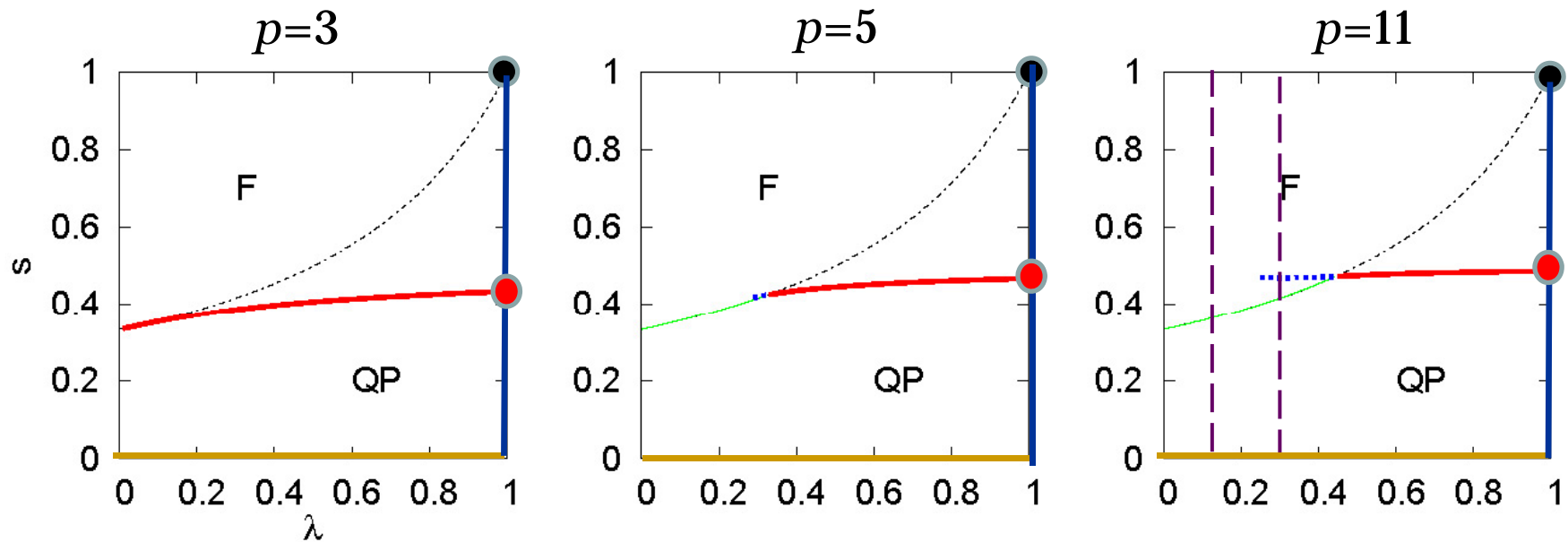
$$|\langle \Psi_{\text{ferro}} | \Psi_{\text{AFF}} \rangle|^2 \propto 1/\sqrt{N}$$



$$|\langle \Psi_{\text{ferro}} | \Psi_{\text{AFF}} \rangle|^2 \propto e^{-aN}$$

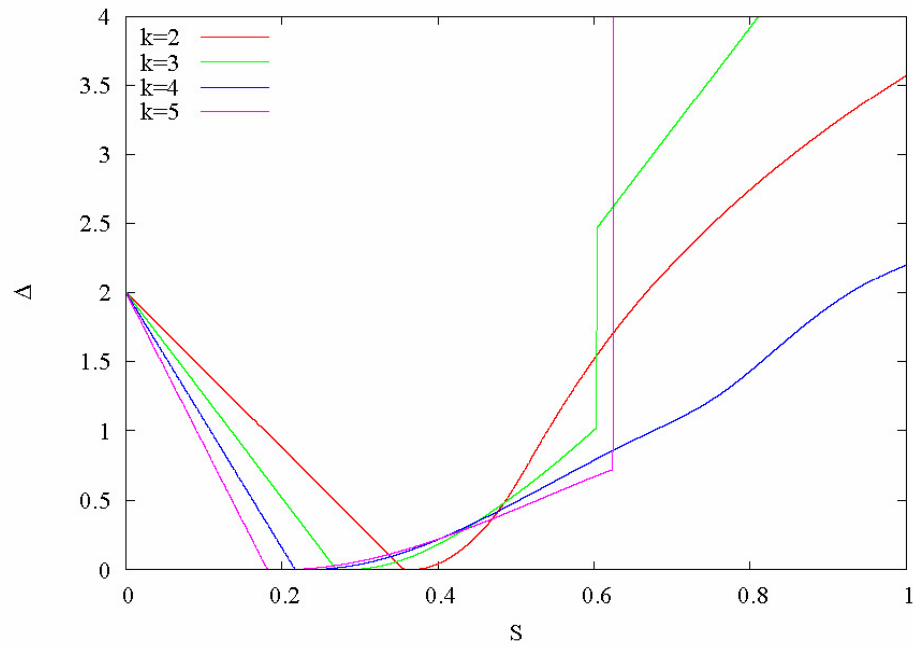
Bapst and Semerjian (2012)

# Energy gap

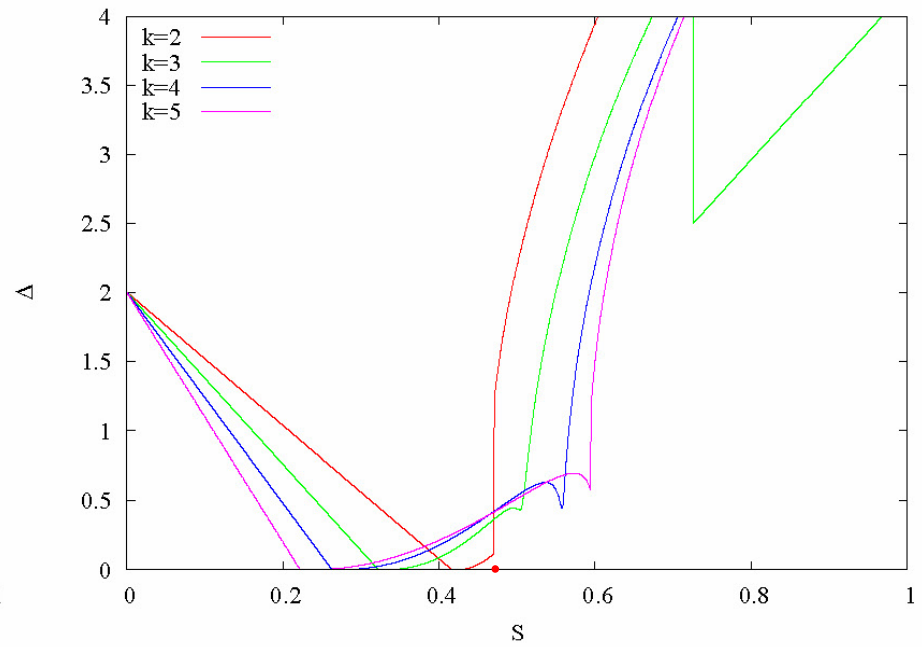


$$H(\lambda, s) = s(\lambda H_0 + (1 - \lambda)H_{\text{AFF}}) + (1 - s)H_{\text{TF}}$$

# Energy gap



$p=11, \lambda=0.1$



$p=11, \lambda=0.3$

# Method

1. Suzuki-Trotter + Static approximation
2. Classical approximation (+ quantum correction)

Botet-Julien (1983), Filiponne et al (2011)

$$\frac{1}{2} \sum_{i=1}^N \sigma_i^z = S^z \approx \frac{N}{2} \cos \theta, \quad \frac{1}{2} \sum_{i=1}^N \sigma_i^x = S^x \approx \frac{N}{2} \sin \theta \cos \varphi$$

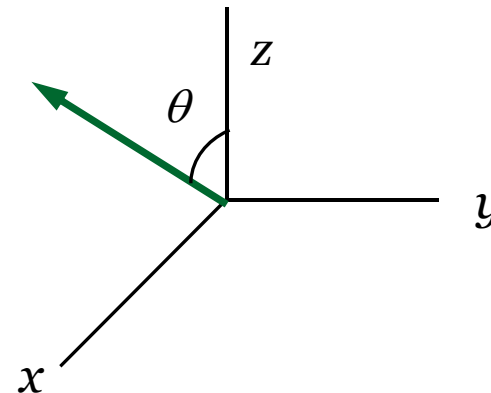
$$H_0 = -N \left( \frac{1}{N} \sum_{i=1}^N \sigma_i^z \right)^p, \quad H_{\text{TF}} = -\sum_{i=1}^N \sigma_i^x$$

$$H_{\text{AFF}} = N \left( \frac{1}{N} \sum_{i=1}^N \sigma_i^x \right)^2$$

$$\frac{H}{N} = -s\lambda(\sin \theta \cos \varphi)^p + s(1-\lambda) \cos^2 \theta - (1-s) \cos \theta$$

Minimize with respect to  $\theta$

$$\varphi = 0$$





# Method (Gap)

Classical approximation (+ **quantum correction**)

Holstein-Primakoff transformation + Bogoliubov transformation

$$S^z = \frac{N}{2} - a^+ a$$

$$S^+ = (N - a^+ a)^{1/2} a \approx Na, \quad S^- = a^+ (N - a^+ a)^{1/2} \approx Na^+$$

$$H = NE_g + c + \Delta(\theta)b^+ b$$

$\Delta(\theta)$ : Gap in the thermodynamic limit

# Summary

# Summary

1. QA works fine as a generic, approximate algorithm.
2. “Better” than SA. Numerical, analytical
3. Different quantum terms:  
Useful to avoid 1<sup>st</sup> order transitions.
4. Useful for **other, more complex, problems?**  
(under investigation)

# Collaborators

- Tadashi Kadowaki
- Helmut G. Katzgraber
- Yoshiki Matsuda
- Satoshi Morita
- Yuya Seki
- Beatriz Seoane