

21th EU/ME meeting x Quantum School

Emerging optimization methods: from metaheuristics to quantum approaches



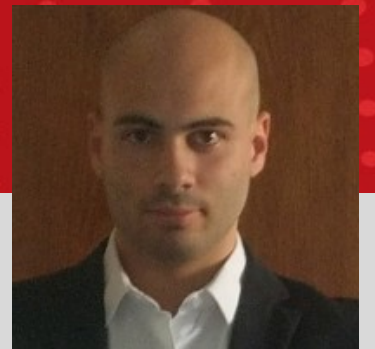
DE LA RECHERCHE À L'INDUSTRIE

# Discussions about High-Quality Embeddings on Quantum Annealers (WIP)

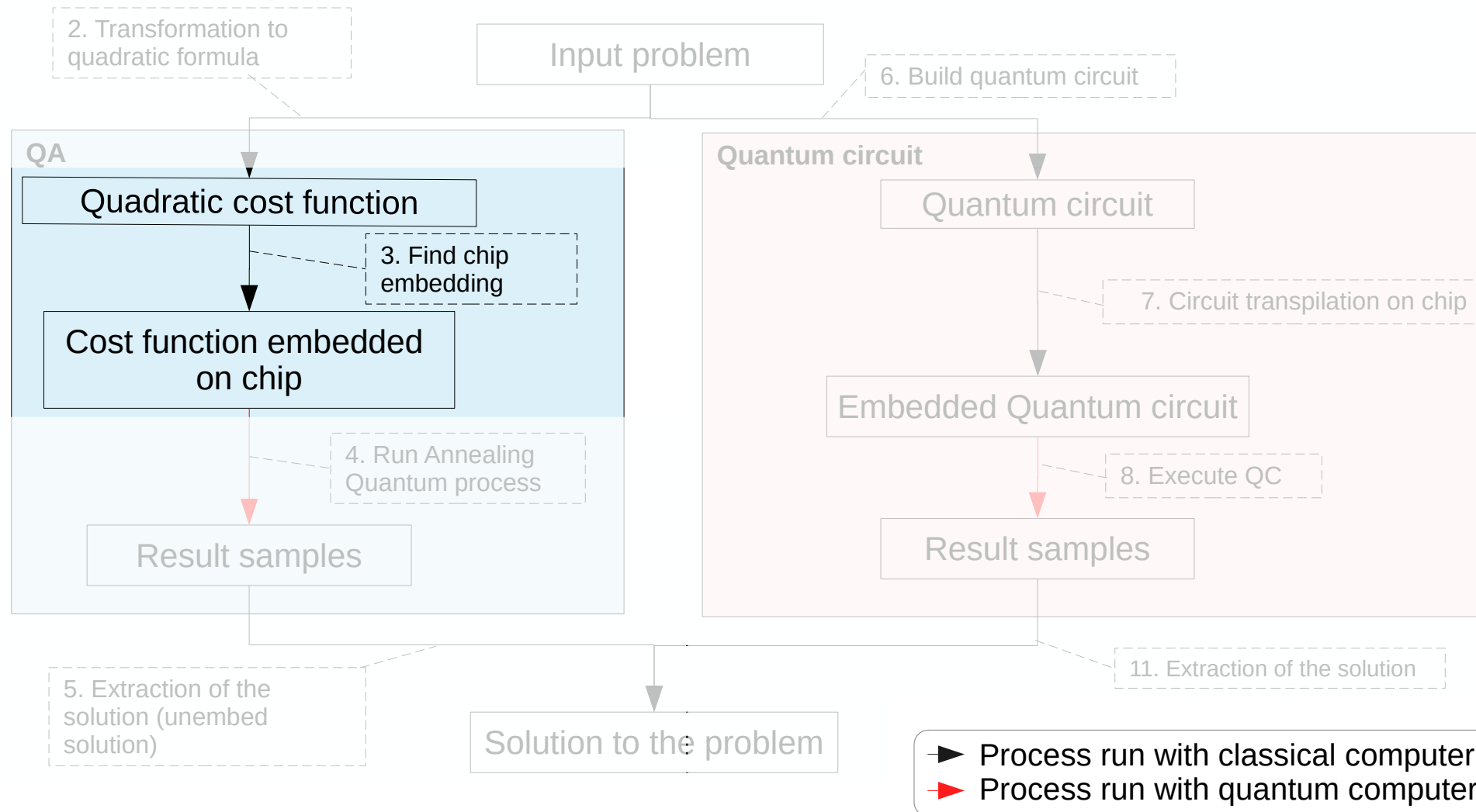
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- **Quantum computing is based on 2 main principles:**
  - Quantum superposition
  - Interference
- **What is limiting the Quantum Advantage:**
  - Quantum noise
  - Quantum chip topologies (require qubit mapping (QA) or swapping strategies (QC))
- **Overview:**
  - The Minor-embedding problem and its context
  - Existing work
  - Proposition of first experiments
  - Perspectives

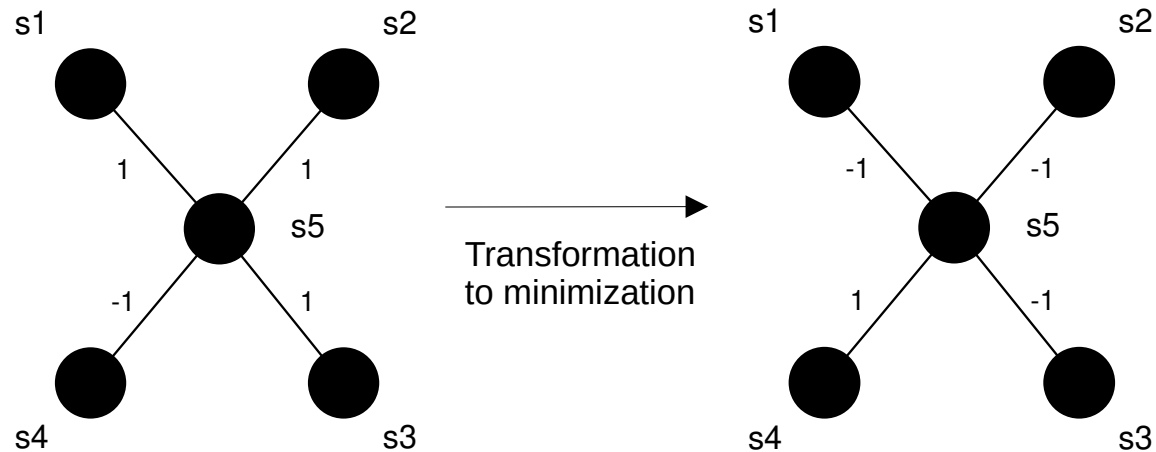


- Ising cost function: Minimization of a quadratic cost function

$$\text{Minimize } - \sum_{i=0}^n h_i s_i - \sum_{i < j} J_{ij} s_i s_j$$

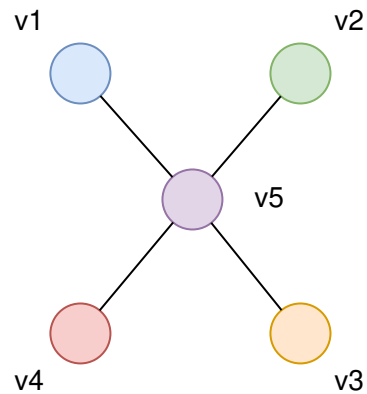
$$s_i, s_j \in \{-1, +1\} \text{ and } h_i, J_{ij} \in \mathbb{R}$$

- Ising problem formulation to solve Max-Cut problem

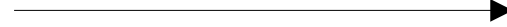


$$\text{Minimize } - (-s_1 s_5 - s_2 s_5 - s_3 s_5 + s_4 s_5)$$

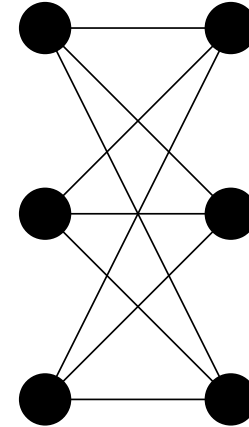
- Graph minor N. Robertson et al. [1]



$$G_s = (V_s, E_s)$$



$$\phi : V_s \rightarrow V_t \times V_t$$



$$G_t = (V_t, E_t)$$

- Rules for the graph minor:

1. Each vertex  $v \in V_s$  is mapped onto a connected subgraph  $\phi(v)$  of  $V_t$
2. Each connected subgraph must be vertex disjoint:  $\phi(v) \cap \phi(v') = \emptyset$  for  $v \neq v'$
3.  $\forall (u, v) \in E_s, \exists u' \in \phi(u), \exists v' \in \phi(v)$  such that  $(u', v') \in E_t$

- **Limitations & Complexity [1]**

- For fixed  $G_s$  finding the minor embedding in  $G_t$  has a polynomial complexity:

$$O(|V_s|^3)$$

**BUT**

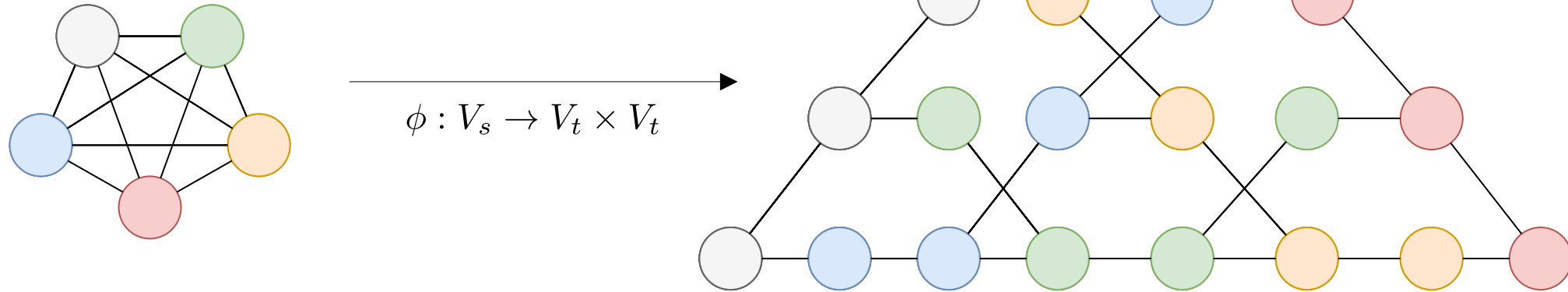
- The **algorithm runtime** is **exponential** in the **size of  $G_s$**

- **Heuristics are required to solve this problem efficiently**

- Mapping of cliques on regular graphs
  - Mapping of random source graphs on random target graphs
  - Other methods

- Mapping of cliques with near-optimal patterns [2, 3]

- Example with TRIAD pattern [2]

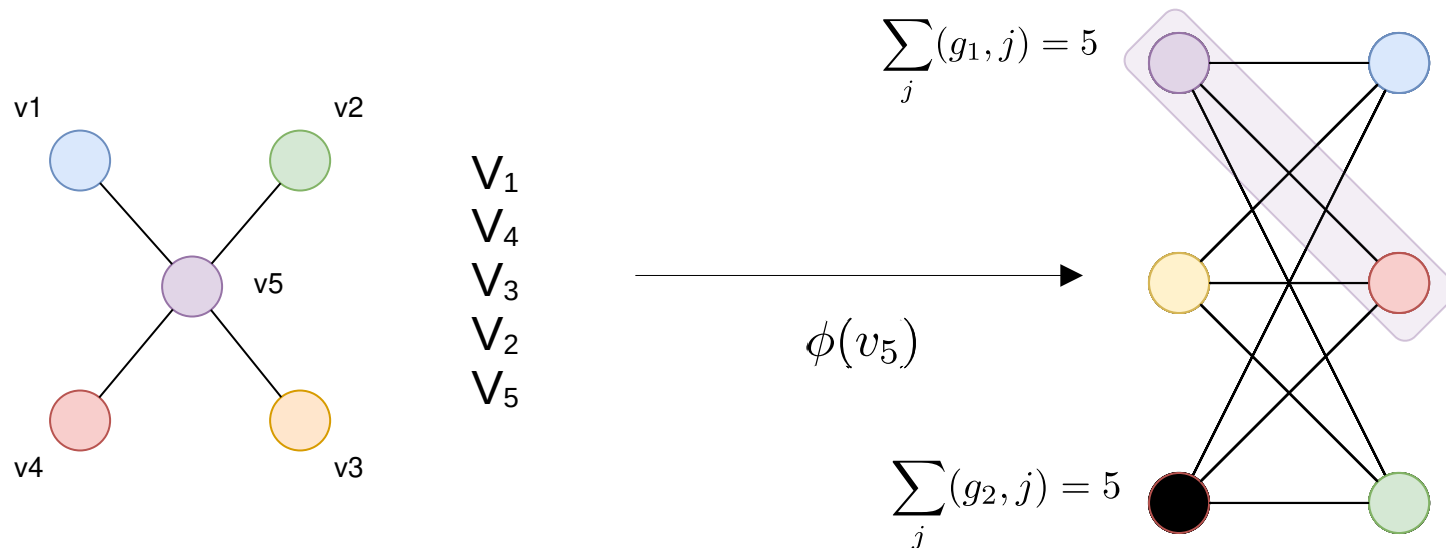


- Strength:
  - Near optimal encoding for dense graphs
- Weaknesses:
  - Does not always consider defective qubits
  - Not adapted for sparse graphs

- Algorithm of J. Cai et al [4, 5]

### Stage 1: Initialization

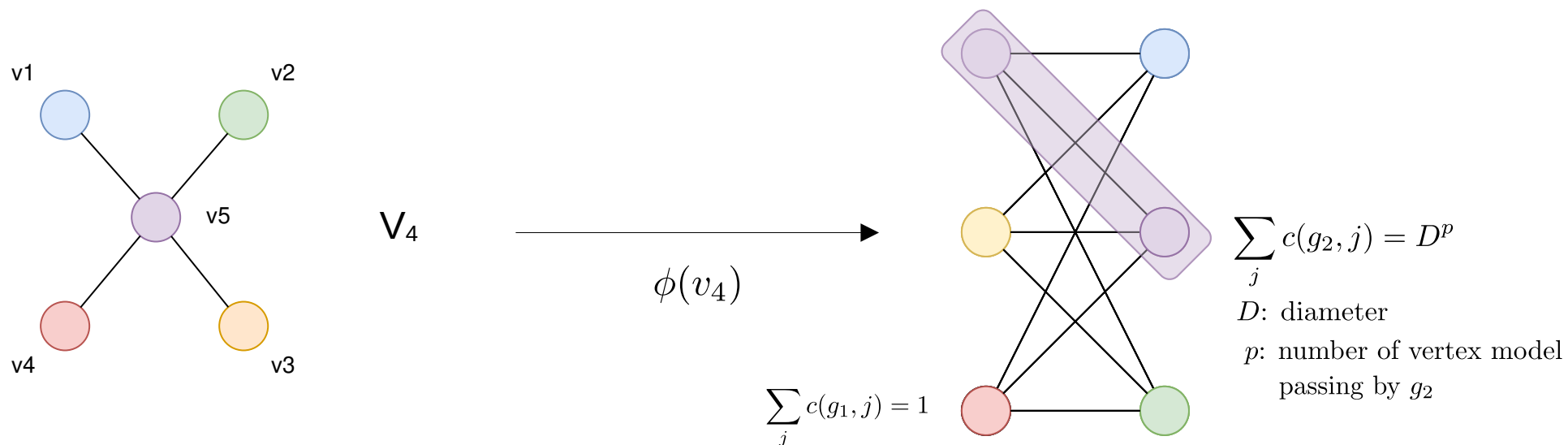
- Set the vertices in random order.
- For each vertex, find the set of vertices in the target graph that minimize the weighted shortest path distance to its neighbours (using Dijkstra's algorithm).





## Stage 2: Refinement

- Go through the vertices of  $G_s$ , remove its mapping and try to find another one, minimizing the weighted shortest path distance of the whole mapping to its neighbors.



- Strength:
  - Work with any kind of target graph
- Weaknesses:
  - Costly algorithm due to the computation of the shortest distance path at each optimization step:
  - Does not work well for dense graphs

$$O(n^3 \log n)$$

- **Other algorithms:**
  - Extensions to the algorithm of J. Cai et al. :
    - Layout-aware minor-embedding [6]
    - Clique-based minor-embedding [7]
  - SA-based approaches [8]
    - Starts with a clique near-optimal encoding (like TRIAD)
    - Run a guided simulated annealing to reduce the number of vertices.
- **Objectives of current methods:**
  - Minimization of the number of qubits
  - Minimize the maximum chain length

- Consequences of the qubit mapping

- Increases the number of qubits (qubit duplications)
- Changes the required precision of couplings and auto-couplings (automated rescaling of weights).

For D-Wave system Advantage6.1:

$$-1 \leq J_{ij} \leq 1 \quad -2 \leq h_i \leq 2$$

$$\text{Minimize} \quad - \sum_{i=0}^n (h_i + \delta h_i) s_i - \sum_{i < j} (J_{ij} + \delta J_{ij}) s_i s_j$$

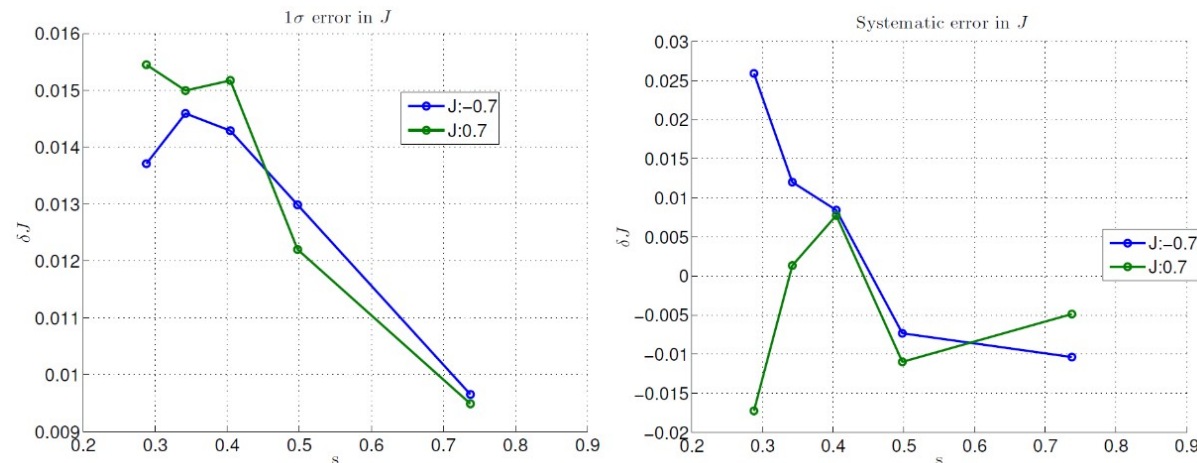


Figure taken from [9]. It Represents the standard deviation (left) and mean error rate (right) of  $J_{ij}$  coefficients with respect to arbitrary unit of annealing time.

- Potentially changes the spectral gap of the problem (i.e., increases or decreases the time to solve the problem)

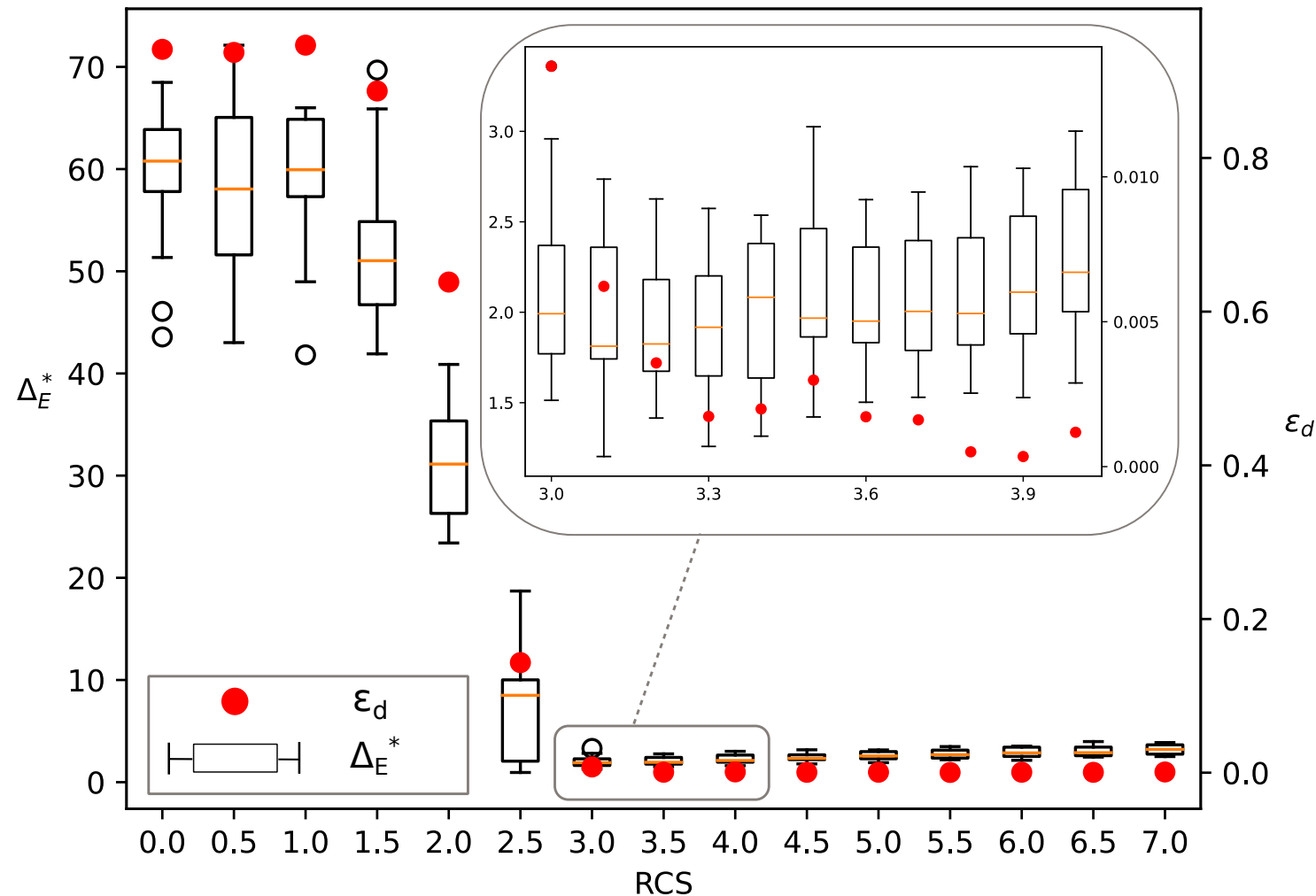
- Phase transition of the energy gap considering the chain strength [2]

Experimental determination of the optimal global chain strength (cs) and the Relative Chain Strength factor (RCS) [10]

$$cs = RCS \times \max(\{h_i\} \cup \{J_{ij}\})$$

$\epsilon_d$  : qubit duplication error rate after the  
( less is better)

$\Delta_E^*$  : energy gap between optimal and the  
solution obtained on D-Wave systems  
( less is better)



- **What is considered High-Quality embedding ?**
  - Minimization of the number of qubits
  - Minimize the maximum chain length
- **What is the best structure for logical qubits ?**
  - Error propagation on logical chains starts at the boundaries of the chain.
  - Errors on logical qubits require less coupling strength when they form a clique [11]
  - What about other topologies as cycles, trees, etc. ?
- **Is there a maximum chain length that shouldn't be reached ?**
  - Definition of bounds on the maximal chain length (requiring extra costs) could be computed considering QA precision and weights distribution.
- **Does chains distribution impact solution finding ?**
  - Study of chains sparsity versus concentration over the chip.
  - Perform experiments over different chain length distributions.

- **Many pieces of algorithms exist for minor-embedding graphs**
  - Major contribution made by J. Cai et al. In 2014.
  - Combinations of heuristics start to exist.
- **Tips and advices owned during our experiments**
  - Chain strength value is crucial so it has to be chosen carefully.
  - When mapping dense problems, use clique embedding methods.
  - Experiments should be done using as many qubits as possible to maximize the bias induced by imperfect couplers.
- **Perspectives**
  - Realization of the listed experiments.
  - Define metrics and bounds to quantify the quality of the embedding.
  - Depending on the result, design of a new heuristic.
  - Benchmarking the heuristic with state of the art methods.

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**Thank you !**

- **AQC, a continuous interpolation of 2 time independent Hamiltonians:**

$$H(t) = \left(1 - \frac{t}{T}\right) H_M + \frac{t}{T} H_C$$

- **The adiabatic theorem:**

“If a quantum state is initialized in the ground state of the Hamiltonian  $H_M$ , and that  $t$  varies slowly enough between 0 and  $T$ , the quantum state will state close to the ground state of  $H(t)$ ”

- **Quantum Annealing, a noisy version of AQC:**