

Quantum Annealers Chain Strengths: A Simple Heuristic to Set Them All

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I- Context & Motivations – Quantum Annealing (QA)

- D-Wave QA Hamiltonian

$$H(s) = A(s)H_M + B(s)H_P$$

$$H_M = \sum_{v \in V_s} \sigma_v^x$$

$$H_P = \sum_{v \in V_s} h_v \sigma_v^z + \sum_{(u,v) \in E_s} J_{uv} \sigma_u^z \sigma_v^z$$

- Eigen energies decomposition of the Hamiltonian as fixed s :

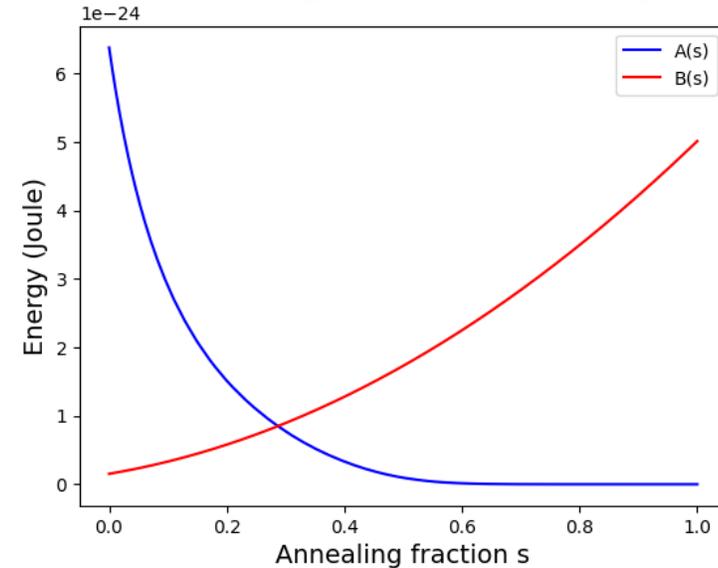
$$H_s |\psi_i(s)\rangle = E_i(s) |\psi_i(s)\rangle \quad \text{with } E_0(s) < E_1(s) < \dots < E_k(s)$$

- Minimum spectral gap (the bigger, the better !)

$$\Delta_{\min} = \min_{0 \leq s \leq 1} (E_1(s) - E_0(s))$$

$$T > O(1/\Delta_{\min}^2)$$

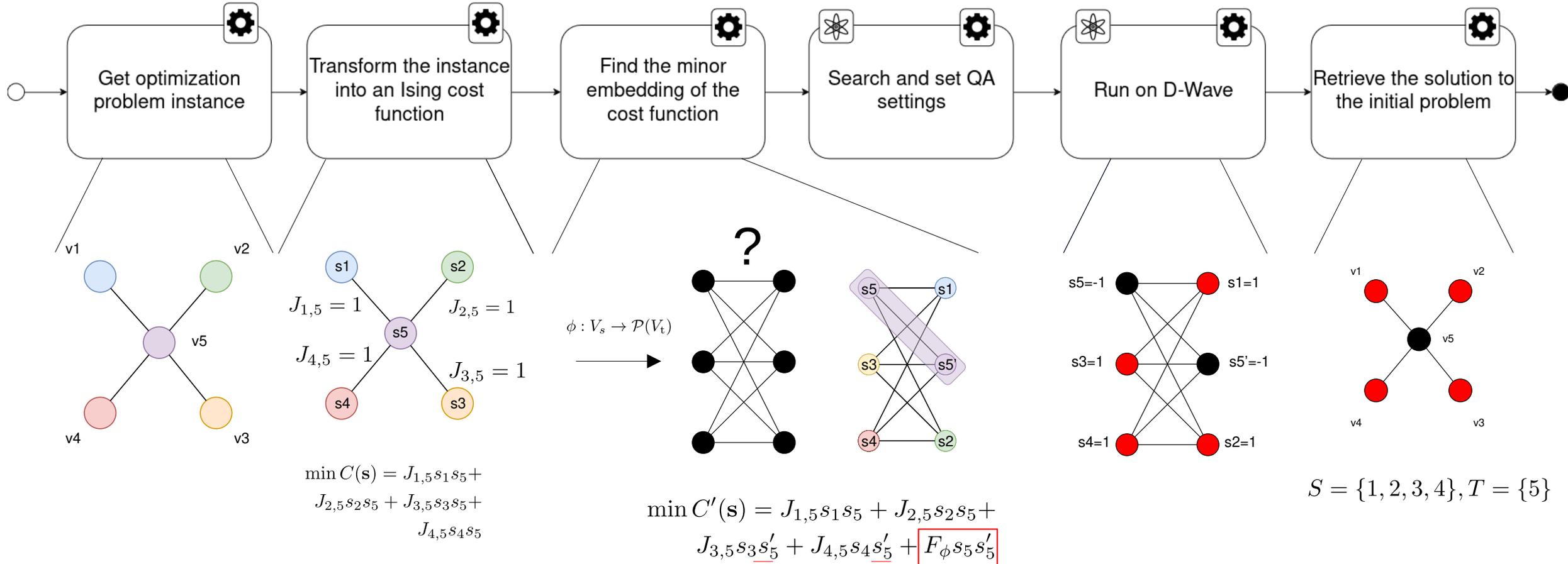
D-Wave Advantage 6.4 Annealing Schedule



I- Context & Motivations – QA programming flow

 Involves classical processing

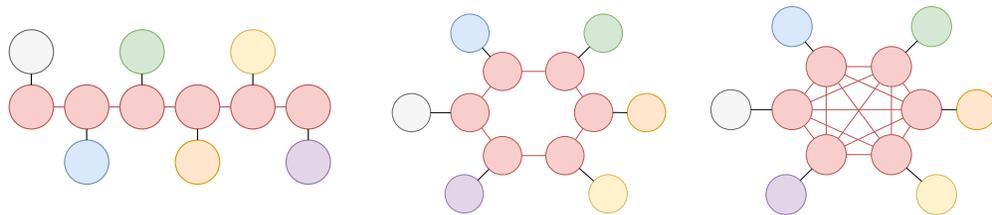
 Involves quantum processing



I- Context & Motivations – Research question

- How does the **topology of physical qubits** associated with a single logical qubit impact the chain strength and the resolution of the problem ?

- How does the arrangement of the physical qubits impact the spectral gap ?



- Which arrangement should be privileged ?

- How to efficiently set the chain strength F_ϕ ?
 - Conservation of the minimum optimal solution

$$\arg \min C = \arg \min C'$$

- Why is this value hard to set ?
 - Sweet spot usually determined experimentally. Strong but not too strong.
 - Imply auto-coupler and coupler strength rescaling.
 - Changes the problem Hamiltonian and can induce Integrated Control Errors (ICE) [1]

II- State of the Art

- Topology of Physical Qubits
 - Exploration of 4-clique minor-embedding [2]
 - Provide stronger qubits couplings
 - Reduce the chain break

- Chain strength F_ϕ
 - Upper-bound on individual qubits [3]:

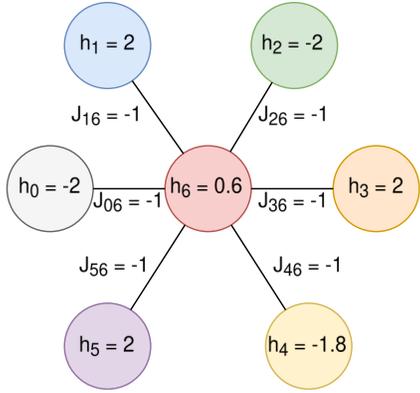
$$F_{\phi(v)} < - \left(|h_i| + \sum_{j \in \text{nbr}(i)} |J_{ij}| \right) \quad O(D)$$

- Other bounds:
 - Tighter bound in $O(DL)$ [3]
 - Exact Bound in $O(D2^L)$ [4]
- Average scaling of spin glass problems (e.g. SK problem [5] and default D-Wave method [1])
$$F_\phi \sim \tau \sqrt{n}$$
- Chain scan method [6] [7]
 - Use of the expectation value to optimize chain strength
- Other optimization approaches (Lagrangian [8] and Genetic-based [9] approaches)

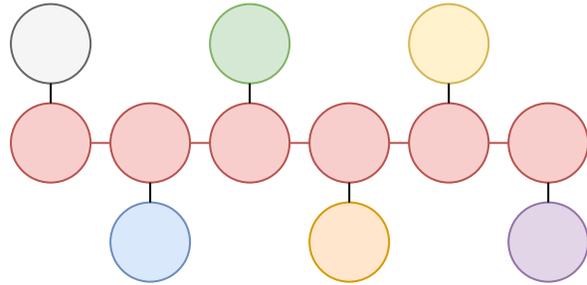
III- Shape of the logical qubit



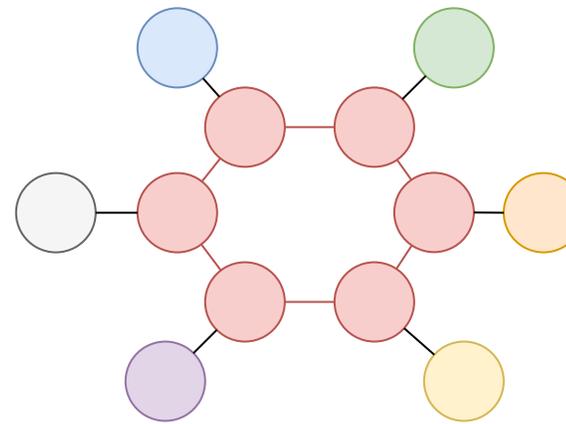
Source graph



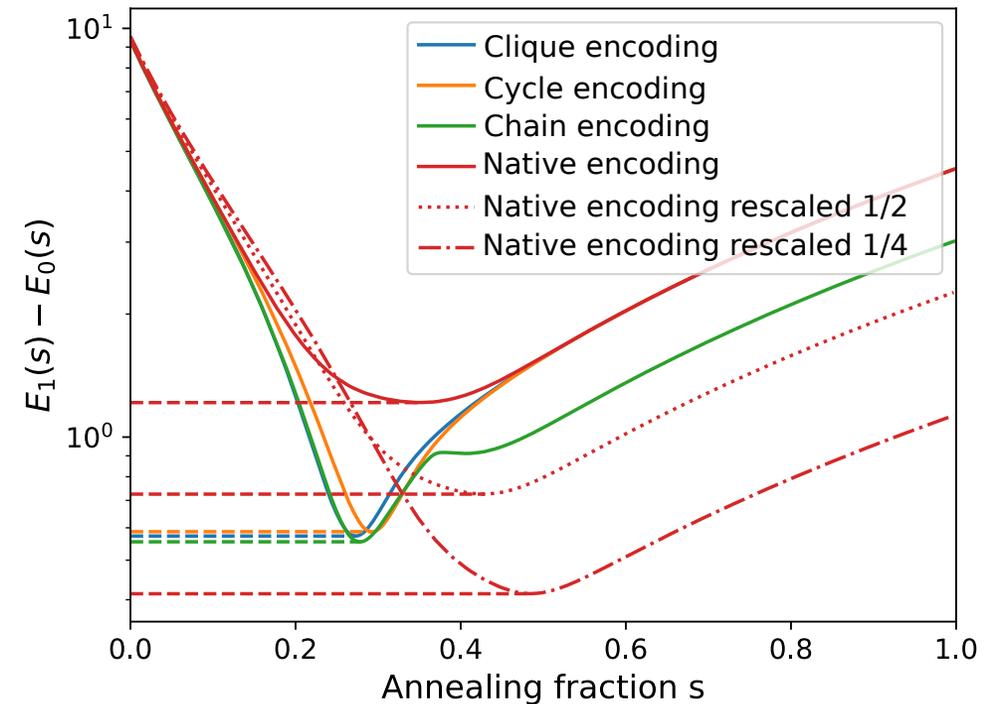
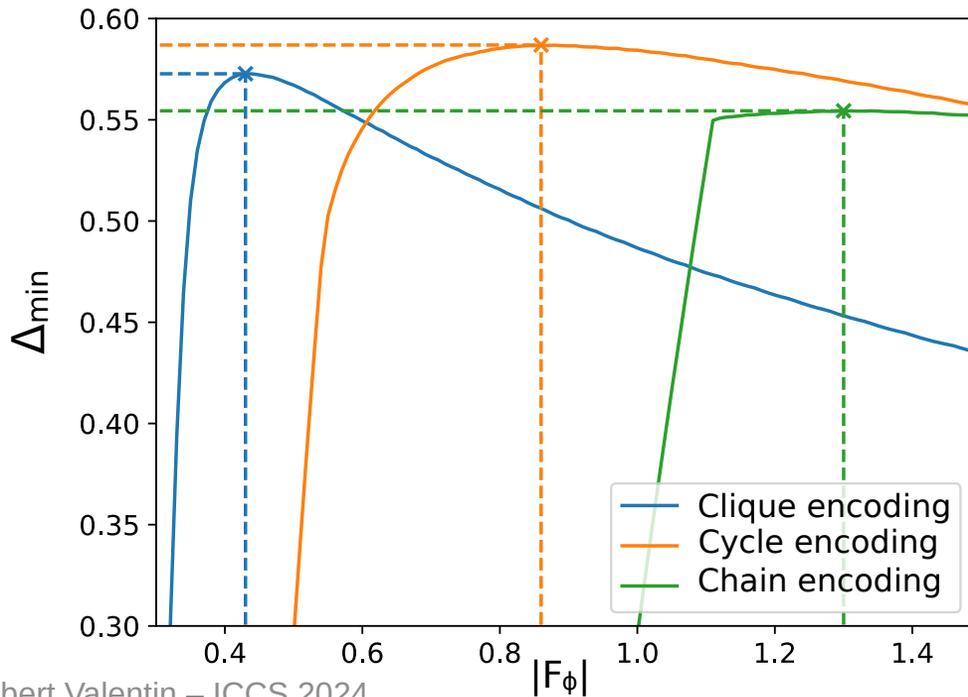
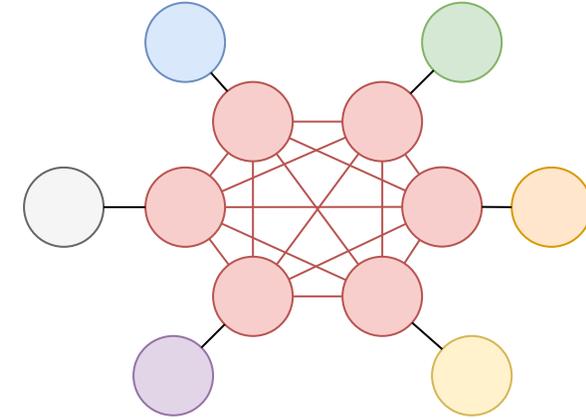
Chain encoding



Cycle encoding

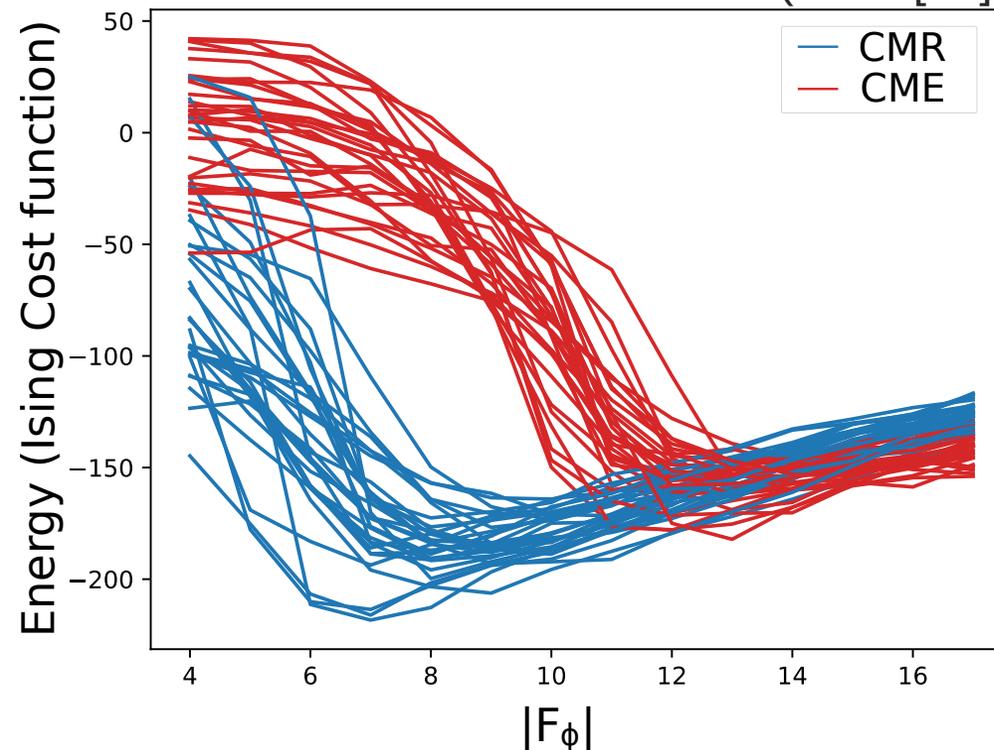


Clique encoding



III- Shape of the logical qubit

- How does the arrangement of physical qubits representing a single logical qubit impact the chain strength and the resolution of the problem ?
 - Denser structures seem to require less coupling strength to maintain ferromagnetic coupling and maximize the minimum spectral gap
 - Spectral gap difference is negligible compared to the effect of weight rescaling
- => Different embeddings with an equal number of qubits may have different optimal chain strengths.
- Chain scan for instances embedded with two different methods (CME [10] & CMR [11]):



IV- Setting the chain strength

- Measure of the chain break error rate.

$p_b(i)$: probability of having a chain break on qubit i

n_{lq} : number of logical qubits

n_s : number of shots

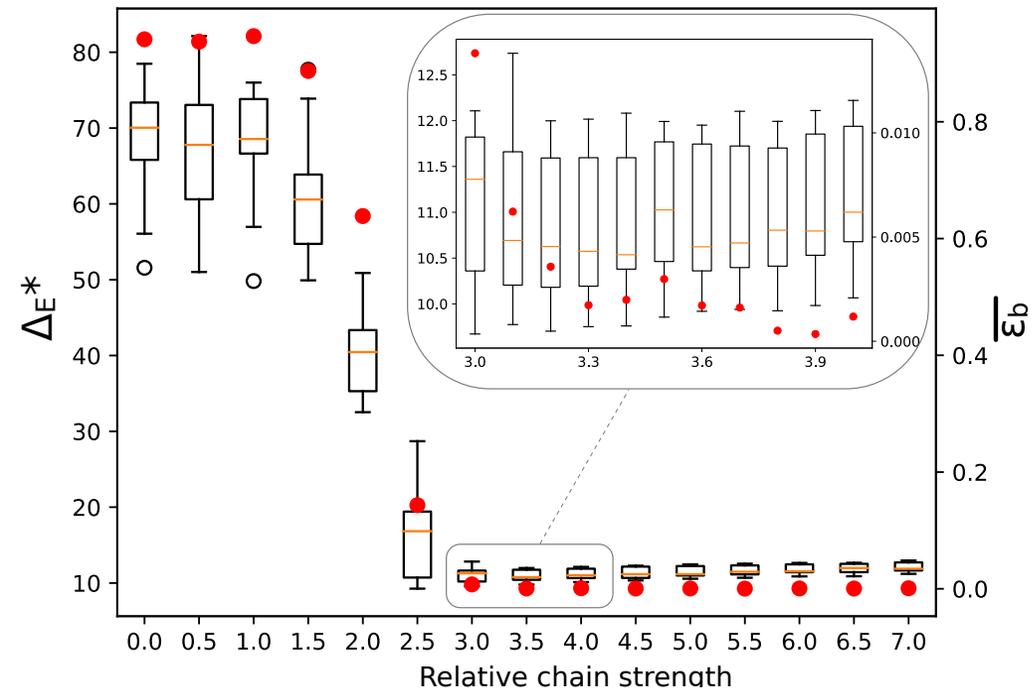
Single shot chain break error rate

$$\epsilon_b = \frac{1}{n_{lq}} \sum_{i=1}^{n_{lq}} p_b(i)$$

Average chain break error rate

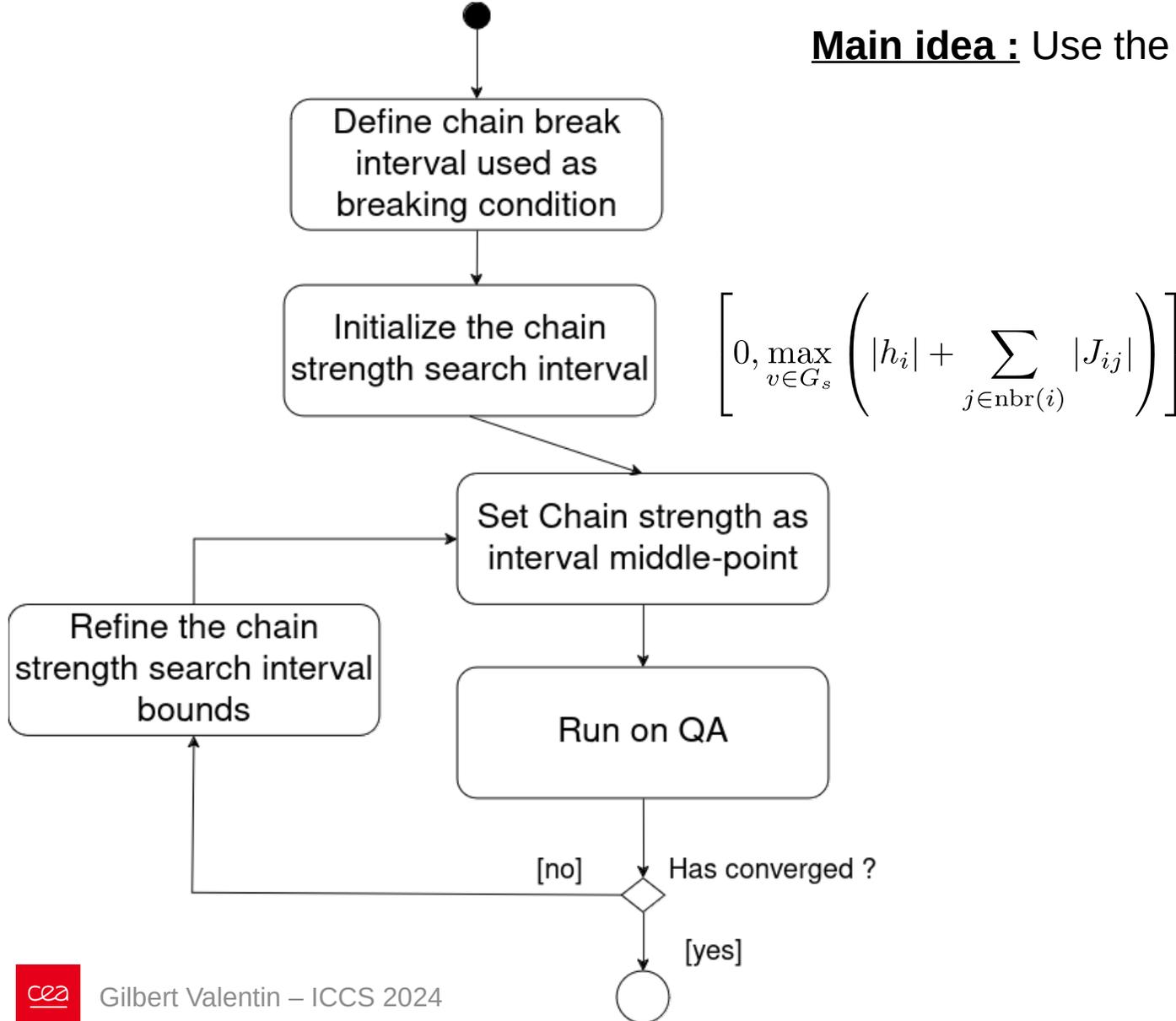
$$\bar{\epsilon}_b = \frac{1}{n_s} \sum_{i=1}^{n_s} \epsilon_b^{(i)}$$

- Correlation between the chains strength and the duplication error rate (Previous study in [12])



IV- Setting the chain strength – Binary search

Main idea : Use the chain break statistics to drive the search



- Pros of the method:
 - Universal method (can be used for any Ising cost function)
 - Fast convergence
 - Require only a few shots per run
- Cons of the method:
 - Require the setting of a precise chain break interval for the breaking condition
 - May depend on the QC
 - Mays depend on the size of the problem

IV- Setting the chain strength - Results



- 30 instances of unweighted max-cut problem for each density

- Chain break intervals:

Advantage2:

$[6 \times 10^{-3}, 2 \times 10^{-2}]$

Advantage6.4:

$[2 \times 10^{-2}, 5 \times 10^{-2}]$

- Shot / pre-processing step:
128

- Final run number of shots:

Advantage2: 3072

Advantage6.4: 4096

Advantage2_prototype2.2			Best cut size	Cut size improvement				Step
Instance size	Density	Embedding		min	max	mean	std	
$n = 40$	0.1	CMR	66.4	+0%	+0%	+0%	0%	5.4
	0.5	CMR	243	+0%	+2%	+0.2%	0%	5.8
	0.9	CMR	362.8	+2.1%	+8.2%	+5%	1.6%	4.6
$n = 80$	0.1	CMR	235.5	+0%	+0%	+0%	0%	4.7
	0.5	CME	804	+9.8%	+17.2%	+12.5%	0.2%	4.2
	0.9	CME	1435	+2%	+4.7%	+3.2%	0.6%	4.2
Advantage6.4			Best Cut size	Cut size improvement				Step
Instance size	Density	Embedding		min	max	mean	std	
$n = 100$	0.1	CMR	355.9	+0%	+0.3%	+0%	0%	4.5
	0.5	CME	1271.4	+5.6%	+14.5%	+8.8%	1.8%	2.7
	0.9	CME	2243	+1.4%	+3.7%	+2.5%	0.5%	3.7
$n = 170$	0.1	CMR	950.8	-2.1%	+0.6%	-0.5%	0.5%	2.1
	0.5	CME	3631.4	+2.8%	+6.2%	+4.5%	0.7%	2.1
	0.9	CME	6519.4	+0.4%	+1.4%	+0.8%	0.2%	3.2

V- Discussion & Perspectives

■ Conclusion

- **Optimal chain strength** seems to depend on the **shape of physical qubits** representing the logical qubit(denser mappings require lower chain strength).
- The chain strength can be optimized only using **duplication error rates**.
- Our method further optimizes the chain strength value with **few pre-processing shots** (depends on the optimal chain break interval).

■ Perspectives

- Study spectral gap tradeoffs between weight rescaling and adding new qubits to avoid the rescale.
- Identify and optimize the optimal chain break interval used by our method as the breaking condition.
- Try this method on other optimization problems (find problems for which the default D-Wave method is not efficient).
- See if the chain strength can be individually optimized
=> We made some attempts that are not conclusive



Thank You !!

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Outline

- Context & Motivation
- State of the Art
- Topology of physical qubits for a single logical qubit
- Setting the chain strength
- Results