

Distributing Multipartite Entanglement over Noisy Quantum Networks

Luís Bugalho^{1,2}, Bruno C. Coutinho², Francisco A. Monteiro^{2,3} and Yasser Omar^{1,4,5}

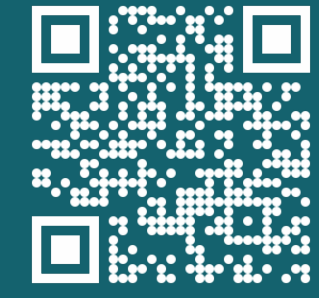
¹Instituto Superior Técnico, Universidade de Lisboa, Portugal

²Instituto de Telecomunicações, Portugal

³ISCTE - Instituto Universitário de Lisboa, Portugal

⁴PQI - Portuguese Quantum Institute, Portugal

⁵CeFEMA - Centro de Física e Engenharia de Materiais Avançados, Portugal



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Quantum Networks

A future **quantum internet** provides support for several quantum technologies, ranging from quantum secure communications to distributed quantum computation and distributed quantum metrology and sensing.

Some applications require the distribution of **multipartite entanglement**, such as quantum conference key agreement (CKA) [1], atomic clocks synchronization [2] and quantum sensor networks [3]. Optimizing its distribution, naturally impacts the applications that utilize this important resource.

Our quantum network model is a graph constituted by a set of nodes and a set of links. Each network link, connecting two neighbouring nodes, contains:

- Noisy quantum channel
- Imperfect quantum memory
- Bipartite entangled state

The protocols for entanglement generation and entanglement swapping are **probabilistic**, and each entangled pair of the network is modelled by a **Werner state**:

$$\rho_W = \gamma|\phi^+\rangle\langle\phi^+| + (1-\gamma)\mathbb{I}$$

where $\gamma = (4F - 1)/3$, and F is the fidelity of the state.

Optimising bipartite entanglement distribution over a quantum network is essentially finding the **shortest-path** in a graph. The definition of shortest-path depends on what we want to optimise for.

For bipartite entanglement we ought to maximise the **fidelity of the state** and **probability of success** and minimise the **communications time**.

However, since the fidelity of an entangled pair will depend on the generation fidelity and the decoherence, we must first optimise four parameters individually and then map them to only three:

$$[p_{m:n}, t_{m:n}, \gamma_{m:n}, \sigma_{m:n}] \mapsto [p_{m:n}, t_{m:n}, F_{m:n}]$$

$$\frac{4F_{m:n} - 1}{3} = \gamma_{m:n} e^{-t_{m:n}/\sigma_{m:n}}$$

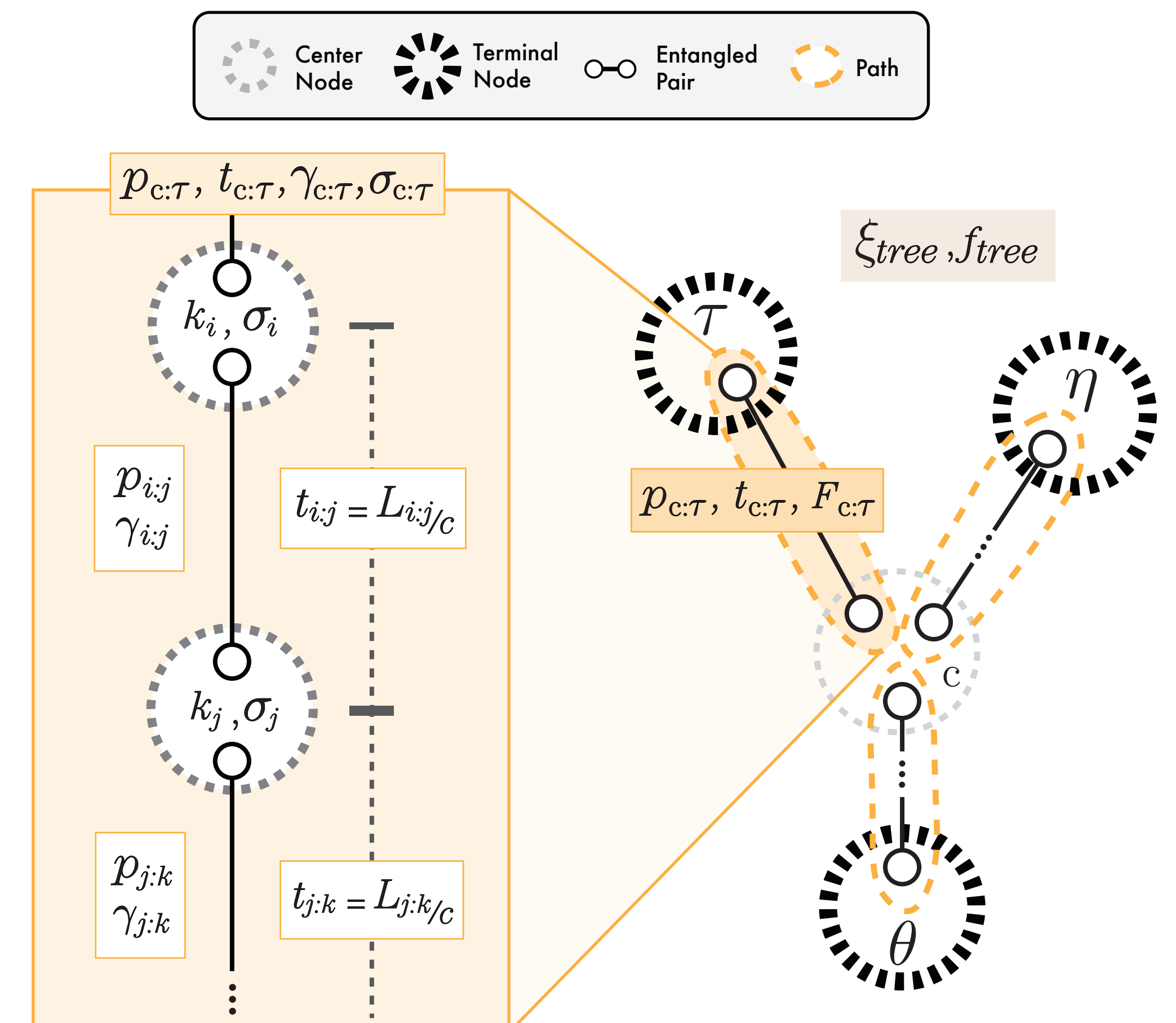


Figure 1. Branch composed by an arbitrary number of links. Each link has a fidelity $\gamma_{i,j}$, a communications time $t_{i,j}$ and probability of successful generation $p_{i,j}$. Moreover, each node has a memory coherence time σ_i and a probability of successful entanglement swapping k_i .

Figure 2. Star composed by 3 different paths, connecting terminal nodes T, η and θ to the center node C . Notice that the rate ξ_{tree} depends on each branch probability of success $p_{C:T}$ and communications time $t_{C:T}$. Moreover, the fidelity of the final multipartite state f_{tree} depends on each branch fidelity $F_{C:T}$.

Figure 4. Simulations of the MOSP algorithm and the Star Algorithm for 3 and 4 qubits GHZ state distribution in Erdős-Rényi networks with average degree 3. Complexity scaling with the number of network nodes.

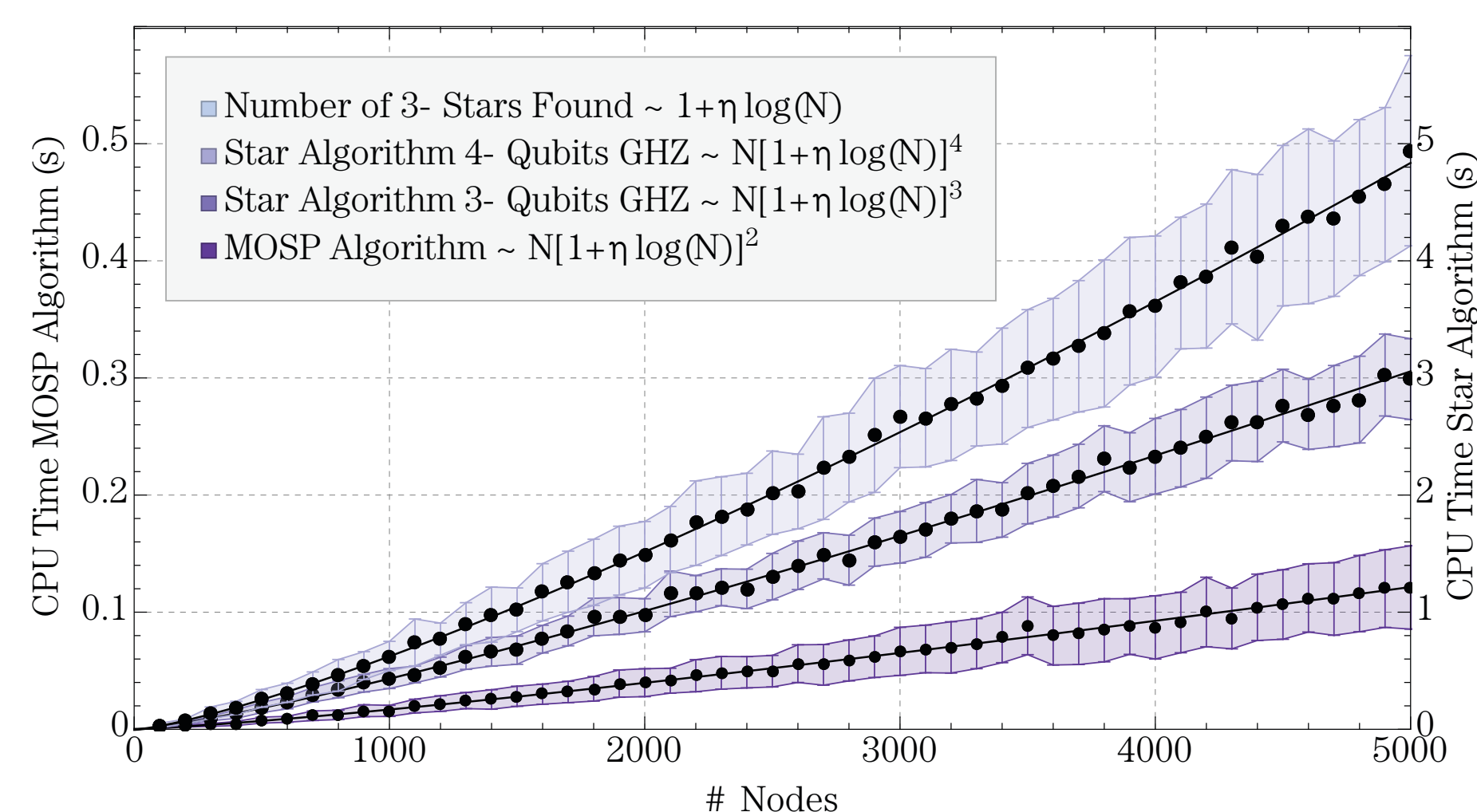
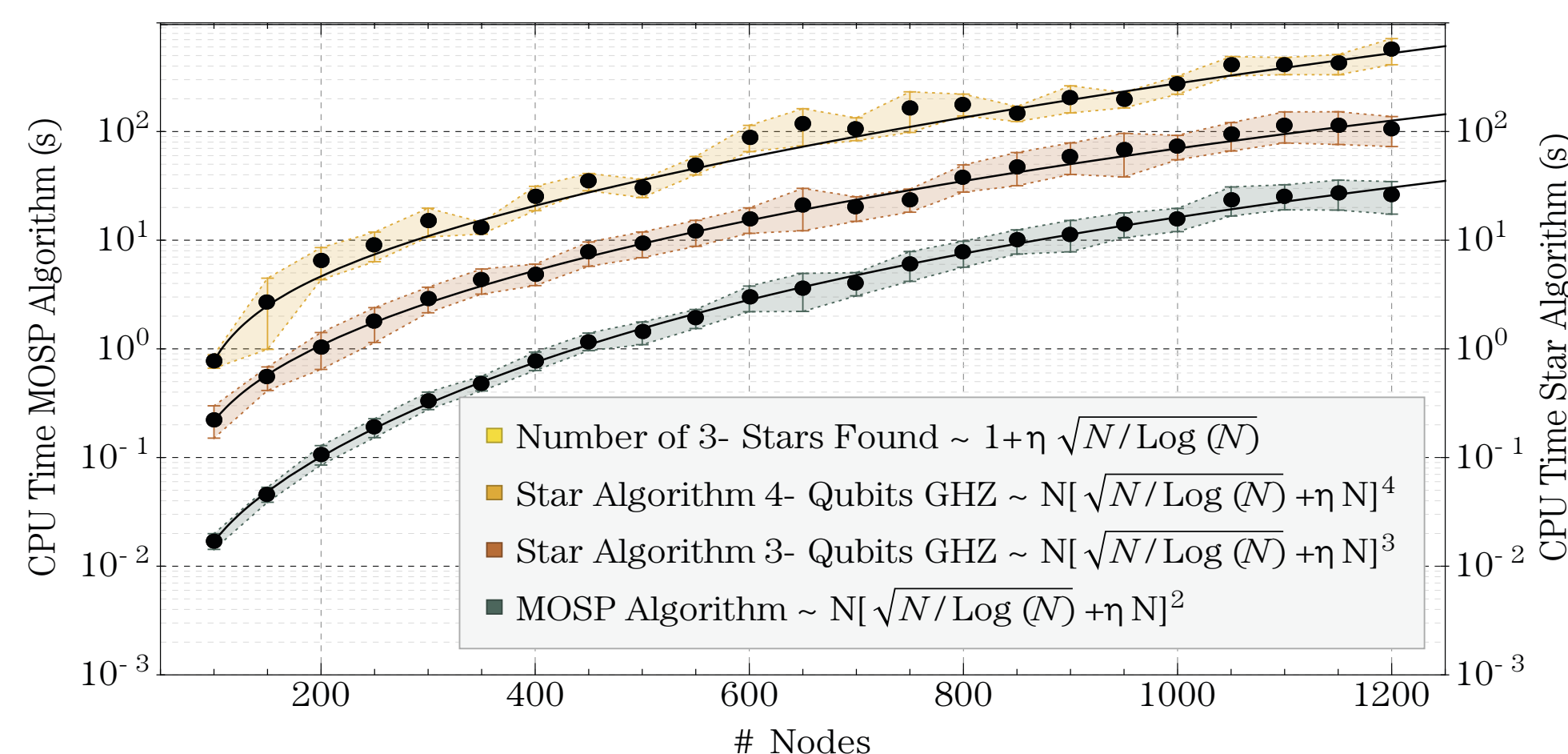


Figure 5. Simulations of the MOSP algorithm and the Star Algorithm for 3 and 4 qubits GHZ state distribution in random geometric networks with average degree 8. Complexity scaling with the number of network nodes.



Multipartite Entanglement Distribution

Similarly to [4], to distribute a 3-qubit GHZ state, first we have to find the **shortest-tree** connecting the set of terminal nodes, and then perform a set of measurements on some qubits. For 3 qubits, the shortest-tree is a **star-graph**, and the scheme is equivalent to first distributing bipartite entanglement between every terminal and a center node and then performing a set of operations on the center node.

Thus, the central idea is **finding the center node** for which this results in the optimal solution. Optimality is defined concerning the **metrics for the distribution** of the final multipartite entangled state, namely the **rate of distribution** and the **fidelity**.

The average time for distribution depends on the maximum of each path's communications time multiplied by the average number of attempts before the first success. The **average rate** would then be the inverse of:

$$T_{tree} = 2 \cdot \frac{\max_{\tau \in \mathcal{T}} \{t_{C:\tau}\}}{\prod_{\tau \in \mathcal{T}} p_{C:\tau}}, \quad \xi_{tree} = \frac{1}{T_{tree}}$$

The form of the final distributed state is equivalent to applying a depolarising channel to each qubit of the GHZ state, correspondent to each branch connecting the center node to a terminal. The **final fidelity of this state** is described by:

$$f_{tree} = \frac{1}{2} \left[\prod_{\tau \in \mathcal{T}} \frac{1 + 2F_{\tau}}{3} + \prod_{\tau \in \mathcal{T}} \frac{2(1 - F_{\tau})}{3} + \prod_{\tau \in \mathcal{T}} \frac{4F_{\tau} - 1}{3} \right]$$

Multi-objective Routing Algorithm

In our problem, we want to optimise two distinct parameters. For that, we require a multi-objective approach. In 1984, Martins [5] introduced **multi-objective shortest-paths** (MOSP) algorithms, defining the dominance relation and using the **Pareto optimality** definition – optimal paths are non-dominated paths. This results in a set of optimal solutions that is usually larger than one – the **Pareto front**.

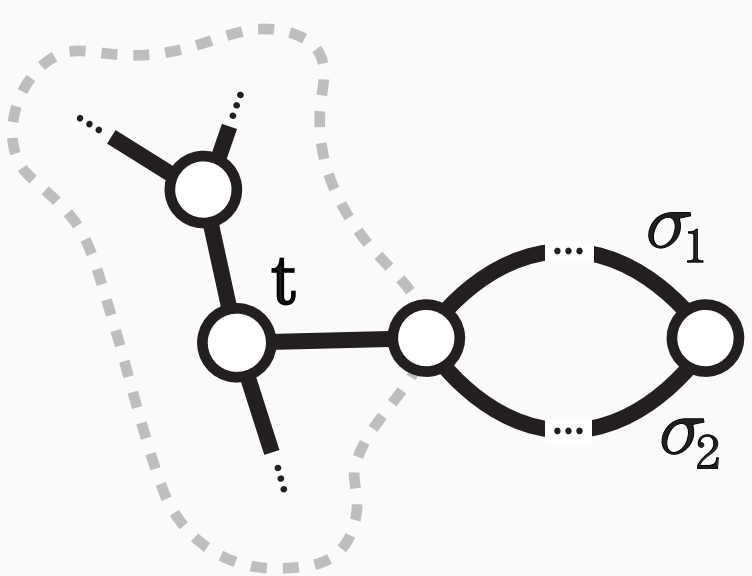
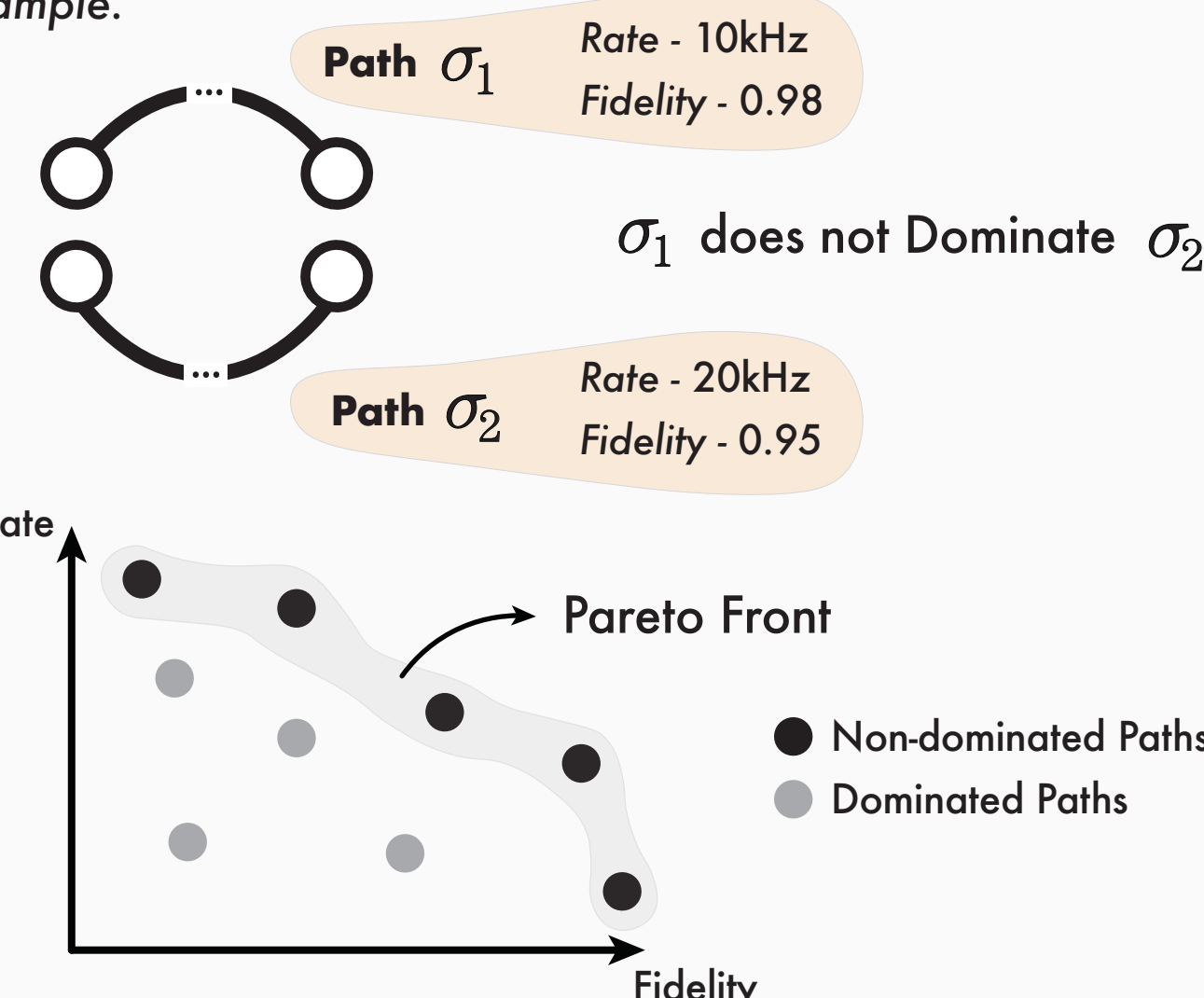


Figure 3. Label-Isotonicity: for every tree and path, when extending a tree with a path, choosing a better path results in a better tree.

Example.



In [6] Sobrinho introduces a more fundamental definition for these metrics involved in routing problems – **algebras for routing**. These algebras incorporate the parameters of the links, paths and how paths are extended with links – **the metrics** –, allowing for a bigger level of abstraction.

Multi-objective routing – **one algebra for each parameter** (rate, fidelity,...). To prove optimality, we expanded these algebras for the case of trees (instead of paths) and introduced **label-isotonicity** (Figure 3).

Results and Conclusions

> We were able to calculate the effect of **imperfect entanglement** throughout quantum networks on the distribution of GHZ and W states, importante resources for distributed applications, such as **communication** and **sensing**;

> We defined a metric for the fidelity and realise an algorithm with **polynomial runtime** for distributing these states optimally, **maximising simultaneously the fidelity and the rate of distribution**;

> Using classical techniques and tools, we can **prove the optimality** and create the necessary formalism to expand the methodology, while also providing the **necessary conditions to attain this optimality**.

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For further details:

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About the Authors:

