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Node Centrality for Continuous-Time Quantum Walks

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Abstract. The study of complex networks has recently attracted increasing interest because of the large variety of systems that can be modeled using graphs. A fundamental operation in the analysis of complex networks is that of measuring the *centrality* of a vertex. In this paper, we propose to measure vertex centrality using a continuous-time quantum walk. More specifically, we relate the importance of a vertex to the influence that its initial phase has on the interference patterns that emerge during the quantum walk evolution. To this end, we make use of the quantum Jensen-Shannon divergence between two suitably defined quantum states. We investigate how the importance varies as we change the initial state of the walk and the Hamiltonian of the system. We find that, for a suitable combination of the two, the importance of a vertex is almost linearly correlated with its degree. Finally, we evaluate the proposed measure on two commonly used networks.

Key words: Vertex Centrality, Complex Network, Quantum Walk, Quantum Jensen-Shannon Divergence

1 Introduction

In recent years, an increasing number of researchers have turned their attention to the study complex networks [1]. Complex network are ubiquitous in a large number of real-world systems. A non-exhaustive list of examples includes metabolic networks [2], protein interactions [3], brain networks [4] and scientific collaboration networks [5]. A fundamental task in complex network analysis is that of measuring the centrality of a vertex, i.e., its importance. To this end, a number of centrality indices have been introduced in the literature [1, 6–9]. Each of these captures different but equally significant aspects of vertex importance.

Perhaps the most intuitive centrality measure is degree centrality [7]. This is defined as the number of links incident upon a node, i.e., the degree of the node. The degree centrality naturally interprets the number of edges incident on a vertex as a measure of its “popularity”, or, alternatively, as the risk of a node being infected in an epidemiological scenario. Closeness centrality [10], on the other hand, links the importance of a vertex to its proximity to the remaining vertices

of the graph. More precisely, the closeness centrality is defined as the inverse of the sum of the distance of a vertex to the remaining nodes of the graph, i.e., $CC(u) = \frac{n-1}{\sum_{v=1}^n d(u,v)}$ where $d(u,v)$ denotes the shortest path distance between nodes u and v . The betweenness centrality [7] is a measure of the extent to which a given vertex lies on the paths between the remaining vertices, where the path may be either that of shortest length or a random walk between the nodes. If $sp(v_1, v_2)$ denotes the number of shortest paths from node v_1 to node v_2 , and $sp(v_1, u, v_2)$ denotes the number of shortest paths from v_1 to v_2 that pass through node u , the betweenness centrality of u is $BC(u) = \sum_{v_1=1}^n \sum_{v_2=1}^n \frac{sp(v_1, u, v_2)}{sp(v_1, v_2)}$. Note that this definition assumes that the communication takes place along the shortest path between two vertices. A number of measures have been introduced to account for alternative scenarios in which the information is allowed to flow through different paths [1, 6–8].

Recently, there has also been a surge of interest in using quantum walks as a primitive for designing novel quantum algorithms on graph structures [11]. Quantum walks on graphs represent the quantum mechanical analogue of the classical random walk on a graph. Despite being similar in their definition, the dynamics of the two walks can be remarkably different. In the classical case the evolution of the walk is governed by a double stochastic matrix, while in the quantum case the evolution is governed by a unitary matrix, thus rendering the walk reversible and non-ergodic. Moreover, the state vector of the classical random walk is real-valued, while in the quantum case the state vector is complex-valued. As there is no constraint on the sign and phase of the amplitudes, different paths are allowed to interfere with each other in both constructive and destructive ways. This in turn gives rise to faster hitting times and reduces the problems of tottering observed in classical random walks [11].

In this paper, we propose to measure the centrality of a vertex using a continuous-time quantum walk. More specifically, we relate the importance of a vertex to the influence that its initial phase has on the evolution of a suitably defined quantum walk. To this end, we make use of the quantum Jensen-Shannon divergence, a recently introduced generalisation of the classical Jensen-Shannon divergence to quantum states [12]. Just as the classical Jensen-Shannon divergence [13], the quantum Jensen-Shannon divergence is symmetric, bounded and always defined. From a physical perspective, the QJSD is computed from density matrices, whose entries are observables. As a consequence, it should be possible, at least in theory, to design a quantum algorithm to compute the QJSD centrality that could benefit from the power of quantum computers. However, the design of such an algorithm is beyond the scope of this paper.

The remainder of this paper is organised as follows: Section 2 provides an essential introduction to the basic terminology required for understanding the proposed quantum mechanical framework. With these notions to hand, we introduce our centrality measure in Section 3 and we study its properties. In Section 4 we apply the proposed measure to the analysis of two commonly used network models, while the conclusions are presented in Section 5.

2 Quantum Mechanical Background

The continuous-time quantum walk [14] is a natural quantum analogue of the classical random walk. Given a graph $G = (V, E)$, classical random walks model a diffusion process over the node set V , and have proven to be a useful tool in the analysis of its structure. Similarly, the continuous-time quantum walk is defined as a dynamical process over the vertices of the graph. By contrast to the classical case, where the state vector is constrained to lie in a probability space, in the quantum case the state of the system is defined through a vector of complex amplitudes over the node set V whose squared norm sums to unity over the nodes of the graph, with no restriction on their sign or complex phase. These phase differences allow interference effects to take place. Moreover, in the quantum case the evolution of the state vector of the walker is governed by a complex valued unitary matrix, whereas the dynamics of the classical random walk is governed by a stochastic matrix. Hence the evolution of the quantum walk is reversible, implying that quantum walks are non-ergodic and do not possess a limiting distribution. As a result, the behaviour of classical and quantum walks differs significantly, and quantum walks possess a number of interesting properties not exhibited by classical random walks.

More formally, using the Dirac notation, we denote the basis state corresponding to the walk being at vertex $u \in V$ as $|u\rangle$. A general state of the walk is a complex linear combination of the basis states, such that the state of the walk at time t is defined as

$$|\psi_t\rangle = \sum_{u \in V} \alpha_u(t) |u\rangle \quad (1)$$

where the amplitude $\alpha_u(t) \in \mathbb{C}$ and $|\psi_t\rangle \in \mathbb{C}^{|V|}$ are both complex.

At each instant in time the probability of the walker being at a particular vertex of the graph is given by the square of the norm of the amplitude of the relative state. Let X^t be a random variable giving the location of the walker at time t . Then the probability of the walker being at the vertex u at time t is given by $\Pr(X^t = u) = \alpha_u(t)\alpha_u^*(t)$, where $\alpha_u^*(t)$ is the complex conjugate of $\alpha_u(t)$. Moreover $\sum_{u \in V} \alpha_u(t)\alpha_u^*(t) = 1$ and $\alpha_u(t)\alpha_u^*(t) \in [0, 1]$, for all $u \in V$, $t \in \mathbb{R}^+$.

The evolution of the walk is then given by the Schrödinger equation, where we take the time-independent Hamiltonian of the system to be the graph Laplacian, yielding

$$\frac{\partial}{\partial t} |\psi_t\rangle = -iL |\psi_t\rangle . \quad (2)$$

Given an initial state $|\psi_0\rangle$, we can solve Eq. (2) to determine the state vector at time t

$$|\psi_t\rangle = e^{-iLt} |\psi_0\rangle . \quad (3)$$

Note that generally one may use any Hermitian operator as the Hamiltonian. Common choices are the graph adjacency matrix, the normalised Laplacian and the signless Laplacian.

Finally, we can compute the spectral decomposition of the graph Laplacian $L = \Phi \Lambda \Phi^\top$, where Φ is the $n \times n$ matrix $\Phi = (\phi_1 | \phi_2 | \dots | \phi_j | \dots | \phi_n)$ with the ordered eigenvectors ϕ_j s of L as columns and $\Lambda = \text{diag}(\lambda_1, \lambda_2, \dots, \lambda_j, \dots, \lambda_n)$ is the $n \times n$ diagonal matrix with the ordered eigenvalues λ_j of L as elements, such that $0 = \lambda_1 \leq \lambda_2 \leq \dots \leq \lambda_n$. Using the spectral decomposition of the graph Laplacian and the fact that $\exp[-iLt] = \Phi \exp[-i\Lambda t] \Phi^\top$ we can then write

$$|\psi_t\rangle = \Phi e^{-i\Lambda t} \Phi^\top |\psi_0\rangle. \quad (4)$$

2.1 Quantum Jensen-Shannon Divergence

The *density operator* (or *density matrix*) is introduced in quantum mechanics to describe a system whose state is an ensemble of pure quantum states $|\psi_i\rangle$, each with probability p_i . The density operator of such a system is defined as

$$\rho = \sum_i p_i |\psi_i\rangle \langle \psi_i|. \quad (5)$$

The von Neumann entropy [15] H_N of a density operator ρ is defined as

$$H_N = -\text{tr}(\rho \log \rho) = -\sum_i \xi_i \ln \xi_i \quad (6)$$

where ξ_1, \dots, ξ_n are the eigenvalues of ρ . If $\langle \psi_i | \rho | \psi_i \rangle = 1$, i.e., the quantum system is a pure state $|\psi_i\rangle$ with probability $p_i = 1$, then the Von Neumann entropy $H_N(\rho) = -\text{tr}(\rho \log \rho)$ is zero. On other hand, for a mixed state described by the density operator σ we have a non zero Von Neumann entropy associated with it.

With the Von Neumann entropy to hand, the quantum Jensen-Shannon divergence between two density operators ρ and σ is defined as

$$D_{JS}(\rho, \sigma) = H_N\left(\frac{\rho + \sigma}{2}\right) - \frac{1}{2}H_N(\rho) - \frac{1}{2}H_N(\sigma) \quad (7)$$

This quantity is always well defined, symmetric and positive definite. Finally, it can also be shown that $D_{JS}(\rho, \sigma)$ is bounded, i.e., $0 \leq D_{JS}(\rho, \sigma) \leq 1$.

3 QJSD Centrality

In order to measure the centrality of vertex v , we define two quantum walks where v is initially set to be in phase and in antiphase with the respect to the remaining nodes. Let the normalised graph Laplacian be the Hamiltonian of our system, and let $|\psi_0^{v-}\rangle = \sum_{u \in V} \alpha_u^{v-}(0) |u\rangle$ and $|\psi_0^{v+}\rangle = \sum_{u \in V} \alpha_u^{v+}(0) |u\rangle$ denote the quantum walks on G with initial amplitudes

$$\alpha_j^{v-}(0) = \begin{cases} -\frac{\sqrt{d_j}}{C} & \text{if } j = v \\ +\frac{\sqrt{d_j}}{C} & \text{otherwise} \end{cases} \quad \alpha_j^{v+}(0) = \begin{cases} +\frac{\sqrt{d_j}}{C} & \forall j \end{cases} \quad (8)$$

where C is the normalisation constant such that probabilities sum to 1. In other words, we define the initial amplitude to be proportional to the square root of the node degrees. Finally, let ρ_{v+} and ρ_{v-} be the density operators which describe the ensembles of quantum states $|\psi_t^{v-}\rangle$ and $|\psi_t^{v+}\rangle$ respectively, i.e.,

$$\rho_{v-} = \lim_{T \rightarrow \infty} \frac{1}{T} \int_0^T |\psi_t^{v-}\rangle \langle \psi_t^{v-}| dt \quad \rho_{v+} = \lim_{T \rightarrow \infty} \frac{1}{T} \int_0^T |\psi_t^{v+}\rangle \langle \psi_t^{v+}| dt \quad (9)$$

Given this setting, we can measure how the initial phase of the vertex v affects the evolution of the quantum walks by computing the distance between the quantum states defined by ρ_{v-} and ρ_{v+} . That is, we define the quantum Jensen-Shannon divergence (QJSD) centrality of a vertex v as

$$C_{QJSD}(v) = D_{JS}(\rho_{v-}, \rho_{v+}) \quad (10)$$

Note that the computational complexity of the QJSD centrality is bounded by that of computing the eigendecomposition of the graph Laplacian, i.e., $O(n^3)$. Let $\Phi \Lambda \Phi^\top$ be the spectral decomposition of the graph normalised Laplacian and let $P_\lambda = \sum_{k=1}^{\mu(\lambda)} \phi_{\lambda,k} \phi_{\lambda,k}^\top$ be the projection operator on the subspace spanned by the $\mu(\lambda)$ eigenvectors $\phi_{\lambda,k}$ associated with the eigenvalue λ of the graph normalised Laplacian. Rossi et al. [16] have shown that $\rho_\infty = \sum_{\lambda=1}^m P_\lambda \rho_0 P_\lambda^\top$, where m denotes the number of unique eigenvalues of the graph normalised Laplacian. Note that as a consequence of Eq. ?? we have that ρ_{v-} and ρ_{v+} are simultaneously diagonalisable. That is, there exist a single invertible matrix M such that $M^{-1} \rho_{v-} M$ and $M^{-1} \rho_{v+} M$ are diagonal. More precisely, here $M = \Phi$, the $n \times n$ matrix $\Phi = (\phi_1 | \phi_2 | \dots | \phi_j | \dots | \phi_n)$ with the ordered eigenvectors ϕ_j s of the Hamiltonian as columns.

3.1 Relation with Degree Centrality

We are now interested in studying the relation between the QJSD centrality and the degree centrality. It has been shown, for example, that the degree and the betweenness centrality are highly correlated [17]. This should not come as a surprise, as we expect high degree vertices to be more often included in the shortest path along a pairs of vertices.

Let the initial states of the walks be defined as in Eq. 8 and let the normalised Laplacian be the Hamiltonian of our system. We start by observing that $|\psi_0^{v+}\rangle = \sum_{u \in V} \alpha_u^{v+}(0) |u\rangle$ corresponds to the eigenvector ϕ_0 associated with the zero eigenvalue of the Hamiltonian, and as a consequence $|\psi_0^{v+}\rangle$ will remain constant over time. In other words, we have that $\rho_{v+} = |\psi_0^{v+}\rangle \langle \psi_0^{v+}|$. Note that the spectrum of ρ_{v+} is composed of a single eigenvector ϕ_0 with eigenvalue equal to 1. Moreover, recall from Eq. ?? that ρ_{v-} and ρ_{v+} are co-diagonalisable matrices. As a result, each eigenvalue of $\rho_{v-} + \rho_{v+}$ is a sum of eigenvalues of ρ_{v-} and ρ_{v+} . More precisely, when the two walks are initialised as in Eq. 8, all the eigenvalues μ_i of $\frac{\rho_{v-} + \rho_{v+}}{2}$ will be equal to the eigenvalues of ρ_{v-} , except for the eigenvalue $\mu_0 + 1$ which is associated to the common eigenvector ϕ_0 .

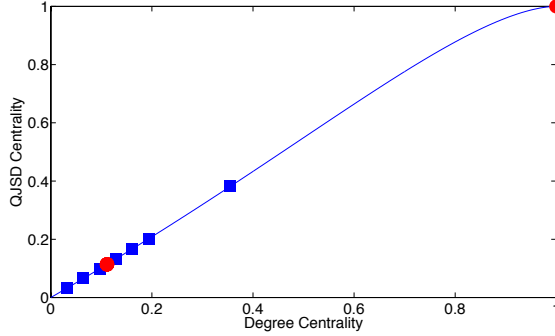


Fig. 1. The correlation between degree and QJSD centrality, for a star graph (red dots) and a scale-free graph (blue squares). The blue line shows the predicted dependency between the two centrality indices.

We now show that, as a consequence of this, the QJSD centrality is proportional to the degree centrality. Note that since ρ_{v+} has a single non-zero eigenvalue which is equal to 1, we have that $H_N(\rho_{v+}) = 0$. As a consequence of this and of Eq. 7, we have that

$$\begin{aligned}
 D_{JS}(\rho_{v-}, \rho_{v+}) &= H_N\left(\frac{\rho_{v-} + \rho_{v+}}{2}\right) - \frac{1}{2}H_N(\rho_{v-}) \\
 &= -\frac{\mu_0 + 1}{2} \log_2 \frac{\mu_0 + 1}{2} - \sum_{i \neq 0} \frac{\mu_i}{2} \log_2 \frac{\mu_i}{2} + \frac{1}{2} \sum_i \mu_i \log_2 \mu_i \\
 &= \frac{\mu_0 + 1}{2} - \frac{\mu_0 + 1}{2} \log_2(\mu_0 + 1) + \sum_{i \neq 0} \frac{\mu_i}{2} - \frac{1}{2} \sum_{i \neq 0} \mu_i \log_2 \mu_i + \frac{1}{2} \sum_i \mu_i \log_2 \mu_i \\
 &= 1 - \frac{1}{2} \log_2(\mu_0 + 1) + \frac{\mu_0}{2} \log_2 \frac{\mu_0}{\mu_0 + 1} \tag{11}
 \end{aligned}$$

where μ_i denotes the i th eigenvalue of ρ_{v-} and we used the fact that $\sum_i \mu_i = 1$. We now proceed to show that μ_0 is proportional to the degree of node v , and therefore the QJSD centrality is proportional to the degree centrality. In fact, we have that

$$\mu_0 = \langle \phi_0 | \rho_0 | \phi_0 \rangle = \langle \phi_0 | \psi_0^{v-} \rangle^2 = \left(1 - \frac{d_v}{|E|}\right)^2 \tag{12}$$

where d_v is the degree of v and $|E|$ denotes the number of edges in the graph.

In other words, when we take the normalised Laplacian as our Hamiltonian and we initialise the walks according to Eq. 8, the QJSD centrality turns out to be quasi-linearly correlated with the degree centrality. Fig. 1 shows the correlation between the QJSD centrality and the degree centrality for a scale-free random graph and a star graph. Recall that the degree centrality is normalised between 0 and 1 by dividing it by $|V|(|V| - 1)$, i.e., the maximum cardinality of the

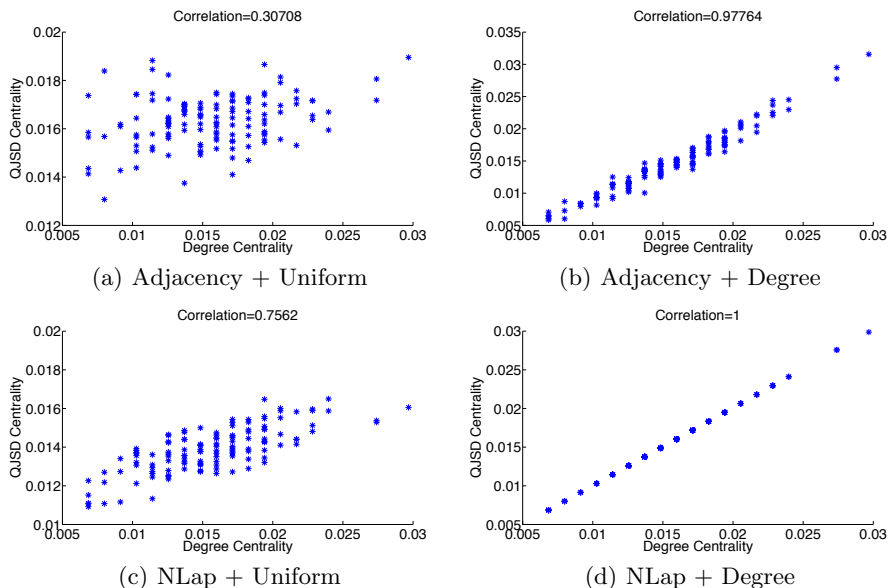


Fig. 2. Correlation between the QJSD centrality and the degree centrality for different choices of the Hamiltonian (adjacency matrix or normalised Laplacian) and of the initial state (normalised uniform distribution or normalised degree distribution).

edge set. Note that the non-linearity of the correlation becomes evident only for those nodes with degree close to $|E|$, for which we have that $\frac{d_v}{|E|} \approx 0$ and thus $\mu_0 \approx 1 + \frac{d_v^2}{|E|}$.

So far we assumed that the Hamiltonian of the quantum walk is the graph normalised Laplacian. However, any Hermitian operator encoding the structure of the graph can be chosen as an alternative. Similarly, there is no constraint on the initial state of the walk, as long as it is a valid amplitude vector. Fig. 2 shows the correlation between the QJSD centrality and the degree centrality computed on a stochastic Kronecker graph for different choices of the initial state and the Hamiltonian. More specifically, we let the Hamiltonian be either the adjacency matrix or the normalised Laplacian of the graph, while the initial state is either proportional to the node degree as in Eq. 8 or uniformly equal to $1/\sqrt{n}$, where n denotes the number of nodes in the graph. As expected, our centrality measure is strongly correlated with the degree centrality when the Hamiltonian is the graph normalised Laplacian and the initial state is proportional to the node degree (see Fig. 2(d)). In general, we see that when the starting state is proportional to the node degree, the correlation tends to be very high, while the choice of a uniform initial state leads to a value of the centrality which is less dependent on the node degree.

Hence, in an attempt to capture structural information which are not trivially revealed by examining the node degree, we explore the consequences of letting the

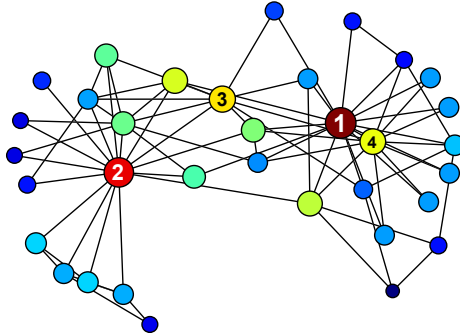


Fig. 3. Zachary’s karate club network, where we have drawn each node with a diameter that is proportional to its QJSD centrality.

walk start from a uniform amplitude vector and choosing the adjacency matrix as the Hamiltonian. Moreover, in order to balance the strength of the positive and negative signals, i.e., the contribution of the node amplitudes with either positive or negative phases, we let the magnitude of the initial amplitude on the node being analysed be equal to the sum of the amplitudes on the remaining nodes, which gives the initial state

$$\alpha_j^{v-}(0) = \begin{cases} -\frac{1}{\sqrt{2}} & \text{if } j = v \\ +\frac{1}{\sqrt{2(|V|-1)}} & \text{otherwise} \end{cases} \quad \alpha_j^{v+}(0) = \begin{cases} +\frac{1}{\sqrt{2}} & \text{if } j = v \\ +\frac{1}{\sqrt{2(|V|-1)}} & \text{otherwise} \end{cases} \cdot \quad (13)$$

4 Experimental Evaluation

We now apply the QJSD centrality to a pair of commonly used network datasets, namely Zachary’s karate club [18] and Padgett’s network of marriages between the 16 most eminent Florentine families in the 15th century [19]. Fig. 3 shows Zachary’s karate club network, where each vertex is drawn with a diameter that is proportional to the QJSD centrality. We see that there are two main actors, node #1 and node #2, which correspond to the instructor and the administrator of the club. Note that using our measure the instructor turns out to be the node with the highest centrality, which is also the most central according to the degree centrality, while the betweenness centrality elects the administrator as the most important node. However, the betweenness centrality indicates as the second most important actor node #3, as this vertex has many contacts with both the members of the administrator cluster and the members of the instructor cluster and thus it is misunderstood as a center by the betweenness centrality. Finally, node #4 is identified as the third most important by the degree centrality, leaving node #3 at the fourth place, although the latter is more central in the sense that it shares many links with both groups.

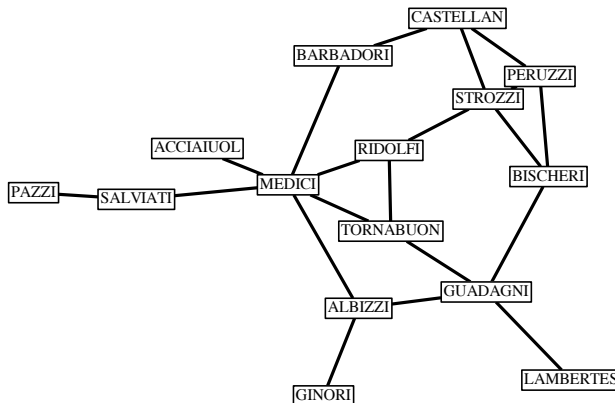


Fig. 4. Padgett’s network of marriages between eminent Florentine families in the 15th century [19]. We omit the Pucci, which had no marriage ties with other families.

Family	Centrality	Family	Centrality	Family	Centrality
Medici	0.4867	Castellan	0.3245	Salviati	0.2248
Ridolfi	0.4619	Barbadori	0.3205	Ginori	0.1993
Strozzi	0.4192	Albizzi	0.3172	Acciaiuol	0.1534
Tornabuon	0.4041	Guadagni	0.3091	Lambertes	0.1267
Bischeri	0.3586	Peruzzi	0.2990	Pazzi	0.1126

Table 1. The QJSD centrality of the families of Padgett’s network [19].

Padgett’s network of marriages is depicted in Fig. 4. In Table 1, we show the ranking of the 15 families according to their QJSD centrality. As expected, the Medici easily outperform the Strozzi, who are their main rivals. This agrees with the historical view that Medici’s supremacy was largely due to their skills in manipulating the marriage network. Interestingly the Pazzi, which is the most loosely connected family of the graph, achieve the lowest centrality. Note also that the Ridolfi family, which connect two of the most influential families at that time, the Medici and the Strozzi, is assigned a high centrality. Moreover, the Tornabuon, which form a tightly connected clique together with the Medici and the Ridolfi, is the fourth most central node of the network.

5 Conclusions

In this paper, we have proposed to measure vertex centrality using a continuous-time quantum walk. We measured the importance of a vertex as the influence that its initial phase has on the interference patterns that emerge during the quantum walk evolution. We have showed that, under particular settings, the resulting centrality measure is almost linearly correlated with degree centrality.

Thus, we have proposed an alternative starting state where the contribution of the node amplitudes with positive and negative phases is equal. Finally, we have evaluated the resulting measure to two commonly used network models.

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