Peer-to-Peer Systems Final Project Report

Andrea Maggiordomo – mggndr89[at]gmail.com Master degree in Computer Science, Università di Pisa

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1 Introduction

The goal of the project was to define Deccen, a decentralized algorithm to compute stress centrality indices in a network based on the framework introduced in Lehmann and Kaufmann [8], and study its behavior using the PeerSim [10] simulator. Additionally – given the high cost of the algorithm – approximation techniques could be proposed in order to reduce the communication overhead and memory requirements, and the algorithm could be extended to also compute closeness and betweenness indices.

This report defines a version of Deccen capable of computing all the three centrality measures mentioned. It also defines a different decentralized algorithm, based on the previous works of Brandes [1, 2] and Eppstein and Wang [3], to approximate centralities by sampling the contributions of a subset of network nodes to their values. This second algorithm, referred to as Multi-BFS, has a significantly smaller communication overhead.

This document is organized as follows: in section 2 some preliminary definitions are given, while Deccen is defined in section 3 and Multi-BFS is defined in section 4. Section 5 reports some experimental results: the aim of the experiments was to compare the performance of the two algorithms, and to evaluate the quality of the approximations obtained with Multi-BFS. Finally, section 6 contains a small overview of the most relevant choices made in the development of the simulator.

2 Preliminary definitions and assumptions

The task is to compute centrality indices for a given undirected graph G = (V, E), which is assumed to be connected. Unless otherwise stated, n denotes the number of nodes |V| and m the number of edges |E|. Each vertex $v \in V$ represents a network node with some given computational power, that can only communicate with its direct neighbors $N_v = \{u \in V : \{u, v\} \in E\}$. The terms *vertex* and *node* will be used interchangeably.

A path of length k from a source $s \in V$ to a destination $t \in V$ is a sequence $\langle v_0, v_1, \ldots, v_k \rangle$ of vertices such that $s = v_0$, $t = v_k$ and $\{v_{i-1}, v_i\} \in E$ for $i = 1, 2, \ldots, k$. The distance d(u, v) between two vertices is the length of the shortest path that connects them (note that d(u, v) = d(v, u) since the graph is undirected) and the diameter Δ is the maximum distance between any pair of vertices. A vertex v is a predecessor of w with respect to s if $\{v, w\} \in E$ and d(s, v) + 1 = d(s, w). The predecessor set $P_s(w)$ of a vertex w is the set of all predecessors of w with respect to s.

The number of different shortest paths that connect two vertices $s, t \in V$ is denoted by σ_{st} , while the quantity $\sigma_{st}(v)$ is the number of shortest paths between s and t that pass through v (this means that if v = s or v = t then $\sigma_{st}(v)$ is always zero).

The centrality indices relevant to this document are the following:

Closeness centrality. The *closeness* centrality $C_C(v)$ of a vertex $v \in V$ is

$$C_C(v) = \frac{\sum_{u \in V} d(v, u)}{n - 1} \tag{1}$$

Stress centrality. The *stress* centrality $S_C(v)$ of a vertex $v \in V$ is

$$S_C(\nu) = \sum_{s \in V} \sum_{t \in V} \sigma_{st}(\nu) \tag{2}$$

Betweenness centrality. The *betweenness* centrality $B_C(v)$ of a vertex $v \in V$ is

$$B_C(v) = \sum_{s \in V} \sum_{t \in V} \frac{\sigma_{st}(v)}{\sigma_{st}}$$
 (3)

The algorithms described in this report all assume an underlying synchronous communication model where the computation evolves in steps: at each step all the network nodes perform their computations independently and autonomously, and the messages they send at step t are delivered to the destinations and processed at step t+1.

3 The Deccen algorithm

The specification of DECCEN is based on the scheme outlined in [8], which uses broadcasts to count the number of shortest paths and distances between any pair of nodes by leveraging the synchronous model and the following

Lemma (Bellman conditions). A node $v \in V$ lies on a shortest path from $s \in V$ to $t \in V$ if and only if d(s,t) = d(v,s) + d(v,t).

After k steps any node $t \in V$ will know all the nodes $s \in V$ such that d(s, t) = k and the number of shortest paths σ_{st} that links it to each of them (taken as

the sum of the number of shortest paths between s and all the predecessors of t with respect to s). This information is stored locally and also reported in broadcast to allow other nodes $v \in V$ to compute the quantity $\sigma_{st}(v)$: the synchronous communication ensures that if a node v lies on a shortest path between s and t and receives a report for such a pair, it has already computed σ_{vs} and σ_{vt} ; then, according to the above conditions $\sigma_{st}(v) = \sigma_{vs} \cdot \sigma_{vt}$.

3.1 Message types

The messages used during the algorithm execution are the following:

DISCOVERY $\langle s, u, d, \sigma_{su} \rangle$ These messages are broadcast from a node to allow others to compute their distance from it and the number of shortest paths. They contain the source $s \in V$ of the broadcast, the sender $u \in V$, the distance d = d(s, u) of u from s and the number of shortest path that connect u to s.

REPORT $\langle (s,t), \sigma_{st}, d_{st} \rangle$ These messages are broadcast by a node t after it has determined the distance from and the number of shortest path to s.

3.2 Node state

Each node v maintains three accumulators C_C , C_B and C_S to compute its closeness, betweenness and stress centrality values, a counter k, a set R of node pairs for which a REPORT message has already been received, and two dictionaries D and S that associate each node $s \in V$ with the discovered distance d(s, v) and the number of shortest paths σ_{SV} respectively.

3.3 Algorithm initialization

Each node $v \in V$ initializes the counter k and the accumulators C_C , C_B and C_S to 0, the set R to the empty set, the dictionary D so that it only contains the entry (v,0) and S so that it only contains the entry (v,1). Furthermore, it sends to all its neighbors a DISCOVERY(v,v,0,1) message.

3.4 Step actions

The actions performed by a node v at each step are the following:

- 1. All the DISCOVERY messages are grouped by source. Any group of messages having a source s appearing in the dictionary D as key is discarded.
- 2. Each group of messages is processed independently. For each group, let s be the source and d be the distance of all the DISCOVERY $\langle s, u, d, \sigma_{su} \rangle$ messages in it (they will be the same for all the messages), then:
 - 2.1. Add the entry (s, d+1) to the dictionary D, so that D[s] = d+1 and increment the counter k.

- 2.2. Let $\sigma_{sv} = \sum_{u} \sigma_{su}$ and add the entry (s, σ_{sv}) to S. (The number of shortest paths from s to v is the sum of the number of shortest paths from s to all the predecessors of v).
- 2.3. Send a REPORT $\langle (s, v), \sigma_{sv}, d+1 \rangle$ message to every neighbor.
- 3. For each REPORT $\langle (s,t), \sigma_{st}, d_{st} \rangle$ message such that $(s,t) \notin R$:
 - 3.1. If s = v then $C_C \leftarrow C_C + d_{st}$.
 - 3.2. If $d_{st} = D[s] + D[t]$ let $\sigma_{st}(v) = \sigma_{sv} \cdot \sigma_{tv}$, then $C_B \leftarrow C_B + \frac{\sigma_{st}(v)}{\sigma_{st}}$ and $C_S \leftarrow C_S + \sigma_{st}(v)$.
 - 3.3. Add the pair (s, t) to R, and forward the REPORT to every neighbor.

At the end of the execution a node v can obtain its closeness centrality index as $\frac{C_C}{k}$, its betweenness centrality as C_B and its stress centrality as C_S .

3.5 Cost analysis

The broadcast of a DISCOVERY or a REPORT requires O(m) messages. Since each node starts a DISCOVERY and generates n-1 reports the total number of messages exchanged is $O(nm+n^2m)$, with REPORT messages inducing the dominant term n^2m . An optimization that can be performed at step 3.3 is to avoid the propagation of a (s,t) REPORT if $\sigma_{st}(\nu)=0$. In this case the report is irrelevant to ν , and any neighbor that may have needed it will have received it earlier from the broadcast of other nodes that lie on shortest st-paths.

Regarding memory consumption, each node will add O(n) entries each of the two dictionaries and $O(n^2)$ pairs to the set R.

4 Approximation of centrality indices

The high cost of Deccen justifies the investigation of approximation techniques in order to reduce communication and memory overhead at the expense of numerical accuracy.

Eppstein and Wang introduced in [3] an approximation algorithm for the computation of closeness centrality. Their approach is to sample the contributions of a subset of nodes to any closeness value, and extrapolate from those an approximated measure; the contribution of a node s to the closeness of v is taken to be the distance d(v,s) and is computed by solving a Single-Source-Shortest-Path problem with s as source. Brandes in [2] extended this scheme to the approximation of betweenness centrality: his method relies on reformulating the betweenness in terms of contributions from other nodes and employs an augmented SSSP to compute those contributions recursively, as he originally proposed in [1]. This approach can be applied in a decentralized way (by adapting the SSSP problem, which in th undirected case becomes a decentralized Breadth-First Search) and extended to the estimation of stress centrality,

leading to the definition of another decentralized algorithm for the computation (and estimation) of centrality indices.

4.1 Reformulating centrality indices

The first step is to rewrite each centrality index as a sum of terms where each term denotes the contribution of a different *source* node.

Contribution of a source to Closeness centrality As already mentioned, the contribution of a source node s to the closeness centrality of v is the distance d(s, v). The closeness of a node is simply the sum of the contributions of all the other nodes to its centrality divided by n-1.

Contribution of a source to Betweenness centrality The contribution of a source s to the betweenness centrality of v is the *dependency* of s on v introduced in [1]:

$$\delta(s|v) = \sum_{t \in V} \frac{\sigma_{st}(v)}{\sigma_{st}},\tag{4}$$

that allows to rewrite the betweenness centrality of ν as

$$BC(v) = \sum_{s \in V} \delta(s|v). \tag{5}$$

Contribution of a source to Stress centrality The contribution to stress centrality is analogous to the contribution to betweenness centrality:

$$\sigma(s|\nu) = \sum_{t \in V} \sigma_{st}(\nu), \tag{6}$$

and the stress centrality of v is rewritten as

$$SC(v) = \sum_{s \in V} \sigma(s|v). \tag{7}$$

4.2 Computing contributions from a single source

A decentralized Breadth First Search can be adapted to compute the contribution of a source to any of the three centrality indices considered.

The closeness centrality contribution is simply the distance from the source to the node and it can be computed directly during the visit.

For betweenness centrality, Brandes [1] has shown that dependencies of a source obey a recursive relation expressed in terms of predecessors set:

Theorem 1 (Brandes, 2001). The dependency of $s \in V$ on any $v \in V$ obeys

$$\delta(s|\nu) = \sum_{w:\nu \in P_s(w)} \frac{\sigma_{s\nu}}{\sigma_{sw}} \cdot (1 + \delta(s|w)). \tag{8}$$

For stress centrality contributions a similar relation holds (the proof is reported in the appendix):

Theorem 2. The stress centrality contribution of $s \in V$ on any $v \in V$ obeys

$$\sigma(s|\nu) = \sum_{w:\nu \in P_s(w)} \sigma_{s\nu} \cdot \left(1 + \frac{\sigma(s|w)}{\sigma_{sw}}\right). \tag{9}$$

The predecessor sets can be easily discovered by the BFS algorithm during the network exploration, while contributions are computed with a backward walk from the frontier of the BF-Tree to the source of the visit.

4.3 Random sampling of source nodes

To let the algorithm operate in a decentralized way, each node independently initiates a visit with a given probability p, which reflects the fraction of the network sampled. Even if the number k of samples is not known beforehand, all the nodes will be able to compute it by counting the number of visits in which they will take part.

The result of sampling from a source node v_i at a node u is modeled with the following random variables (that rely on the contributions of v_i to u)

$$X_i(u) = \frac{n}{n-1} \cdot d(v_i, u), \qquad Y_i(u) = n \cdot \delta(v_i|u), \qquad Z_i(u) = n \cdot \sigma(v_i|u),$$

which are then used to define the estimators of the three centrality indices as

$$\widetilde{C}_C(u) = \sum_{i=1}^k \frac{X_i(u)}{k}, \qquad \widetilde{C}_B(u) = \sum_{i=1}^k \frac{Y_i(u)}{k}, \qquad \widetilde{C}_S(u) = \sum_{i=1}^k \frac{Z_i(u)}{k}.$$

The proof of the next theorem – which ensures that the centrality estimators are unbiased – is reported in the appendix.

Theorem 3 (Unbiased estimators). *The expected values of the centrality estimators are the actual centrality indices:*

- (a) $\mathbf{E}[\widetilde{C}_C(u)] = C_C(u)$,
- (b) $\mathbf{E}[\widetilde{C}_S(u)] = C_S(u)$,
- (c) $\mathbf{E}[\widetilde{C}_B(u)] = C_B(u)$.

4.4 Multi-BFS algorithm specification

The only parameter of the algorithm is the probability p of a node becoming the source of a decentralized BFS and having its centrality contributions sampled at every location. The network size n is assumed to be known in advance to all the nodes, otherwise it can be easily computed during the backtracking phase of the augmented BFS and broadcast by one or more sources.

Note that if this algorithm is executed with p = 1 it is basically the decentralized version of the algorithm described in [1], and at the end of the execution the estimators will yield the exact centrality values.

4.4.1 Message types

The messages used during the algorithm execution are the following:

DISCOVERY $\langle s, u, d, \sigma_{su} \rangle$ These messages are used during the visit from a source to build the predecessors sets at each node. The fields are the source s of the visit, the sender u, the distance d of the sender to the source (that is, d = d(s, u)) and the number of shortest path from the source to the sender σ_{su} .

REPORT $\langle s, v, \delta(s|v), \sigma(s|v), \sigma_{sv} \rangle$ These messages are sent by a node v as part of the backward walk to inform its predecessors of the computed contributions and to allow them to compute their own by applying the recursive relations introduced in section 4.2.

4.4.2 Visit states

Visit states are parametric with respect to the discovery from a source $s \in V$. A node v is in state:

WAITING(s) if it has not yet received any DISCOVERY having s as source.

ACTIVE(s) if it has received one or more DISCOVERY messages with source s and has not yet computed the contributions of s to its centrality indices.

COMPLETED(s) if it has computed the contributions of s to its centrality indices and reported them back to each predecessor in $P_s(v)$.

4.4.3 Node state

Each node v maintains three centrality accumulators C_C , C_B and C_S like in Deccen, and a counter k to track the number of sample nodes involved in the procedure. Furthermore, while a node is in state ACTIVE(s) it will need to partition the set of neighbors N_v in three subsets: the set of predecessors $P_s(v)$, the set of siblings $S_s(v)$ and the set of children $C_s(v)$, and to track the contributions of s to its centrality values with three parametric accumulators $C_C^{(s)}$, $C_B^{(s)}$ and $C_S^{(s)}$.

4.4.4 Protocol initialization

4.4.5 Step actions

The actions performed by a node v at each step are:

- 1. The messages received at the current step are divided by type and then DISCOVERY messages are grouped by source.
- 2. For each group of DISCOVERY $\langle s, u, d, \sigma_{su} \rangle$ messages having s as a common source (at distance d), with v in state WAITING(s):
 - 2.1. Change state to ACTIVE(s) and initialize $C_C^{(s)} \leftarrow d+1$, $C_B^{(s)} \leftarrow 0$ and $C_S^{(s)} \leftarrow 0$; then collect all the senders u in the predecessor set $P_s(v)$ and compute $\sigma_{sv} = \sum_{u \in P_s(v)} \sigma_{su}$.
 - 2.2. If $P_s(v) = N_v$ send a REPORT $\langle s, v, 0, 0, \sigma_{sv} \rangle$ message to all the predecessors $u \in P_s(v)$ and change state to COMPLETED(s).
 - 2.3. Otherwise, send to all $w \in N_{\nu} \setminus P_{s}(\nu)$ a DISCOVERY $\langle s, \nu, d+1, \sigma_{s\nu} \rangle$ message and change state to ACTIVE(s).
- 3. For each group of DISCOVERY $\langle s, u, d, \sigma_{su} \rangle$ messages having source s and distance d with v in state ACTIVE(s):
 - 3.1. Under the synchronous model assumption v can only receive such messages from nodes u such that d(s,u) = d(s,v). Collect these nodes in the set $S_s(v)$ of the siblings of v with respect to s. (Note that these messages are received exactly one step after v has been contacted by its predecessors).
- 4. For each REPORT $\langle s, w, \delta(s|w), \sigma(s|w), \sigma_{sw} \rangle$ message received when v is in state ACTIVE(s):
 - 4.1. Add w to the children set $C_s(v)$.
 - 4.2. Update $C_B^{(s)} \leftarrow C_B^{(s)} + \frac{\sigma_{sv}}{\sigma_{sw}} \cdot (1 + \delta(s|w))$.
 - 4.3. Update $C_S^{(s)} \leftarrow C_S^{(s)} + \sigma_{sw} \cdot (1 + \frac{\sigma(s|w)}{\sigma_{sw}})$.
- 5. For any $s \in V$ with v is in state ACTIVE(s), if $P_s(v) \cup S_s(v) \cup C_s(v) = N_v$ then:
 - 5.1. If $s \neq v$ update the accumulators:
 - $C_C \leftarrow C_C + C_C^{(s)}$
 - $-C_B \leftarrow C_B + C_B^{(s)}$
 - $C_S \leftarrow C_S + C_S^{(s)}$

and send a REPORT $\langle s, v, \delta(s|v), \sigma(s|v), \sigma_{sv} \rangle$ message to all the predecessors contained in $P_s(v)$.

5.2. Increment the counter k and change state to COMPLETED(s).

A node ν can obtain the value of the estimators in the following way:

$$\widetilde{C}_C(v) = \frac{nC_C}{k(n-1)}, \qquad \widetilde{C}_B(v) = \frac{nC_B}{k}, \qquad \widetilde{C}_S(v) = \frac{nC_S}{k}.$$

4.5 Cost analysis

Each independent visit requires O(m) messages for the DISCOVERY phase and O(m) messages to backtrack with REPORT messages. The parameter p determines the fraction of nodes pn that initiate a DISCOVERY search, so the total number of messages is $O(\lceil pn \rceil m)$.

In terms of memory consumption, note that for each $v \in V$ there are at most $\lceil pn \rceil$ other nodes s for which v is in state ACTIVE(s) and needs to partition its neighbor set N_v in the three subsets $P_s(v)$, $S_s(v)$ and $C_s(v)$, so the cost is $O(\lceil pn \rceil N_v)$.

5 Experiments

The aim of the experiments was to compare the number of messages required by the algorithms to complete the decentralized computation of the indices, and to measure the error of the estimates of MULTI-BFS with different values of the parameter *p*.

The results were obtained by running simulations on the following networks (taken from the KONECT [7] repository):

dolphins A social network of bottlenose dolphins where nodes are dolphins and edges represent frequent associations between individuals. (n = 62, m = 159, source [9]).

surf A network that represents interpersonal contacts between windsurfers in southern California during the fall of 1986. (n = 62, m = 336, source [4]).

macaques The undirected version of a directed network that represent dominance behavior in a colony of 62 adult female Japanese macaques – obtained by forcing the edges to be symmetric. (n = 62, m = 1167, original source [11]).

train A network that contains contacts between suspected terrorists involved in the train bombing of Madrid on March 11, 2004 as reconstructed from newspapers. (n = 62, m = 243, source [6]).

email The email communication network at the University Rovira i Virgili in Tarragona in the south of Catalonia in Spain. (n = 1133, m = 5451, source [5]).

powergrid The high-voltage power grid in the Western States of the United States of America. (n = 4941, m = 6594, source [12]).

				Deccen		Multi-BFS	
Network	n	m	Δ	Steps	Messages	Steps	Messages
surf	43	336	3	6	181374	7	28896
dolphins	62	159	8	16	140051	17	19716
macaques	62	1167	2	4	1458368	5	144708
train	64	243	6	12	272673	13	31104

Table 1: Steps required to complete and number of exchanged messages by Deccen and Multi-BFS executed with p = 1.

5.1 Performance evaluation

The performance measures used to evaluate the algorithms are the number of steps required to complete the computation and the number of messages that the network nodes generate.

Table 1 reports the results obtained by running Deccen and Multi-BFS with p=1 (in order to compute exact values). Due to the high memory requirements of Deccen the simulations could only be performed on small networks, but the algorithms can be expected to behave similarly in any other scenario since all the messages are generated in a deterministic way.

As expected, both algorithms require a number of steps to complete the execution that is twice the diameter Δ of the network, since each of the two "discovery" phases takes at most Δ steps to reach any destination from any given source, and the same also holds for the report phase of Deccen or the backtracking in Multi-BFS. In the test instances Multi-BFS exhibits an extra delay of one step between reaching the farthest child nodes and the beginning of the backtracking phase: this is bound to happen whenever nodes at distance Δ from a source have a non-empty siblings set, as the siblings need to be identified before reporting back.

The experimental results also show that DECCEN requires a significantly larger amount of messages than MULTI-BFS to terminate the execution, as anticipated in the previous sections. This is a consequence of the different way in which the algorithms handle the report phase: while in DECCEN new reports are broadcast at every step of a discovery, MULTI-BFS uses the distance from the source of the discovery to rank the nodes, and only generates reports from the most distant nodes (that are then routed back to the source using the predecessor sets).

5.2 Multi-BFS approximations

The quality of the approximations obtained with Multi-BFS were evaluated by running simulations with increasing values of p (from 0.05 to 1.0 at steps of 0.05). The goal was to measure the numerical error of the estimates generated

by the algorithm, and to evaluate how close the "ranking" of the nodes using these estimates is to the actual ranking of the nodes according to the exact centrality values, since nodes could be ranked correctly even if the estimates are subject to high error (this is especially interesting if a node wants to make a "greedy" choice knowing only its and it's neighbors index's value, as it gives an idea of how often this choice may be the wrong one). Each experiment was repeated ten times.

Figure 1 reports the average relative error ϵ_r yielded by the centrality estimators in the dolphins, email and powergrid networks. The data shows satisfactory results for the estimation of closeness centrality for small values of p, while stress and betweenness estimates exhibit large numerical error even for values of p close to 1.

The accuracy of the ranking induced by the approximated indices is measured by counting, among all the pairs of nodes, the fraction of pairs in which the nodes are wrongly ordered with respect to the ranking imposed by the exact centrality values; results are reported in figure 2. In this case the results are encouraging: even for small values of *p* the fraction of pairs in wrong order is relatively small.

6 Project code overview

The project code is based on the *event driven* engine of the PeerSim [10] simulator.

As already stated stated, the algorithms presented in this report assume the communication is synchronized, which means that a computation step is not executed until all the messages sent at the previous step have reached the destination. This can be achieved with the event driven engine by letting the nodes exchange messages using a reliable Transport protocol and having them execute a periodic action (the computation step) at intervals larger than the maximum delay required by the Transport to deliver a message. Since messages are only sent during computation steps, this configuration guarantees that all the messages sent at one step are delivered before the next one is executed.

6.1 Messages

The nodes communicate by exchanging objects of the Message class. This class provides factory methods to generate Message objects that comply with the specification given in the algorithms description.

6.2 Protocol classes

The implementation of the PeerSim protocols for DECCEN and MULTI-BFS is derived from the common base class CentralityProtocol. This class implements

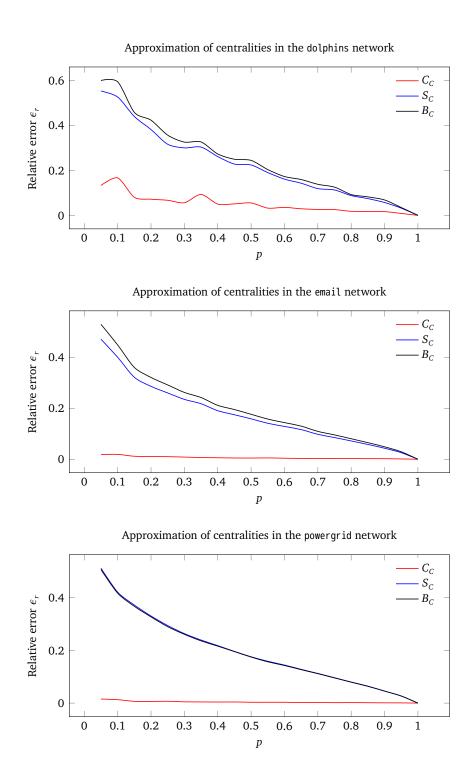


Figure 1: Approximation error in the estimation of centrality indices.

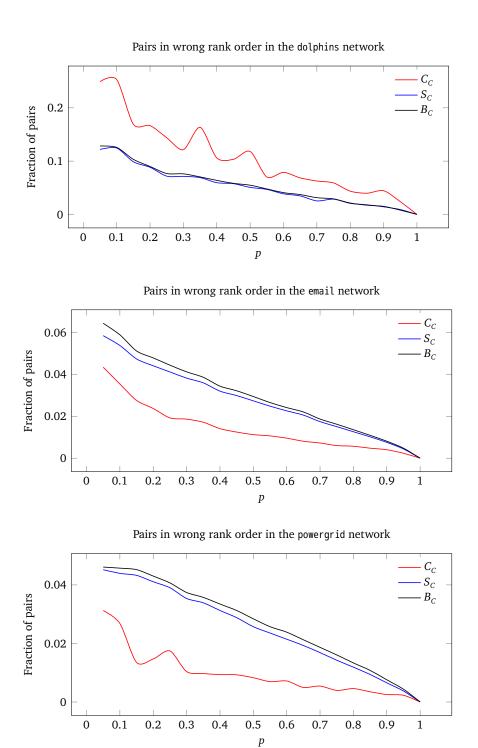


Figure 2: Pairs in wrong rank order with approximated centralities

the common mechanisms needed by the synchronized protocols, in particular the support for the execution of periodic actions and the handling of the "communication windows" during which the network nodes exchange messages. Here is an excerpt of the class:

The implementation of the EDProtocol interface is required by the event driven engine. The EDProtocol.processEvent method simply stores a newly received message in the incoming list, and notifies the CentralitySimulation control that a new message has been received (this control stops the simulation if no messages are sent in a time window):

```
public void processEvent(Node node, int pid, Object event) {
  incoming.add((Message) event);
  CentralitySimulation.newMessage();
}
```

Messages are delivered by a Transport protocol, and can be sent by calling the send method which is a convenience wrapper for Transport.send:

```
protected void send(Node from, Node to, Message m, int pid) {
   Transport tr = (Transport) from.getProtocol(transportProtocolID);
   tr.send(from, to, m, pid);
}
```

The class also implements the CDProtocol interface to support the execution of periodic actions. This feature can be enabled by including in the simulation (through the configuration file) a CDScheduler component and configuring it to run a CDProtocol at regular steps, which causes the simulator to trigger the CDProtocol.nextCycle method on all the nodes that are running it.

When performing the experiments, the synchronous communication was configured using a UniformRandomTransport protocol to deliver messages with

mindelay and maxdelay options of 15 and 200 milliseconds respectively, while the computation steps (implemented by the subclasses in the nextCycle method) were triggered by a CDScheduler at intervals of 500 milliseconds.

6.2.1 Deccen protocol implementation

The Deccen class implements a DECCEN node in the simulator. A node state consists of the shortest path information it collects and of the reports it has handled during the execution:

```
public class Deccen extends SynchronousCentralityProtocol {
   private static class ShortestPathData {
     public final int count;
     public final int length;
     public ShortestPathData(int count, int length) { ... }
   }

   private static class OrderedPair<T1,T2> {
     public final T1 first;
     public final T2 second;
     public OrderedPair(T1 first, T2 second) { ... }
        ...
   }
   ...
   private Map<Node,ShortestPathData> shortestPathMap;
   private Set<OrderedPair<Long,Long>> handledReports;
   ...
   public void nextCycle(Node self, int protocolID) { ... }
   ...
}
```

Whenever new DISCOVERY messages are received, additional information about the number and length of shortest paths toward a particular source becomes available: this information is stored in the shortestPathMap structure by adding a new ShortestPathData entry, paired with the source of the discovery. As the algorithm computes all the centrality indices, it needs both the distance from the source and the number of different paths counted (the length and count fields respectively). Reports about a source—destination pair (s,t) are tracked by adding to the handledReports set an OrderedPair of Node IDs.

The nextCycle method implements the actions specified by this report in section 3.4:

```
public void nextCycle(Node self, int protocolID) {
   Map<Node,List<Message>> discoveryMap =
        new HashMap<Node,List<Message>>();
```

```
List<Message> reportList = new LinkedList<Message>();
parseIncomingMessages(discoveryMap, reportList);
if (!discoveryMap.isEmpty())
   processDiscoveryMessages(self, protocolID, discoveryMap);
if (!reportList.isEmpty())
   processReportMessages(self, protocolID, reportList);
}
```

Messages received in the previous communication window are parsed by the parseIncomingMessages method: discovery messages with the same source are grouped together in the discoveryMap structure (using the source node as key), and reports are collected in the reportList.

6.2.2 Multi-BFS protocol implementation

The state of a Multi-BFS node consists of a Map to keep track of the various visits that are performed by the source nodes during the algorithm execution:

```
public class MultiBFS extends SynchronousCentralityProtocol {
   private static class VisitState {
     ...
   }
   ...
   private Map<Node, VisitState> activeVisits;
   private Set<Node> completed;
   ...
   public void nextCycle(Node self, int protocolID) { ... }
   ...
   public boolean isWaiting(Node source) { ... }
   public boolean isActive(Node source) { ... }
   public boolean isCompleted(Node source) { ... }
}
```

Recall that the state of a node is parametric with respect to each source of a visit. The state is WAITING(s) if the Node instance s does not appear as key in the activeVisits map and neither is contained in the completed set; this is the initial state for any source since both structures are empty at the start of the protocol. When a node is in state ACTIVE(s) an entry with key s is present in the activeVisits map. After a node has reported back to all the predecessors, it changes state COMPLETED(s) by removing the mapping from activeVisits and inserting s in the completed set.

Entries in the activeVisits map are used to store data while a node is in the "active" phase. A VisitState object is used to keep track of the predecessors and children sets, and to incrementally compute the values to be reported to the predecessor nodes:

```
private static class VisitState {
  public Set<Node> predecessors;
  public Set<Node> siblings;
  public Set<Node> children;
  public int distanceFromSource;
  public int sigma;
  public long contributionSC;
  public double contributionBC;
  ...
  public void accumulate(
     Node child, double bcc, long scc, int numSP) { ... }
}
```

The accumulate method updates the local contributions of the source (stored in the contributionSC and contributionBC fields) by applying the recursive relations (8) and (9), and it's invoked whenever a child node sends a report.

Finally, the nextCycle method which again follows the specification given earlier (in section 4.4.5):

```
public void nextCycle(Node self, int protocolID) {
   Map<Node,List<Message>> discoveryMap =
        new HashMap<Node,List<Message>>();
   List<Message> reportList = new LinkedList<Message>();
   parseIncomingMessages(discoveryMap, reportList);
   if (!discoveryMap.isEmpty())
      processDiscoveryMessages(self, protocolID, discoveryMap);
   if (!reportList.isEmpty())
      processReportMessages(reportList);
   reportNewCompleted(self, protocolID);
}
```

Active visits for which all child nodes have reported are finalized with the reportNewCompleted method by integrating the contributions in the centrality accumulators and sending reports to the predecessors.

Appendix: Theorem proofs

Theorem 2. The stress centrality contribution of $s \in V$ on any $v \in V$ obeys

$$\sigma(s|\nu) = \sum_{w:\nu \in P_s(w)} \sigma_{s\nu} \cdot \left(1 + \frac{\sigma(s|w)}{\sigma_{sw}}\right). \tag{9}$$

Proof. The proof is similar to the one used by Brandes [1] to prove Theorem 1. Let $\sigma_{st}(v, e)$ be the number of shortest paths between s and t that pass through v and across edge e. Observe that if a shortest path from s to t passes through

v, then after v it must immediately reach some other node w that has v in its predecessor set $P_s(w)$, so equation (6) can be rewritten as

$$\sigma(s|v) = \sum_{t \in V} \sigma_{st}(v) = \sum_{t \in V} \sum_{w:v \in P_s(w)} \sigma_{st}(v, \{v, w\})$$
$$= \sum_{w:v \in P_s(w)} \sum_{t \in V} \sigma_{st}(v, \{v, w\}).$$

Let w be any node with $v \in P_s(w)$, then

$$\sigma_{st}(v, \{v, w\}) = \begin{cases} \sigma_{sv} & \text{if } t = w \\ \frac{\sigma_{sv}}{\sigma_{sw}} \cdot \sigma_{st}(w) & \text{if } t \neq w \end{cases}$$

and substituting it in the previous expression yields

$$\sigma(s|v) = \sum_{w:v \in P_s(w)} \sum_{t \in V} \sigma_{st}(v, \{v, w\})$$

$$= \sum_{w:v \in P_s(w)} \left(\sigma_{sv} + \sum_{t \neq w} \frac{\sigma_{sv}}{\sigma_{sw}} \cdot \sigma_{st}(w)\right)$$

$$= \sum_{w:v \in P_s(w)} \sigma_{sv} \cdot \left(1 + \frac{\sigma(s|w)}{\sigma_{sw}}\right).$$

Theorem 3 (Unbiased estimators). *The expected values of the centrality estimators are the actual centrality indices:*

- (a) $\mathbf{E}[\widetilde{C}_C(u)] = C_C(u)$,
- (b) $\mathbf{E}[\widetilde{C}_S(u)] = C_S(u)$,
- (c) $\mathbf{E}[\widetilde{C}_B(u)] = C_B(u)$.

Proof. (a) Recall that for the estimation of closeness centrality, the result of sampling from a source v_i at a node u is modeled with the random variable $X_i(u) = \frac{n}{n-1} \cdot d(v_i, u)$. The derivation exploits the linearity of the expected value operator **E** and the fact that source nodes are random (that is, each node has equal probability 1/n of being a source).

$$\mathbf{E}[\widetilde{C}_{C}(u)] = \mathbf{E}\left[\sum_{i=1}^{k} \frac{X_{i}(u)}{k}\right] = \mathbf{E}\left[\sum_{i=1}^{k} \frac{n \cdot d(v_{i}, u)}{k(n-1)}\right]$$

$$= \frac{n}{k(n-1)} \sum_{i=1}^{k} \mathbf{E}[d(v_{i}, u)] \qquad \text{(by linearity of expectation)}$$

$$= \frac{n}{k(n-1)} \cdot k \cdot \frac{1}{n} \sum_{v \in V} d(v, u) \qquad \text{(random source selection)}$$

$$= \frac{\sum_{v \in V} d(v, u)}{n-1} = C_{C}(u)$$

(b) To estimate stress centrality, the result of sampling from v_i is modeled at u with the random variable $Y_i(u) = n \cdot \sigma(v_i|u)$. The derivation relies again on the linearity of **E** and the uniform distribution of source nodes.

$$\mathbf{E}[\widetilde{C}_{S}(u)] = \mathbf{E}\left[\sum_{i=1}^{k} \frac{Y_{i}(u)}{k}\right] = \mathbf{E}\left[\sum_{i=1}^{k} \frac{n \cdot \sigma(v_{i}|u)}{k}\right]$$
$$= \frac{n}{k} \sum_{i=1}^{k} \mathbf{E}[\sigma(v_{i}|u)] = \frac{n}{k} \cdot k \cdot \frac{1}{n} \sum_{v \in V} \sigma(v|u) = C_{S}(u)$$

(c) The proof is the same as (b) with the substitution of $Z_i(u) = n \cdot \delta(v_i|u)$ for $Y_i(u)$.

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