Fault-Tolerant Spanners: Better and Simpler*

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ABSTRACT

A natural requirement for many distributed structures is fault-tolerance: after some failures in the underlying network, whatever remains from the structure should still be effective for whatever remains from the network. In this paper we examine spanners of general graphs that are tolerant to vertex failures, and significantly improve their dependence on the number of faults r for all stretch bounds.

For stretch $k \geq 3$ we design a simple transformation that converts every k-spanner construction with at most f(n) edges into an r-fault-tolerant k-spanner construction with at most $O(r^3 \log n) \cdot f(2n/r)$ edges. Applying this to standard greedy spanner constructions gives r-fault tolerant k-spanners with $\tilde{O}(r^2n^{1+\frac{2}{k+1}})$ edges. The previous construction by Chechik, Langberg, Peleg, and Roddity [CLPR09] depends similarly on n but exponentially on r (approximately like k^r).

For the case of k=2 and unit edge-lengths, an $O(r\log n)$ -approximation is known from recent work of Dinitz and Krauthgamer [DK11], in which several spanner results are obtained using a common approach of rounding a natural flow-based linear programming relaxation. Here we use a different (stronger) LP relaxation and improve the approximation ratio to $O(\log n)$, which is, notably, independent of the number of faults r. We further strengthen this bound in terms of the maximum degree by using the Lovász Local Lemma.

Finally, we show that most of our constructions are inherently local by designing equivalent distributed algorithms in the \mathcal{LOCAL} model of distributed computation.

Categories and Subject Descriptors

G.2.2 [Discrete Mathematics]: Graph Theory—Graph Algorithms

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1. INTRODUCTION

Let G=(V,E) be a graph, possibly with edge-lengths $\ell: E \to \mathbb{R}_{\geq 0}$. A k-spanner of G, for $k \geq 1$, is a subgraph G'=(V,E') that preserves all pairwise distances within factor k, i.e. for all $u,v \in V$,

$$d_{G'}(u,v) \le k \cdot d_G(u,v). \tag{1}$$

Here and throughout, d_H denotes the shortest-path distance in a graph H, and n=|V|. The distance preservation factor k is called the stretch of the spanner. It is easy to see that requiring (1) only for edges $\{u,v\} \in E$ suffices. This definition also extends naturally to directed graphs. Obviously G is a 1-spanner of itself, so usually the goal is to compute a "small" spanner. One traditional notion of "small" is the number of edges in G', called the size of G'.

This notion of graph spanners, first introduced by Peleg and Schäffer [PS89] and Peleg and Ullman [PU89], has been studied extensively, with applications ranging from routing in networks (e.g. [AP95, TZ05]) to solving linear systems (e.g. [ST04, EEST08]). Many of these applications, especially in distributed computing, arise by modeling computer networks or distributed systems as graphs. But one aspect of distributed systems that is not captured by the above spanner definition is the possibility of failure. We would like our spanner to be robust to failures, so that even if some nodes fail we still have a spanner of what remains. More formally, G' is an r-fault tolerant k-spanner of G if for every set $F \subseteq V$ with $|F| \le r$, the spanner condition holds for $G \setminus F$, i.e. for all $u, v \in V \setminus F$ we have $d_{G' \setminus F}(u, v) \le k \cdot d_{G \setminus F}(u, v)$.

This notion of fault-tolerant spanners was first introduced by Levcopoulos, Narasimhan, and Smid [LNS98] in the context of geometric spanners (the special case when the vertices are in Euclidean space and the distance between two points is the Euclidean distance). They provided both size and weight bounds for $(1+\epsilon)$ -spanners, which were later improved by Lukovski [Luk99] and Czumaj and Zhao [CZ03]. The first result on fault-tolerant spanners for general graphs, by Chechik, Langberg, Peleg, and Roditty [CLPR09], constructs r-fault tolerant (2k-1)-spanners with size $O(r^2k^{r+1} \cdot n^{1+1/k} \log^{1-1/k} n)$, for any integer $k \geq 1$. Since it has long been known how to construct (2k-1)-spanners with size

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 $O(n^{1+1/k})$ (see e.g [ADD⁺93]), this means that the extra cost of r-fault tolerance is $O(r^2k^{r+1})$. While this is independent of n, it grows rapidly as the number of faults r gets large. We address an important question left open in [CLPR09] of improving this dependence on r from exponential to polynomial.

Nontrivial absolute bounds on the size of a k-spanner are possible only when the stretch $k \geq 3$. For k = 2, there are graphs with $\Omega(n^2)$ edges for which every edge must be included in the spanner (e.g. a complete bipartite graph). So the common approach is to provide relative bounds, namely, design approximation algorithms for the problem of computing a minimum size r-fault tolerant 2-spanner. In this context one typically assumes that all edges have unit length. Without fault tolerance, the problem is reasonably well understood: there are algorithms that provide an $O(\log n)$ -approximation [KP94, EP01] (or, with some extra effort, an $O(\log(|E|/|V|))$ -approximation), and the problem is NP-hard to approximate better than $\Omega(\log n)$ [Kor01]. For the r-fault tolerant 2-spanner problem, the first nontrivial approximation was recently given by Dinitz and Krauthgamer [DK11], who designed an $O(r \log n)$ approximation. However, they did not provide evidence that this loss of r was necessary, an issue that we address in this paper.

1.1 Results and Techniques

Stretch bounds k > 3.

Our main result for general $k \geq 3$ is a new r-fault tolerant k-spanner with size that depends only polynomially on r, thereby improving over the exponential dependence by Chechik et al. [CLPR09].

Theorem 1.1. For every graph G = (V, E) with positive edge-lengths and odd $k \geq 3$, there is an r-fault tolerant k-spanner with size $O(r^{2-\frac{2}{k+1}}n^{1+\frac{2}{k+1}}\log n)$.

In fact, we prove something slightly stronger: a general conversion theorem that turns any algorithm for constructing k-spanners with size f(n) into an algorithm for constructing r-fault tolerant k-spanners with size $O(r^3 \log n \cdot f(2n/r)$. Applying this conversion to the well-known greedy spanner algorithm (see e.g. [ADD⁺93]) immediately yields Theorem 1.1.

At a high level, Chechik et al. [CLPR09] apply the spanner construction of Thorup and Zwick [TZ05] to every possible fault set, eventually taking the union of all of these spanners. They show, through a rather involved analysis that relies on specific properties of the Thorup-Zwick construction, that taking a union over as many as $O(n^r)$ spanners increases the size bound only by an $O(r^2k^r)$ factor. Our conversion technique, on the other hand, is extremely general. Inspired by the color-coding technique of Alon, Yuster, and Zwick [AYZ95] and its recent incarnation in designing data structures and oracles (e.g. [WY10]), we randomly sample nodes to act as a fault set, and then apply a generic spanner algorithm on what remains. Our sampling dramatically oversamples nodes — instead of fault sets of size r, we end up with sampled fault sets of size approximately $(1-\frac{1}{n})n$. This allows us to satisfy many fault sets of size r with a single iteration of the generic algorithm. The size bound follows almost immediately.

Stretch k = 2 (and assuming unit edge-lengths).

Here, our main result is an approximation algorithm with ratio that is independent of r. Our algorithm actually works in an even more general setting, where the graph is directed and edges have arbitrary nonnegative costs $c_e: E \to \mathbb{R}_{\geq 0}$. The goal is to find an r-fault tolerant 2-spanner of minimum total cost. We refer to this problem as MINIMUM COST r-FAULT TOLERANT 2-SPANNER. Since our algorithms are randomized, we will say that they provide an α -approximation if, with probability at least 2/3, they return a valid r-fault 2-spanner of cost at most α times the minimum cost r-fault tolerant 2-spanner. Our algorithms will in fact be a little stronger: they return a valid r-fault tolerant 2-spanner with high probability, and this spanner will be an α -approximation with probability at least 3/4.

Theorem 1.2. There is a (randomized) $O(\log n)$ approximation algorithm for Minimum Cost r-Fault Tolerant 2-Spanner for all $r \leq n$ (even on directed graphs).

This removes the dependence on r from the previously known $O(r \log n)$ -approximation [DK11]. Similarly to this previous approximation [DK11], we design a flow-based linear programming (LP) relaxation of the problem and then apply a rounding scheme that uses randomization at the vertices, rather than naively at the edges. However, the relaxation used by [DK11] is not strong enough to achieve approximation factor independent of r; even simple graphs (such as the complete graph with unit costs) have integrality gaps of $\Omega(r)$. We thus design a different relaxation, and add to it a large family of constraints that are essentially the knapsack-cover inequalities of Carr, Fleischer, Leung, and Phillips [CFLP00], adapted to our context. With these additional constraints, we are able to show that the simple rounding scheme devised in [DK11] now achieves an $O(\log n)$ -approximation.

We further show that in the special case where all edge costs are unit the integrality gap is at most $O(\log \Delta)$, where Δ is the maximum degree of the graph. Note that this bound is at least as good as the $O(\log n)$ bound (and possibly better). We prove this by a more careful analysis of essentially the same randomized rounding scheme using the Lovász Local Lemma. By using constructive versions of the LLL [MT10] we can achieve this gap algorithmically, giving a better approximation algorithm.

Distributed versions of our algorithms.

Finally, one feature that is shared by both the k=2 and the $k\geq 3$ case is that the algorithms are local (assuming that the generic algorithm used by the conversion theorem is itself local). To show this formally, we provide distributed versions of the algorithm in the \mathcal{LOCAL} model of distributed computation. The \mathcal{LOCAL} model is a standard message-passing model in which in each round, every node is allowed to send an unbounded-size message to each of its neighbors [Pel00]. While the unbounded message-size assumption may not be realistic, this model captures locality in the sense that in t rounds, each node has knowledge of, and is influenced by, only the nodes that are within hop-distance t of it.

Assuming that the underlying generic spanner algorithm is distributed in this sense, our general conversion theorem trivially provides a distributed algorithm since the failure sampling is done independently by every node. Designing

a distributed version of the r-fault tolerant 2-spanner algorithm is not quite as simple, since our centralized algorithm uses the Ellipsoid method to solve a linear program that has an exponential number of constraints. While there is a significant amount of literature on solving linear programs in a distributed manner, much of the time strong assumptions are made about the structure of the linear program. In particular, it is common to assume that the LP is a positive (i.e. a packing/covering) LP. Unfortunately the LP relaxation that we use is not positive, even for r=0, so we cannot simply use an off-the-shelf distributed LP solver. Instead, we leverage the fact that the LP itself is "mostly" local we partition the graph into clusters, solve the LP separately on each cluster, and then repeat this process several times, eventually taking the average values. This technique is quite similar to the work of Kuhn, Moscibroda, and Wattenhofer [KMW06], who showed how to approximately solve positive LPs using the graph decompositions of Linial and Saks [LS93]. We construct padded decompositions using a variant of the methods developed by Bartal [Bar96] and by Linial and Saks [LS93]. Combining this distributed methodology for solving the LP relaxation together with the obvious distributed implementation of the aforementioned rounding scheme, we obtain a distributed $O(\log n)$ -approximation. While the approximation works on directed graphs, we need to assume that communication along edges is bidirectional.

THEOREM 1.3. There is a distributed randomized algorithm in the \mathcal{LOCAL} model that takes $O(\log^2 n)$ rounds and is an $O(\log n)$ -approximation for MINIMUM COST r-FAULT TOLERANT 2-SPANNER (even on directed graphs).

2. GENERAL STRETCH k

In this section we give our construction of r-vertex-tolerant k-spanners (with arbitrary edge-lengths). For each $F\subseteq V$ with $|F|\leq r$, we let E_F denote the edges of $G\setminus F$, i.e. $E_F=\{\{u,v\}\in E: u,v\not\in F\}$. We first give a general conversion theorem that turns any k-spanner construction into an r-fault tolerant k-spanner construction at an extra cost of at most $poly(r)\cdot \log n$. This conversion actually works fine even when the underlying spanner construction is randomized, but since good deterministic constructions exist we will assume for simplicity that the underlying construction is deterministic. We say that an event happens with high probability if it happens with probability at least $1-\frac{1}{n^C}$ for constant C that can be made arbitrarily large (at the cost of increasing the constants hidden by $O(\cdot)$ notation).

THEOREM 2.1. If there is an algorithm A that on every graph builds a k-spanner of size f(n), then there is an algorithm that on any graph builds with high probability an r-fault tolerant k-spanner of size $O(r^3 \log n \cdot f(\frac{2\pi}{n}))$.

PROOF. Our algorithm is simple: in each iteration, we independently add each vertex to a set J with probability p=1-1/r, and then use the given algorithm A to build a k-spanner on the remaining graph $G\setminus J$. If r=1 then we can set p=1/2, which will just affect the constants in the $O(\cdot)$. We do this for $\alpha=\Theta(r^3\log n)$ iterations, each independent of the others. Let H be the graph obtained by taking the union of the iterations.

We first bound the size of H. Without loss of generality we can assume that $r \leq n^{2/3}$, since when $r > n^{2/3}$ the claimed size bound is larger than $r^3 \geq n^2$, and it is trivially

true that H has at most $O(n^2)$ edges since H is a graph on n vertices. In each iteration, the expected number of vertices in $G \setminus J$ is n/r. By a simple Chernoff bound, the probability that a given iteration has more than 2n/r vertices in $G \setminus J$ is at most $e^{-(1/3)n/r} \leq e^{-(1/3)n^{1/3}}$. Since there are only $\alpha = O(r^3 \log n)$ iterations, we can take a union bound over the iterations and get that with high probability the number of vertices in $G \setminus J$ is at most 2n/r in every iteration. Thus the total size of H is at most $O(\alpha \cdot f(\frac{2n}{r}))$. Now we just need to prove that this algorithm results in a valid r-fault tolerant k-spanner for $\alpha = O(r^3 \log n)$.

tolerant k-spanner for $\alpha = O(r^3 \log n)$. For each $F \subseteq V$ with $|F| \le r$, let E_F' be the edges in E_F for which the shortest path in $G \setminus F$ between the endpoints is just the edge. More formally, $E_F' = \{\{u,v\} \in E_F: d_{G \setminus F}(u,v) = \ell(\{u,v\})\}$. It is easy to see that it is sufficient for there to be a path of length at most $k \cdot \ell(\{u,v\})$ between u and v in $G \setminus F$ for every $F \subseteq V$ with $|F| \le r$ and $\{u,v\} \in E_F'$. To show this, consider some fault set F and some $u,v \notin F$. Let $P \subseteq E_F$ be the shortest path from u to v in $G \setminus F$. Since this is a shortest path, every edge in P is actually in E_F' . Thus for every edge $\{x,y\} \in P$ there is a path of length at most $k \cdot \ell(\{x,y\})$ in $G \setminus F$, so by concatenating these paths together we get a path from u to v in $G \setminus F$ of length at most $k \cdot \sum_{e \in P} \ell(e) = k \cdot d_{G \setminus F}(u,v)$.

So we consider a particular such F and $\{u,v\}$ and upper bound the probability that there is no stretch-k path between u and v in $G \setminus F$. Suppose that in some iteration neither u nor v is in J, but all of F is in J. Then since $\{u,v\} \in E_F'$, the spanner that we build on $G \setminus J$ contains a path between u and v of length at most $k \cdot d_{G \setminus J}(u,v) = k \cdot \ell(\{u,v\}) = k \cdot d_{G \setminus F}(u,v)$. Obviously this path also exists in $G \setminus F$, since $F \subseteq J$. So if this happens then H is valid for $\{u,v\}$ and F. The probability that this happens in a particular iteration is clearly $(1-p)^2 \cdot p^r$, which is at least $1/(4r^2)$ as long as $r \geq 2$ (if r = 1 then this probability it 1/8, which does not significantly affect the results). Thus the probability that this never happens in any iteration is at most $(1-\frac{1}{4r^2})^{\alpha} \leq e^{-\alpha/4r^2}$, so if we set $\alpha = \Theta(r^3 \log n)$ this becomes less than $1/n^{C(r+2)}$ for arbitrarily large constant C. Now taking a union bound over all $\{u,v\}$ and F gives the theorem, since there are less than $O(n^r)$ fault sets and at most n^2 edges. \square

COROLLARY 2.2. For every graph G=(V,E) with nonnegative edge lengths $\ell:E\to\mathbb{R}_{\geq 0}$ and every odd $k\geq 1$, there is a polynomial time algorithm that with high probability constructs an r-vertex-tolerant k-spanner with at most $O(r^{2-\frac{2}{k+1}}n^{1+\frac{2}{k+1}}\log n)$ edges.

PROOF. Althöfer et al. [ADD⁺93] showed that the simple greedy spanner construction has size at most $O(n^{1+\frac{2}{k+1}})$. Applying Theorem 2.1 to this construction completes the proof. \square

Since Theorem 2.1 applies to any k-spanner construction, we can apply it to distributed spanner constructions:

THEOREM 2.3. If there is a distributed algorithm A that on every graph builds a k-spanner of size f(n) in t(n) rounds, then there is a distributed algorithm that on any graph builds with high probability an r-fault tolerant k-spanner of size $O(r^3 \log n \cdot f(2n/r))$ in $O(r^3 \log n \cdot t(n))$ rounds.

PROOF. The algorithm is simple: $O(r^3 \log n)$ times, each node independently decides whether or not to join J with

probability 1-1/r, and then A is run on the remainder. This obviously takes at most $O(r^3 \log n \cdot t(n))$ rounds, and the analysis of Theorem 2.1 proves the desired size bound. \square

COROLLARY 2.4. There is a distributed algorithm in the \mathcal{LOCAL} model that in $O(kr^3 \log n)$ rounds constructs with high probability an r-fault tolerant k-spanner with at most $O(kr^{2-\frac{2}{k+1}}n^{1+\frac{2}{k+1}}\log n)$ edges.

PROOF. Apply Theorem 2.3 to the distributed deterministic spanner construction of Derbel, Gavoille, Peleg, and Viennot [DGPV08], which has size $O(kn^{1+\frac{2}{k+1}})$ and runs in O(k) rounds. \square

3. UNIT-LENGTH r-FAULT TOLERANT 2-SPANNER

We now move from general k to the specific case of k =2. It is easy to see (and has long been known) that no non-trivial absolute bounds on the size of a 2-spanner are possible, so following previous work, we instead consider the approximation version. In this section we will mostly work in the directed setting in which every edge e has an arbitrary cost $c_e \geq 0$. This is more general in some ways than the setting in Section 2 (which is undirected and has unit edge costs), but in other ways it is much less general (we now assume that k=2 and edge lengths are unit). Recent work of Dinitz and Krauthgamer [DK11] achieves approximation ratio $O(r \log n)$ for Minimum-Cost r-Fault Tolerant 2-Spanner, and an $O(r \log \Delta)$ ratio when all edge costs are 1 (where Δ is the maximum degree). Here we improve these results to $O(\log n)$ and $O(\log \Delta)$ (for all r) via a different LP relaxation, and also provide a distributed implementation.

3.1 The Previous LP Relaxation

The relaxation in [DK11] uses, at a high level, a characterization of r-fault tolerant 2-spanners based on flows where for every set of r faults, it is possible to send one unit of (integral) flow from u to v along paths of length at most 2 for any edge (u,v) still present in the graph once the faults have been removed. More formally, for each $(u,v) \in E$ let $\mathcal{P}_{u,v}$ denote the paths of length exactly two from u to v, so $\mathcal{P}_{u,v} \cup \{(u,v)\}$ is the set of all paths of length at most 2. Let $\mathcal{F} = \{F \subseteq V : |F| \leq r\}$ be the set of possible fault sets, and as in Section 2, for any possible fault set $F \in \mathcal{F}$ let E_F be the set of edges in E with neither endpoint in F. Let $\mathcal{P}_{u,v}^F$ be the subset of $\mathcal{P}_{u,v} \cup \{(u,v)\}$ that still survives in E_F . The integer program (IP) used by Dinitz and Krauthgamer [DK11] is presented below as IP (2).

min
$$\sum_{e \in E} c_e x_e$$
s.t.
$$\sum_{P \in \mathcal{P}_{u,v}^F : e \in P} f_P^F \le x_e \quad \forall F \in \mathcal{F}, \ \forall (u,v) \in E_F,$$

$$\sum_{P \in \mathcal{P}_{u,v}^F} f_P^F \ge 1 \qquad \forall F \in \mathcal{F}, \ \forall (u,v) \in E_F$$

$$x_e \in \{0,1\} \qquad \forall e \in E$$

$$f_P^F \in \{0,1\} \qquad \forall F \in \mathcal{F}, \ \forall (u,v) \in E_F,$$

$$\forall P \in \mathcal{P}_{u,v}^F$$

This formulation has capacity variables x_e for every edge e, flow variables f_P^F for every possible fault set F and every path $P \in \mathcal{P}_{u,v}^F$ (for every $(u,v) \in E$), and constraints that require flows to obey the capacities and still send one unit of flow for every possible fault set. Even though there are an exponential number of both constraints and variables, it can be solved in polynomial time [DK11].

While IP (2) is the obvious integer programming formulation, its straightforward relaxation to a linear program is not strong enough to give an approximation that is independent of r (despite having an exponential number of both constraints and variables). An easy way to see this is by considering the complete graph. On the complete graph, every vertex obviously needs at least r incoming and outgoing edges, or else it could be isolated with less than r faults. So on K_n the optimum spanner has size at least rn. On the other hand, when we relax the integrality constraints we can set the capacity of every edge to 1/(n-r-2) and still have enough capacity to send one unit of flow from any vertex to any other even after r of them have failed. So the linear program has cost of only $n^2/(n-r-2)$, which is O(n) as long as r < cn for some constant c < 1. Thus the integrality gap of the relaxation is $\Omega(r)$ for an extremely wide range of

3.2 A New LP Relaxation

To get around this problem, we will use a different relaxation based on weighted flow. Before we give our formulation, we first prove a simple and useful characterization of r-fault-tolerant 2-spanners:

LEMMA 3.1. For every (directed) graph G=(V,E), a subgraph H=(V,E') is an r-fault tolerant 2-spanner if and only if for every (u,v) in E either $(u,v) \in E'$ or there are at least r+1 paths of length 2 from u to v in E'

PROOF. Let H be an r-fault tolerant 2-spanner of G, and for the sake of contradiction assume that there is some $(u,v)\in E$ that is not in E' and for which there are at most r paths of length 2 from u to v. Let $W\subseteq V$ be the vertices that are the midpoints of these paths. Then if we let our fault set F be W, in the remaining graph $H\setminus W$ there is no u-v path, while in $G\setminus W$ the edge (u,v) still exists. Thus H is not an r-fault tolerant 2-spanner, giving the contradiction.

For the other direction, suppose that for every $(u,v) \in E$ either $(u,v) \in E'$ or there are at least r+1 paths of length 2 from u to v. Let $F \subseteq V$ with $|F| \le r$ be some fault set. We need to show that H is a valid 2-spanner for $G \setminus F$. As usual, it will suffice for us to show that $d_{H \setminus F}(u,v) \le 2 \cdot d_{G \setminus F}(u,v)$ for all $(u,v) \in E_F$, so since we are assuming unit lengths we just need to show that $d_{H \setminus F}(u,v) \le 2$ for all $(u,v) \in E_F$. Let $(u,v) \in E_F$ be an arbitrary edge in $G \setminus F$. If $(u,v) \in E'$ then obviously $d_{H \setminus F}(u,v) = 1$, and if $(u,v) \not\in E'$ then by assumption there are at least r+1 paths from u to v of length 2 in H. At most r of the intermediate vertices on those paths can be in F, so in $H \setminus F$ there is at least one such path remaining and thus $d_{H \setminus F}(u,v) \le 2$, as required. \square

With this lemma in hand, it is easy to see that the following integer program is an exact formulation of the r-fault tolerant 2-spanner problem. It simply forces the conditions of Lemma 3.1 to be true, i.e. every edge e is either included (has x_e value of 1) or there are r+1 paths of length 2 that are included.

2)

min
$$\sum_{e \in E} c_e x_e$$
s.t.
$$\sum_{P \in \mathcal{P}_{u,v}: e \in P} f_P \le x_e \qquad \forall (u,v) \in E, \\ \forall e \in E$$

$$(r+1)x_{(u,v)} + \sum_{P \in \mathcal{P}_{u,v}} f_P \ge r+1 \quad \forall (u,v) \in E$$

$$x_e \in \{0,1\} \qquad \forall e \in E$$

$$f_P \in \{0,1\} \qquad \forall (u,v) \in E, \\ \forall P \in \mathcal{P}_{u,v}$$
(3)

So now we have a different IP formulation than the one that was used in [DK11] to get an $O(r \log n)$ -approximation. Unfortunately, it is still not strong enough to yield an approximation ratio independent of r; there are still simple examples that give an integrality gap of $\Omega(r)$. For example, consider a graph with nodes u and v and an edge of cost M from u to v (for some arbitrarily large M), together with r nodes w_1, \ldots, w_r and an edge of cost 1 from u to w_i and from w_i to v for all $i \in [r]$. The set of all w_i nodes is a valid fault set, so the optimum spanner needs to include the (u, v) edge in order to still be valid. So the optimum spanner has cost at least M. On the other hand, the LP can set x_e to 1 for all edges e incident on some w_i , and set $x_{(u,v)} = 1/(r+1)$. This has cost of only M/(r+1) + 2r. By setting M large enough, we get a gap of $\Omega(r)$.

We will strengthen the relaxation by adding a set of valid inequalities that are essentially the knapsack-cover inequalities of Carr et al. [CFLP00] applied to this IP. Let $(u,v) \in E$, and let $\mathcal{W}_{u,v} = \{W \subseteq P_{u,v} : |W| \le r\}$. Consider some arbitrary subset $W \in \mathcal{W}_{u,v}$. If $x_{(u,v)} = 0$, then the covering inequality for (u,v) implies that $\sum_{P \in \mathcal{P}_{u,v}} f_P \ge r+1$, and thus $\sum_{P \in \mathcal{P}_{u,v} \setminus W} f_P \ge r+1 - |W|$. On the other hand, if $x_{(u,v)} = 1$ then clearly $(r+1-|W|)x_{(u,v)} \ge r+1 - |W|$. So for all $(u,v) \in E$ and all $W \subseteq \mathcal{P}_{u,v}$ with $|W| \le r$, we can add the constraint

$$(r+1-|W|)x_{(u,v)} + \sum_{P \in \mathcal{P}_{u,v} \setminus W} f_P \ge r+1-|W|.$$

To simplify notation, for any such W let $r_W = r + 1 - |W|$. These are the knapsack-cover inequalities, and when we add them to our IP formulation and relax the integrality constraints we get the following LP relaxation:

$$\min \sum_{e \in E} c_e x_e$$
s.t.
$$\sum_{P \in \mathcal{P}_{u,v}: e \in P} f_P \le x_e \qquad \forall (u,v) \in E, \\
\forall e \in E$$

$$r_W x_{(u,v)} + \sum_{P \in \mathcal{P}_{u,v} \setminus W} f_P \ge r_W \quad \forall (u,v) \in E, \\
\forall W \in \mathcal{W}_{u,v}$$

$$0 \le x_e \le 1 \qquad \forall e \in E$$

$$f_P \ge 0 \qquad \forall (u,v) \in E, \\
\forall P \in \mathcal{P}_{u,v}$$
(4

We refer to the first type of constraints as *capacity* constraints, the second type as *knapsack-cover* constraints (or

inequalities), and the third as *multiplicity* constraints. This relaxation has a polynomial number of variables but a possibly exponential number of constraints, so we first need to show that we can solve it. To do this we construct a separation oracle, which allows us to solve it in polynomial time by using the Ellipsoid algorithm.

LEMMA 3.2. There is a polynomial time algorithm that solves LP (4).

PROOF. We want to construct a separation oracle. Note that there are only a polynomial number of capacity constraints and multiplicity constraints, so we can check them all in polynomial time. To find a violated knapsack-cover inequality, note that if there is some $(u,v) \in E$ and some $W \subseteq \mathcal{P}_{u,v}$ that violates the inequality, then the set W' which consists of the |W| paths in $\mathcal{P}_{u,v}$ with the largest f_P value also violates the inequality. So for every $(u,v) \in E$, for every $i \in [0,r]$, it suffices to check the constraint for (u,v) and the i paths in $\mathcal{P}_{u,v}$ with largest flow. Since $r \leq n$, this takes only polynomial time. \square

3.3 O(log n)-approximation

We now give the main result of this section.

Theorem 3.3. There is a randomized $O(\log n)$ approximation for Minimum Cost r-Fault Tolerant 2-Spanner on directed graphs that works for all r.

PROOF. The first step of the algorithm is to solve LP (4) using Lemma 3.2. We then round the solution using Algorithm 1 below. (This rounding algorithm was designed in [DK11] but was used to round a different relaxation, hence they were forced to set $\alpha = \Theta(r \log n)$ and the analysis therein is not applicable here.)

Algorithm 1: Rounding algorithm for r-fault tolerant 2-spanner.

Input: Graph G = (V, E), fractional solution $\langle x, f \rangle$ to LP (4)

- 1 Set $\alpha = C \ln n$ (for a large enough constant C).
- **2** For every $v \in V$ choose independently a random threshold $T_v \in [0, 1]$.
- **3** Output $E' = \{(u, v) \in E : \min\{T_u, T_v\} \le \alpha \cdot x_{(u,v)}\}.$

We first show that the cost of the solution is likely to be at most 6α times the LP value. The probability that some edge e is selected to be in E' is at most $2\alpha x_e$, so the expected cost of the solution E' is $\sum_{e\in E} c_e \cdot 2\alpha x_e = 2\alpha \sum_e c_e x_e$. By Markov's inequality, the cost of the solution E' exceeds $8\alpha \sum_e c_e x_e$ with probability at most 1/4.

We now argue that this algorithm returns a valid r-fault tolerant 2-spanner with high probability. We say that E' satisfies an edge (u,v) if either $(u,v) \in E'$ or E' contains at least r+1 length 2 paths from u to v. By Lemma 3.1, if E' satisfies all edges then it is a valid r-fault tolerant 2-spanner. Consider some edge $(u,v) \in E$. Order the paths in $\mathcal{P}_{u,v}$ in nonincreasing order by their flow values in the LP solution, so P_i is the path with the ith largest flow. Let $W_i = \{P_1, P_2, \ldots, P_i\}$, and let $i^* = \max\{i : f_{P_i} \ge 1/\alpha\}$. If $i^* > r$ then r+1 paths have flow value at least $1/\alpha$, so both of the edges in each path have x value at least $1/\alpha$, so they are included in E' with probability 1. Thus (u,v) is satisfied with probability 1.

On the other hand, suppose that $i^* \leq r$. Let us denote $r' = r + 1 - i^* \geq 1$. By the knapsack-cover constraint for (u, v) and W_{i^*} , we know that

$$r'x_{(u,v)} + \sum_{P \in \mathcal{P}_{u,v} \setminus W_{i*}} f_P \ge r'$$

If $r'x_{(u,v)} \geq r'/2$ then $x_{(u,v)} \geq 1/2$ and thus (u,v) is included in E' with probability 1, satisfying (u,v). Otherwise it must be the case that $\sum_{P \in \mathcal{P}u,v \setminus W_{i^*}} f_P \geq r'/2$. For $P \in \mathcal{P}_{u,v}$, let I_P be an indicator for the event that the T value of the middle vertex is at most α times the flow value f_P (formally, if P = (u,z,v) then $I_P = 1_{\{T_z \leq \alpha f_P\}}$), and observe that this event implies that both edges of P are included in E' (because then we have $T_z \leq \alpha \cdot \min\{x_{(u,z)}, x_{(z,v)}\}$). Note that for $P \in W_{i^*}$, we have $I_P = 1$ with probability 1. For $P \in \mathcal{P}_{u,v} \setminus W_{i^*}$, we have $I_P = 1$ with probability at least $\alpha f_P \in [0,1]$. The number of paths from $\mathcal{P}_{u,v} \setminus W_{i^*}$ included in E' is clearly at least $\sum_{P \in \mathcal{P}_{u,v} \setminus W_{i^*}} I_P$, and we can bound that last quantity (which is a sum of independent indicators) by a Chernoff bound (see e.g. [MR95, DP09]). Its expectation is

$$\mathbb{E}\left[\sum_{P\in\mathcal{P}_{u,v}\setminus W_{i^*}}I_P\right] \geq \sum_{P\in\mathcal{P}_{u,v}\setminus W_{i^*}}\alpha f_P \geq \alpha r'/2,$$

so by our choice of $\alpha = C \log n$ for a large enough C,

$$\Pr\left[\sum_{P \in \mathcal{P}_{u,v} \setminus W_{i^*}} I_P \le \alpha r'/4\right] \le e^{-\Omega(\alpha r')} \le 1/n^{\Omega(C)} \le 1/n^3.$$

Thus with high probability the total number of length 2 paths between u and v included in E' is at least $i^* + \alpha r'/4 \ge r+1$, and so (u,v) is satisfied. The theorem follows by taking a union bound over these events for all edges (u,v), and the aforementioned event that the solution's cost exceeds 6α times the LP value. \square

3.4 Bounded-Degree Graphs

When the maximum degree of the graph is bounded by Δ and the edge costs c_e are all 1, we can improve Theorem 3.3 slightly and give an $O(\log \Delta)$ -approximation. We simply change the inflation parameter α in Algorithm 1 to be $O(\log \Delta)$ instead of $O(\log n)$. We then need a more careful analysis, using an algorithmic version of the Lovász Local Lemma.

Theorem 3.4. There is a randomized $O(\log \Delta)$ approximation for Minimum Cost r-Fault Tolerant 2-Spanner on directed graphs in which $c_e=1$ for all $e\in E$ and the maximum (in and out) degree is at most $\Delta\geq 2$.

We shall use the following constructive version of the symmetric Lovász Local Lemma, which is an immediate corollary of the nonsymmetric version proved by Moser and Tardos [MT10].

Lemma 3.5 (Moser and Tardos [MT10]). Let \mathcal{P} be a finite set of mutually independent random variables in a probability space. Let \mathcal{A} be a finite set of events determined by the variables in \mathcal{P} . Suppose that each $A \in \mathcal{A}$ is mutually independent of all but at most d other events in \mathcal{A} , and suppose that $\Pr[A] \leq p$ for all $A \in \mathcal{A}$. If $ep(d+1) \leq 1$ then there exists an assignment of values to the variables \mathcal{P}

such that no event $A \in \mathcal{A}$ occurs. Moreover, there is a randomized algorithm that finds such an assignment in expected time $O(|\mathcal{P}| + |\mathcal{A}| \cdot |\mathcal{P}|/d)$.

PROOF OF THEOREM 3.4. Consider a directed graph G with unit edge costs $c_e = 1$ and vertex degrees bounded by Δ . Consider a solution to the LP relaxation (4), and apply Algorithm 1 to it but with inflation factor $\alpha = C \log \Delta$ instead of $C \log n$.

For an edge $(u, v) \in E$, let $A_{u,v}$ be the event that E' does not satisfy this edge, i.e. $(u, v) \notin E'$ and the graph G' = (V, E') has less than r + 1 paths of length 2 from u to v. The analysis of Theorem 3.3 shows (after modifying (5) with our new value of α), that

$$\Pr[A_{u,v}] \le e^{-\Omega(\alpha)} \le 1/\Delta^{\Omega(C)}.$$

Furthermore, note that $A_{u,v}$ depends only on the random variables T_z for $z \in (N^+(u) \cap N^-(v)) \cup \{u,v\}$. Here and throughout, $N^+(u)$ and $N^-(u)$ denote the out-neighbors and in-neighbors respectively of $u \in V$. Observe that $A_{u,v}$ is independent of all but Δ^3 other events $A_{u',v'}$, simply because there are at most Δ choices for each of z, u', and v'.

We could now apply Lemma 3.5 to these events. The underlying mutually independent random variables \mathcal{P} would be the T_u variables, and the "bad events" \mathcal{A} would be the events $A_{u,v}$. This would give us an algorithm that in polynomial time returned a valid r-fault tolerant k-spanner, but we also need a bound on the cost of this spanner. The analysis via Markov's inequality in Theorem 3.3 is too weak now, because when we apply the algorithm of Lemma 3.5 we change the overall distribution in a way that might destroy the cost bound. We need to integrate the cost analysis into the events that Lemma 3.5 is applied to, so at a high level we employ a more local approach where the cost of E' is split among the vertices and events bounding the cost are added to the $A_{u,v}$ events. More specifically, we shall create many events, each of which controls how the cost of E' compares locally with the cost of the LP, and then apply the Local Lemma to the new events together with the $\{A_{u,v}\}$ events. A formal argument follows.

For each vertex $u \in V$, let the random variable Z_u^+ be the number of outgoing edges (u,v) for which $T_v \leq \alpha \cdot x_{(u,v)}$, and let Z_u^- be the number of incoming edges (v,u) for which $T_v \leq \alpha \cdot x_{(u,v)}$. Informally, $Z_u^+ + Z_u^-$ is the number of edges incident to u whose inclusion in E' can be charged to their other endpoint. The algorithm's cost is $|E'| \leq \sum_{u \in V} (Z_u^+ + Z_u^-)$, since every edge (u,v) included in E' adds 1 to either Z_u^+ or Z_v^- (or both).

For each vertex $v \in V$, let B_u be the event that $Z_u^+ + Z_u^- > 4\alpha(\sum_{(u,v)\in E} x_{(u,v)} + \sum_{(v,u)\in E} x_{(v,u)})$. We would like to show that this event happens only with small probability. Note that $\mathbb{E}[Z_u^+] = \sum_{(u,v)\in E} \min\{\alpha x_{(u,v)}, 1\} \leq \alpha \sum_{(u,v)\in E} x_{(u,v)}$, so by a Chernoff bound (see e.g. [MR95, DP09]) we get

$$\Pr\left[Z_u^+ > 2\alpha \sum_{(u,v) \in E} x_{(u,v)}\right] \le e^{-(1/3)(C\ln\Delta) \sum_{(u,v) \in E} x_{(u,v)}}$$

$$< \Delta^{-C/3},$$

where in the final inequality we assume there is at least one outgoing edge from u and thus $\sum_{(u,v)\in E} x_{(u,v)} \geq 1$ (since otherwise $Z_u^+=0$ with probability 1). We can use a similar

argument to get the same bound for Z_u^- , and by combining them we get that

$$\Pr[B_u] \le \Pr\left[Z_u^+ > 2\alpha \sum_{(u,v) \in E} x_{(u,v)}\right]$$
$$+ \Pr\left[Z_u^- > 2\alpha \sum_{(v,u) \in E} x_{(v,u)}\right]$$
$$\le 2\Delta^{-C/3}.$$

We now apply Lemma 3.5 to the events $A_{u,v}$ and B_u . Note that B_u depends only on the random variables T_z for $z \in N^+(u) \cup N^-(u)$, and recall that $A_{u,v}$ depends only on T_z for $z \in N^+(u) \cap N^-(v)$. Thus each event is mutually independent of all but $O(\Delta^3)$ other events — for an event $A_{u,v}$ we exclude at most Δ^3 events $A_{u',v'}$ and at most $2\Delta^2$ events $B_{u'}$; for an event B_u we exclude at most $4\Delta^2$ events $B_{u'}$ and at most $2\Delta^3$ events $A_{u',v'}$. We can thus apply Lemma 3.5 with dependency parameter $d = O(\Delta^3)$, because by setting sufficiently large C, the probability of each event is at most a suitable $p = \Delta^{-\Omega(C)} < 1/e(d+1)$. Since the number of events is at most $O(n^2)$ and the number of underlying variables is only n, we conclude that there is a polynomial time algorithm to find the underlying variables T_u so that none of the events $A_{u,v}$ and B_u occur. This implies that G' = (V, E') is an r-fault tolerant 2-spanner of G of cost

$$|E'| \le \sum_{u \in V} \left(Z_u^+ + Z_u^- \right) \le 8\alpha \sum_{(u,v) \in E} c_{u,v} x_{(u,v)}$$

$$< O(\log \Delta) \cdot \text{LP},$$

which proves Theorem 3.4. \square

3.5 Distributed Construction

We now show how to adapt and use the $O(\log n)$ approximation that we designed in Section 3.3 to give a distributed $O(\log n)$ -approximation. We will assume that communication along an edge is bidirectional, even if the graph is directed. The main problem that we run into when trying to design a distributed algorithm based on Algorithm 1 is solving the linear program. If we had a solution, and every vertex knew the x_e value of its incident edges, then we would be done; the rounding scheme in Algorithm 1 is entirely local, so every vertex $v \in V$ would just locally pick its threshold T_v and include the appropriate edges. If we want both endpoints of an edge to know that it has been included in the spanner, we can then just have every vertex tell all of its neighbors (in a single round) which edges it bought based on its threshold.

In order to (approximately) solve the LP we partition the graph into clusters, solve the LP separately on each cluster, and then repeat this process several times, eventually taking the average. This technique is quite similar to the work of Kuhn, Moscibroda, and Wattenhofer [KMW06], who showed how to approximately solve *positive* LPs using the graph decompositions of Linial and Saks [LS93].

The fundamental tool that we will use is the ability to quickly compute a good padded decomposition, which is a basic tool in metric embeddings that has found numerous applications in approximation and online algorithms (e.g. for network design problems). This notion is essentially a version of low-diameter decompositions, such as a sparse covers [AP90]. This specific version was (probably) introduced by Rao [Rao99], who observed that it can be derived from

an earlier construction of Klein, Plotkin and Rao [KPR93]. An explicit formulation of padded decompositions appeared only later, in [KL03, GKL03], and used a construction of Bartal [Bar96]. The definition given below is actually a special case of the usual notion, where the so-called padding requirement is a unit radius around each vertex, i.e. just the vertex's neighborhood.

Let $\mathcal{T} = \mathcal{T}(V)$ denote the set of all partitions of V (irrespective of the graph structure). For a partition $P \in \mathcal{T}$, we call each set $C \in P$ a cluster. Let G' be the undirected graph corresponding to G, and define the diameter of C to be $\operatorname{diam}(C) = \max_{u,v \in C} d_{G'}(u,v)$ (this is usually called weak diameter, because it corresponds to the shortest u-v path in G', possibly going out of C along the way). Finally, for $x \in V$ and a partition $P \in \mathcal{T}$, we let P(x) denote the cluster of P that contains x. Since G' is undirected, we let N(x) denote the set of all neighbors of x in G'

DEFINITION 3.6. A padded decomposition of G is a probability measure μ on $\mathcal T$ that satisfies the following two conditions:

- 1. For every $P \in \text{supp}(\mu)$ and every $C \in P$ we have $\text{diam}(C) \leq O(\log n)$.
- 2. For every $x \in V$ we have $\Pr_{P \sim \mu}[N(x) \subseteq P(x)] \ge 1/2$.

It is known that every metric space admits such a padded decomposition, and there are polynomial-time randomized algorithms to sample from such a decomposition [Bar96, FRT04]. It is convenient to assign to each cluster a vertex, called the *cluster center*. One could always choose an arbitrary vertex in the cluster (e.g. one whose identifier is the smallest), but in many constructions there is a natural center point. The next lemma is a straightforward adaptation of the construction of Bartal [Bar96] to the distributed context; it can also be viewed as a slight modification of the graph decompositions of Linial and Saks [LS93].

LEMMA 3.7. There is a distributed algorithm that runs in $O(\log n)$ rounds and with high probability samples from a padded decomposition, so that every vertex knows the cluster containing it, meaning all other vertices in the same cluster. Every cluster C also has a cluster center $v \in V$ (which is not necessarily in the cluster) with the property that $\operatorname{diam}(C \cup \{v\}) \leq O(\log n)$.

PROOF. The construction of Bartal [Bar96] is simple, and is usually described iteratively. (As mentioned above, the padding property is not formally proved there, but it can be derived from the analysis therein, see also [KL03, GKL03]). Working in the metric completion of G' (so removing vertices does not change distances), repeat the following procedure until every vertex has been assigned to some cluster: Pick an arbitrary vertex u from those that have not yet been assigned a cluster. Randomly pick a radius r_u from the geometric distribution with some constant parameter p>0. Create a new cluster consisting of u and all unclustered vertices that are within distance r_u of u.

While this procedure is phrased iteratively, it quite obviously can be made distributed with only minor changes. First, every vertex $u \in V$ locally chooses a value r_u from the geometric distribution with parameter p. Then every node u simultaneously sends a message containing the ID of u to all nodes within distance $\min\{r_u, O(\log n)\}$ of u. Note

that this take only $O(\log n)$ rounds, and with high probability $\max_u\{r_u\} \leq O(\log n)$ (the analysis of [KL03] shows that this truncation at $O(\log n)$ does not significantly affect the padding probability). Now every node chooses as a cluster center the sender with the smallest ID (i.e. the sender that comes earliest in the lexicographic ordering) of the vertices whose messages it received. The only difference between the output of this algorithm and Bartal's algorithm is that in Bartal's algorithm only unclustered nodes can be the center of a new cluster, while in our variation every vertex (in lexicographic order) gets the chance to create a cluster (which it might not be a member of itself). It is well known (see e.g. [KL03, GKL03]) that this change does not affect anything in the analysis.

We remark that the construction above has a natural choice of cluster centers. Under this choice, a cluster C might not contain its center $v \in V$, but $\operatorname{diam}(C \cup \{v\}) \leq O(\log n)$, which is sufficient for our purposes. \square

Now that we can construct padded decompositions, we want to use them to decompose LP (4) into "local" parts. Let P be a partition sampled from μ . For each cluster $C \in P$, let N(C) denote the set of vertices in $V \setminus C$ that are adjacent (in G') to at least one vertex in C, let $\delta(C) \subseteq E$ be all edges with one endpoint in C and one endpoint not in C, and let $E(C) \subseteq E$ be the set of edges with both endpoints in C. Let G(C) be the subgraph of G induced by $C \cup N(C)$. We define LP(C) to be LP (4) for G(C), but where edges in $\delta(C)$ are modified to have cost 0.

Let LP^* be the value of an optimal solution to LP (4), and let $LP^*(C)$ be the value of an optimal solution to LP(C).

LEMMA 3.8. $\sum_{C \in P} LP^*(C) \leq LP^*$ for every partition $P \in \mathcal{T}$.

PROOF. Let $\langle x,f\rangle$ be an optimal fractional solution to LP (4). We want to use this solution to build fractional solutions to LP(C) for all $C\in P$ whose total cost is at most LP*. For each cluster $C\in P$, define a solution $\langle x^C,f^C\rangle$ for LP(C) as follows: Let $x_e^C=x_e$ if $e\in E(C)$ and let $x_e^C=1$ if $e\in \delta(C)$. Note that this already satisfies all of the knapsack-cover constraints for edges in $\delta(C)$. For edges $(u,v)\in E(C)$, note that every path in $\mathcal{P}_{u,v}$ appears in G(C), so we can set $f_P^C=f_P$ for these paths. Since these flows satisfy all of the knapsack-cover constraints in LP (4), they also satisfy all of the knapsack-cover constraints in LP(C). All other flows f_P^C (e.g. between vertices in N(C)) are set to 0. Obviously the capacity constraints are satisfied, and thus $\langle x^C,f^C\rangle$ is a feasible solution to LP(C).

Since in LP(C) the edges in $\delta(C)$ have cost 0, and every edge of E is in E(C) for at most one cluster C,

$$\sum_{C \in P} \mathrm{LP}^*(C) = \sum_{C \in P} \sum_{e \in E(C)} c_e x_e^C \le \sum_{e \in E} c_e x_e = \mathrm{LP}^*,$$

which proves the lemma. \Box

With this lemma in hand, we can now design a distributed approximation algorithm and prove the main theorem of this section.

Theorem 3.9. There is a distributed algorithm that terminates in $O(\log^2 n)$ rounds and is an $O(\log n)$ approximation to Minimum Cost r-Fault Tolerant 2-Spanner.

Algorithm 2: Distributed algorithm for r-fault tolerant 2-spanner.

```
1 for i\leftarrow 1 to t=O(\log n) do
2 | Sample a partition P_i from \mu using Lemma 3.7  
// we assume the center of each cluster C\in P_i knows G(C)
3 | The center of each cluster C\in P_i solves \operatorname{LP}(C) using Lemma 3.2, and sends the solution \langle x^{C,i}, f^{C,i} \rangle to all vertices in C
4 For each edge (u,v)\in E, let \mathcal{I}_{(u,v)}=\{i:P_i(u)=P_i(v)\} // these are the iterations in which both endpoints are in same cluster
5 \tilde{x}_e\leftarrow \min\{1,\frac{4}{t}\sum_{i\in\mathcal{I}_e}x_e^{P_i(e),i}\} // P_i(e) is the cluster of P_i containing both endpoints of e
6 Round \tilde{x}_e using Algorithm 1 // each edge is rounded by its endpoints
```

PROOF. Our algorithm is show as Algorithm 2. We first prove the time bound. Lemma 3.7 implies that sampling from μ takes only $O(\log n)$ rounds, and since the diameter of every cluster is at most $O(\log n)$ the other two steps of the loop also take only $O(\log n)$ rounds. Since we execute the loop $O(\log n)$ times, the number of rounds needed to complete the loop is at most $O(\log^2 n)$. After the loop, each vertex can compute x_e for all incident edges e without any extra communication (since each endpoint of an edge e knows \mathcal{I}_e and the LP values for that iteration). Finally, as already pointed out, the rounding of Algorithm 1 can be done locally, with one extra round used to make sure that both endpoints of an edge know if the edge was included by the rounding. Thus the total number of rounds is $O(\log^2 n)$, as claimed.

To prove that this algorithm returns an $O(\log n)$ approximation, we will show that with high probability the \tilde{x}_e values it computes form a feasible solution to LP (4) (when appropriate flow values \tilde{f}_P are chosen) of cost at most $O(\text{LP}^*)$. Once we have this, the analysis of Theorem 3.3 implies that the rounding step outputs an r-fault tolerant 2-spanner G' = (V, E') whose cost is $O(\log n) \sum_e c_e \tilde{x}_e \leq O(\log n) \cdot \text{LP}^*$ with probability at least 2/3. To bound the cost, note that the $\tilde{x}_e/4$ values are just the averages of the LP(C) values for all rounds in which the edge e does not have cost 0. In other words, $\sum_e c_e \tilde{x}_e \leq \frac{4}{t} \sum_{i=1}^t \sum_{C \in P_i} \text{LP}^*(C) \leq 4 \text{LP}^*$, where the final inequality is from Lemma 3.8. So it just remains to show that the \tilde{x}_e 's form a feasible solution to LP (4).

To prove this, consider an edge e = (u, v), and let $\mathcal{I}'_e \subseteq \mathcal{I}_e$ be the set of iterations i in which $N(u) \cup \{u\}$ is all in the same cluster of P_i . By the second property of padded decompositions, the probability that $N(u) \cup \{u\}$ is all in the same cluster is at least 1/2. The iterations are independent, so a straightforward Chernoff bound implies that $\Pr[|\mathcal{I}'_e| \geq t/4] \geq 1 - 1/n^3$. For a path $P \in \mathcal{P}_{u,v}$, set $\tilde{f}_P = \frac{1}{|\mathcal{I}'_{(u,v)}|} \sum_{i \in \mathcal{I}'_{(u,v)}} f_P^{P_i(u,v),i}$. In other words, the flow along a path from u to v is equal to the average flow along it in the LP solutions that were computed in iterations when $N(u) \cup \{u\}$ were all in the same cluster.

The capacity constraints are obviously satisfied, since each iteration satisfies the capacity constraints, and the edge capacities are scaled by 4/t while flows are scaled by $1/|\mathcal{I}'_{u,v}| \leq$

4/t. Note that here we depend on the fact that *all* of N(u) is in the same cluster as u; if some vertex $z \in N(u)$ were in a different cluster, then in the LP solution for the cluster containing u and v there could be flow sent from u to v through z. This flow would not have the corresponding capacity added to the \tilde{x}_e variables, which would be a problem.

Similarly, consider the knapsack-cover constraint for some $(u, v) \in E$ and some $W \in \mathcal{W}_{u,v}$. Then since we could send enough flow in each iteration in $\mathcal{I}'_{(u,v)}$, when we take the average we can still send enough flow, i.e.

$$r_{W}\tilde{x}_{(u,v)} + \sum_{P \in \mathcal{P}_{u,v} \setminus W} \tilde{f}_{P}$$

$$\geq \sum_{i \in \mathcal{I}'_{(u,v)}} \left(\frac{4}{t} r_{W} x_{u,v}^{i} + \sum_{P \in \mathcal{P}_{u,v} \setminus W} \frac{1}{|\mathcal{I}'_{(u,v)}|} f_{P}^{P_{i}(u,v),i} \right)$$

$$\geq \frac{1}{|\mathcal{I}'_{(u,v)}|} \sum_{i \in \mathcal{I}'_{(u,v)}} \left(r_{W} x_{u,v}^{i} + \sum_{P \in \mathcal{P}_{u,v} \setminus W} f_{P}^{P_{i}(u,v),i} \right)$$

$$= r_{W}$$

where the last inequality is by the knapsack-cover constraint for the cluster $P_i(u,v)$. Thus we have a valid LP solution, so rounding using Algorithm 1 gives an $O(\log n)$ approximation. \square

Remark.

While for our purposes it was enough to solve the LP to within a constant factor (since we lose an $O(\log n)$ factor in the rounding anyway), it is easy to see that we could in fact solve the LP to within a $(1+\epsilon)$ factor. First, we could change the padded decomposition to have the padding property (u and N(u) are all in the same cluster) hold with probability at least $1-\epsilon$, which would require increasing the diameter of the clusters, and thus the number of rounds it takes to solve the LP, by an $O(1/\epsilon)$ factor. Second, when we apply the Chernoff bound, instead of asking the number of times the padding event occurs to be at least t/4, we could ask that it is at least $(1-\epsilon)^2t$. By increasing t by an $O(1/\epsilon^2)$ factor, we still get the right probabilities. Overall, the number of rounds would now be $O(\epsilon^{-3}\log n)$.

4. CONCLUSIONS AND FUTURE WORK

This paper considers the problem of constructing r-fault tolerant spanners and gives two basic constructions. For general stretch bounds $k \geq 3$, we show how to construct r-fault tolerant k-spanners whose size is at most polynomially (in r) larger than spanners without fault tolerance, improving over the previous exponential dependency (on r) of [CLPR09]. Our main technique is oversampling failure sets, in order to handle many of them in one iteration. An interesting open question is to provide nontrivial lower bounds on the size of the best r-fault tolerant k-spanner; to the best of our knowledge, no such bounds are known other than the trivial bound of $\Omega(rn)$ and those that apply even when r=0.

For k=2 and unit edge lengths we design an $O(\log n)$ -approximation algorithm (for all r), improving over the previous $O(r \log n)$ factor of [DK11] and showing that the approximation ratio can be independent of the desired amount of fault tolerance r. Our main technique here is to design a new linear programming relaxation that includes the exponentially many knapsack-cover inequalities of [CFLP00].

We also provided a distributed version of the algorithm, and showed that when all edge costs are 1 the approximation can be improved to $O(\log \Delta)$. An interesting open question is to improve this ratio to $O(\log(|E|/|V|))$, which would match the approximation known for the non-fault tolerant version.

5. REFERENCES

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