

The Fractional Prize-Collecting Steiner Tree Problem on Trees^{*}

Extended Abstract

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Abstract. We consider the fractional prize-collecting Steiner tree problem on trees. This problem asks for a subtree T containing the root of a given tree $G = (V, E)$ maximizing the ratio of the vertex profits $\sum_{v \in V(T)} p(v)$ and the edge costs $\sum_{e \in E(T)} c(e)$ plus a fixed cost c_0 and arises in energy supply management. We experimentally compare three algorithms based on parametric search: the binary search method, Newton's method, and a new algorithm based on Megiddo's parametric search method. We show improved bounds on the running time for the latter two algorithms. The best theoretical worst case running time, namely $O(|V| \log |V|)$, is achieved by our new algorithm. A surprising result of our experiments is the fact that the simple Newton method is the clear winner of the tested algorithms.

1 Introduction

We consider a variant of the well-studied *Steiner tree problem in graphs*, namely the *prize-collecting Steiner tree problem*. This problem, where we want to find a subtree of a graph that maximizes an objective function that depends on the profits of the vertices and the costs of the edges, arises in the design of supply networks like district heating systems. It was first mentioned by Segev [14] where it appears as a special case of the node-weighted Steiner tree problem and is called the *Single Point Weighted Steiner Tree problem*. The author proves NP-hardness of the problem, presents integer linear programming formulations, and uses Lagrangean relaxation and heuristics to compute lower and upper bounds for these formulations, respectively. In [4], Duin and Volgenant relate the node-weighted (and thus also the prize-collecting) variant to the classical Steiner tree problem. They adapt reduction techniques and show how the

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rooted prize-collecting Steiner tree problem can be transformed into the directed version of the classical Steiner tree problem.

In [6], Fischetti studies the facial structure of a generalization of the problem, the so called *Steiner arborescence problem*. Goemans studies the polyhedral structure of the node-weighted Steiner tree problem [7] and shows that his characterization is complete in case the input graph is series-parallel.

Approximation results are given by Bienstock *et al.* [1] and by Goemans and Williamson [8]; the latter present a purely combinatorial $O(n^2 \log n)$ -time primal-dual $(2 - \frac{1}{n-1})$ -approximation algorithm, where n denotes the number of vertices in the graph and the objective is to minimize the edge costs plus the prizes of the nodes not spanned. For the more realistic objective to maximize the sum of the profits minus the costs, Feigenbaum *et al.* [5] prove that it is NP-hard to approximate the problem to a constant factor.

In this paper, we look at the special case where the potential network is a tree and instead of the linear objective function, we look at the fractional version of the problem which maximizes the ratio of the sum of the profits and the sum of the (fixed and variable) costs.

Section 2 contains some preliminaries including the description of a linear time algorithm for optimizing the linear objective function. In Section 3, we present three different algorithms that use the parametric formulation: a binary search algorithm, Newton's method and our new variant based on Megiddo's method for parametric search. We show a worst case running time of Newton's method of $O(|V|^2)$, and of our new algorithm of $O(|V| \log |V|)$. In Section 4, we report on extensive computational experiments. Surprisingly for us, our experiments show that Newton's method, although having worst case running time of $O(|V|^2)$, outperforms the two other methods on our benchmark set. Finally, in Section 5 we summarize the results.

2 Preliminaries

In this section, we provide some basic definitions and describe a linear time algorithm for solving the linear version of the prize-collecting Steiner tree problem (PCST problem). A closely related dynamic programming algorithm can also be found in [15] (where trees with only node-weights are considered).

Let $G = (V, E)$ be an undirected graph, $r \in V$ a root vertex of G , $p : V \rightarrow \mathbb{R}^+ \cup \{0\}$ a profit function on the vertices, and $c : E \rightarrow \mathbb{R}^+ \cup \{0\}$ a cost function on the edges.

The *Fractional Prize Collecting Steiner Tree problem* (FPCST) consists of finding a connected subgraph $T = (V', E')$ of G with $r \in V'$ that maximizes the ratio of the profits and the costs:

$$\text{profit}(T) = \frac{\sum_{v \in V'} p(v)}{c_0 + \sum_{e \in E'} c(e)},$$

In the construction of a district heating network, the edges correspond to the potential pipes and the vertices to customers or forks in the pipe network. The

costs of the edges are the costs of the pipes, the profits of the vertices the revenue generated by the customers and the fixed cost c_0 the cost for building the heating plant.

The *Linear Prize Collecting Steiner Tree problem* (LPCST) consists of finding a connected subgraph $T = (V', E')$ of G with $r \in V'$ that maximizes the difference of the profits and the costs:

$$\text{profit}(T) = \sum_{v \in V'} p(v) - \sum_{e \in E'} c(e) .$$

Note that fixed costs are irrelevant if we optimize a linear objective function.

If $T = (V, E)$ is a tree with root r , then the function $\text{parent}(v)$ assigns every vertex $v \in V \setminus \{r\}$ a unique vertex u which is the vertex following v on the path from v to r . The *subtree* rooted at v consists of all vertices and edges reachable from v without passing the vertex $\text{parent}(v)$. The set $C(v)$ of *children* of v is the set that contains all vertices u with $\text{parent}(u) = v$. A subtree of T is *optimal*, if there is no other subtree of T with a higher profit. We recursively define a value $l(v)$ and a subtree $T(v)$ for each vertex $v \in V$ as

$$l(v) = p(v) + \sum_{u \in C(v)} \max\{0, l(u) - c(u, v)\} . \tag{1}$$

The subtree $T(v) = (V(v), E(v))$ with profit $l(v)$ is defined in the following way:

$$V(v) = \{v\} \cup \bigcup_{u \in C(v)} \{V(u) \mid l(u) - c(u, v) \geq 0\}$$

$$E(v) = \bigcup_{u \in C(v)} \{(u, v) \cup E(u) \mid l(u) - c(u, v) \geq 0\} .$$

If $c(u, v) > l(u)$ for a vertex u with $\text{parent}(u) = v$ it does not pay off to include the subtree rooted at u via edge (u, v) (the only possible connection towards r), and we decide to cut off the edge (u, v) together with the corresponding subtree. This decision can be made locally, as soon as the value $l(u)$ is known. It is not hard to construct an algorithm for LPCST that uses these facts and runs in linear time (see [10] for details). The optimal subtree rooted at v is $T(v)$ with $l(v)$ as its profit (the correctness of this algorithm follows easily by induction).

When solving FPCST on trees, in contrast to the linear case, we cannot make local decisions anymore without looking at the whole problem. The following section presents the parametric formulation of the problem that allows us to decide in linear time if a given value t is smaller, equal, or greater than the value of an optimal solution of FPCST.

3 Algorithms Based on Parametric Formulation

To solve FPCST, we first formulate LPCST with an additional parameter. Then we show how this enables us to solve FPCST using our algorithm for LPCST.

The connection between a parametric formulation and the fractional version of the same problem has already been established by Dinkelbach [3].

Let \mathcal{T} be the set of all connected subgraphs $T = (V', E')$ of G that contain the root. We are looking for a graph in \mathcal{T} that maximizes the expression

$$\frac{\sum_{v \in V'} p(v)}{c_0 + \sum_{e \in E'} c(e)}.$$

Now consider the following function $o(t)$:

$$o : \mathbb{R}^+ \rightarrow \mathbb{R}, o(t) = \max_{T=(V',E') \in \mathcal{T}} \sum_{v \in V'} p(v) - t(c_0 + \sum_{e \in E'} c(e)).$$

Let t^* be the value of the optimal solution of FPCST on G and $t \in \mathbb{R}$. Then we have:

$$o(t) = 0 \Leftrightarrow t = t^*, \quad o(t) < 0 \Leftrightarrow t > t^*, \quad o(t) > 0 \Leftrightarrow t < t^* .$$

Using the algorithm for LPCST, we can test for any t in linear time if it is smaller, equal, or greater than the optimal solution for FPCST. This fact can be used to construct different search algorithms that solve the problem.

There is also a geometric interpretation of our problem. Let \mathcal{T} be again the set of all non-empty subtrees of G . Each $T = (V_T, E_T) \in \mathcal{T}$ defines a linear function $f_T : \mathbb{R}^+ \rightarrow \mathbb{R}$ in the following way:

$$f_T(t) = \sum_{v \in V_T} p(v) - t(c_0 + \sum_{e \in E_T} c(e)) .$$

Since all vertex profits and edge costs are non-negative, and c_0 is positive, all these linear functions have negative slope. In this geometric interpretation, the function o defined above is the maximum of these functions. Hence it is a piecewise linear, convex, monotonously decreasing function. What we are looking for is the point where o crosses the x -axis. The functions f_T that contain this point correspond to optimal subtrees for the given profits and costs.

3.1 Binary Search

An easy way of building an algorithm for the FPCST problem that uses the parametric formulation of the previous section is binary search. We start with an interval (t_l, t_h) that contains t^* . Then we test the mid point t of this interval using the algorithm for the linear problem. This will give us either a proof that t equals t^* or a new upper or lower bound and will halve the size of the interval.

It is important to choose the right terminating conditions to achieve good performance. In our case, these conditions rely on the fact that $o(t)$ is the maximum of linear functions (see [10] for details). Since the running time of the algorithm depends to a great degree on the values for the profits and costs, a meaningful upper bound for the worst case running time that depends only on the size of the input graph cannot be given.

3.2 Newton’s Method

We use the adaptation of Newton’s iterative method described for example by Radzik [12]. Let \mathcal{T} be the set of all subtrees of G that contain the root. We start with $t_0 = 0$. In iteration i , we compute

$$o(t_i) = \max_{T=(V',E') \in \mathcal{T}} \sum_{v \in V'} p(v) - t_i(c_0 + \sum_{e \in E'} c(e))$$

together with the optimal tree $T_i = (V_i, E_i)$ for parameter t_i using the linear algorithm from Section 2. As long as $o(t_i)$ is greater than 0, we compute t_{i+1} as the fractional objective value of T_i . So we have:

$$t_{i+1} = \frac{\sum_{v \in V_i} p(v)}{c_0 + \sum_{e \in E_i} c(e)} .$$

In the course of this algorithm, t_i increases monotonically until t^* is reached. Let l be the index with $t_l = t^*$. Radzik shows in [13] for general fractional optimization problems where all weights are non-negative that $l = O(p^2 \log^2 p)$ where p is the size of the problem (in our case the number of vertices of the problem graph G).

For our specific problem, we can prove a stronger bound for l :

Theorem 1. *Newton’s method applied to the fractional prize-collecting Steiner tree problem with fixed costs takes at most $n + 2$ iterations where n is the number of vertices of the input tree T .*

To proof the theorem, we show that for each iteration of Newton’s method on our problem, there is an edge that was contained in the previous solution but is not contained in the current solution. This implies that the number of iterations is linear (see [10] for a detailed proof).

Since we can solve the problem for the linear objective function in linear time using the algorithm from Section 2, Newton’s Method has a worst case running time of $O(|V|^2)$ for our problem.

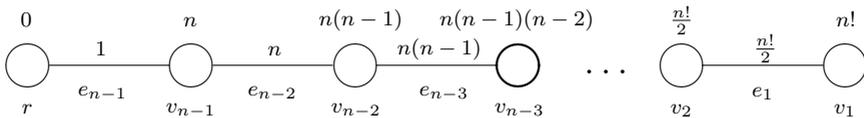


Fig. 1. Worst case example for Newton’s Method. The edge costs and vertex profits are above the path while the names of the vertices and edges are below

Figure 1 shows an example where this worst case running time is reached. If we define the fixed costs $c_0 = 1$, we can show by a coarse estimation of the objective function value for each path starting at r that the solution of Newton’s method shrinks only by one vertex in every iteration and that the

optimal solution is the root together with vertex v_{n-1} . Therefore, the algorithm executes $n - 1$ iterations and since each iteration has linear running time, the total running time of Newton's method on this example is $\Theta(n^2)$.

3.3 A New Algorithm Based on Megiddo's Parametric Search

In this section, we present our new algorithm for the FPCST problem which is a variant of parametric search introduced by Megiddo [11]. Furthermore, we suggest an improvement that guarantees a worst case running time of $O(n \log n)$ for any tree G with n vertices.

The idea of the basic algorithm is to simulate the execution of the algorithm A for LPCST on the unknown edge cost parameter t^* (the objective value of an optimal solution). During the simulation, we keep an interval (t_l, t_h) that contains t^* and that is initialized to $(0, \infty)$. Whenever A has to decide if a certain edge (u, v) is included in the solution, this decision is based on the evaluation of the maximum in (1) and depends on the root r_d of a linear function in t given by $l(u) - t \cdot c(u, v)$.

The decision is clear if r_d is outside (t_l, t_h) . Otherwise, we multiply all edge costs of the tree with r_d and execute A on the resulting problem. The sign of the linear objective function value $o(r_d)$ determines the decision (which enables us to continue the simulation of A) and r_d either becomes the new upper or lower bound of (t_l, t_h) .

There are two possibilities for the algorithm to terminate. The first is that one of the roots we test is t^* . In this case, we can stop without completing the simulation of A . If we have to simulate A completely, we end up with an interval for t^* . In this case, we perform depth first search on the edges that we have not cut during the simulation to obtain an optimal subtree.

Just as in the algorithm for the linear problem, our algorithm assigns labels to the vertices, but these labels are now linear functions that depend on the parameter t . The algorithm uses a copy G' of the problem tree G . In each phase, all leaves of G' are deleted after the necessary information has been propagated to the parents of the leaves. When the algorithm starts, the label of every vertex is set to the constant function equal to its profit. In the course of the algorithm, these labels change and will correspond to linear functions over the parameter t .

When we look at a certain leaf v with label $f_v(t)$ during a phase we compute the linear function $\tilde{f}_v(t) = f_v(t) - t \cdot c(e_v)$ where e_v is the edge incident to v . Let r_v be the root of $\tilde{f}_v(t)$. For all current leaves, we collect the values r_v , sort them and perform binary search on the roots using the linear algorithm to decide if the value t^* is smaller, greater, or equal than a certain root. Note that we do not have to include the roots in the binary search that are outside the current interval for t^* . If there are roots that are inside the current interval, we either find t^* or we end up with a smaller interval.

After the binary search, we know for each leaf v if its root r_v is smaller or greater than t^* (if it is equal, we have already found the solution and the algorithm has stopped). We delete all leaves whose root is smaller than t^* from G' . For all other leaves v , we add the function $\tilde{f}_v(t)$ to the label of its parent

and delete v , too. Now the next phase of the algorithm starts with the vertices that have become leaves because of the deletion of the current leaves (see [10] for a pseudo code.

The correctness of the algorithm follows from the general principle of Megiddo’s method [11]. The running time of the algorithm is dominated by the calls to the linear algorithm. The binary search is performed by solving $O(\log(|B|))$ instances of LPCST with profits and costs determined by the parameter t . The set B is the set of leafs of the current working graph G' . Since it may happen that the graph contains only one leaf in every iteration (G may be a path) the number of iterations can be n . The worst case example for Newton’s method in Section 3.2 is also a worst case example for this algorithm. Thus the overall running time of the algorithm is $O(|V|^2)$.

Improvement Through Path Contraction. If there is no vertex in G with degree two, our algorithm already has a running time of $O(n \log n)$ for a tree with n vertices: In this case we delete at least half the vertices of the graph in every iteration by deleting all leaves. It will follow from the proof of Theorem 2 that this property is sufficient for the improved running time.

We will remove the remaining obstacles in the graph, namely vertices of degree two, by performing a reduction of all paths in the tree. This must be done in every iteration since the removal of all leaves at the end of the previous iteration may generate new paths. The idea of the reduction is based on the fact that the subtree situated at the end of a path can only contribute to the optimal solution if the complete path is also included. Otherwise, only a connected subset of the path can be in the optimal solution.

More formally, a subset of V is a path denoted by $P := \{v_0, v_1, \dots, v_m, v_{m+1}\}$ if v_0 has degree greater two or is the root, v_{m+1} does not have degree two and all other vertices are of degree two. To fix the orientation we assume that v_0 is included in the path from v_1 to r . Since we want to contract the m vertices of the path to a single vertex, trivial cases can be excluded by assuming $m \geq 2$. In an optimal solution either there exists a vertex $v_q \in P$ such that v_1, \dots, v_q are the only vertices of P in the solution, or P is completely contained in the solution and connects a possible subtree rooted at v_{m+1} to r .

The procedure CONTRACTPATH (see Algorithm 1) determines the best possible candidate for v_q and contracts the path by adding an artificial edge from v_0 to v_q with cost equal to the value of the complete subpath including v_1, \dots, v_{q-1} , and a second artificial edge from v_q to v_{m+1} that models the cost of traversing the vertices v_{q+1}, \dots, v_m . The path contraction is invoked at the beginning of every iteration in our algorithm for FPCST.

The main theoretical result of this paper is stated in the following theorem:

Theorem 2. *The running time of Algorithm the algorithm with CONTRACTPATH is in $O(n \log n)$.*

Proof. (Sketch) To find v_q , we need to compute the maximum of m linear functions, which can be done in time $O(m \log m)$ (see [2] for a proof). The resulting

Data : A labeled tree $T = (V, E)$ with fixed root r ;
a path in T $v_0, v_1, \dots, v_m, v_{m+1}$, $m > 2$

Result : A labeled tree $T = (V, E)$ with fixed root r

$end[1] = 0$;

for $j = 1$ *to* m **do**
 $end[j] := end[j - 1] + l(v_j) + c(v_{j-1}, v_j)$;

end

$f(t) = \max_{j=1}^m end[j]$;

$B = \{t \in (t_l, t_h) \mid t \text{ is breakpoint of } f(t)\} \cup \{t_l, t_h\}$;

Perform binary search on B using the modified linear algorithm and update t_l and t_h ;

choose q s.t. $end[q] = f(t)$ for $t \in (t_l, t_h)$;

$c(v_0, v_q) := \sum_{k=1}^{q-1} (l(v_k) + c(v_{k-1}, v_k)) + c(v_{q-1}, v_q)$;

$c(v_q, v_{m+1}) = \sum_{k=q+1}^m (l(v_k) + c(v_{k-1}, v_k)) + c(v_m, v_{m+1})$;

Remove vertices $v_1, \dots, v_{q-1}, v_{q+1}, \dots, v_m$ from T ;

Algorithm 1: Algorithm CONTRACTPATH to remove all nontrivial paths from a tree

piecewise linear function has at most m breakpoints. In every iteration there is a number of breakpoints from CONTRACTPATH and a number of leaves with corresponding root values to be considered. We use binary search in each iteration to find a new interval (t_l, t_h) including neither breakpoints nor roots thus resolving the selection of v_q and the final decision on all leaves.

If k is the size of the graph at the beginning of an iteration, then the binary search performs a logarithmic number of calls to the algorithm that solves LPCST. Therefore, a single iteration takes time $O(k \log k)$. It can be shown that applying the procedure CONTRACTPATH to every non trivial path guarantees that our algorithm together with CONTRACTPATH deletes at least one third of the vertices in each iteration. Since the size of the graph is reduced by a constant fraction after each iteration, the total running time sums up to $O(n \log n)$. See [10] for a detailed proof. □

4 Computational Experiments

We generated two different test sets of graphs to test the performance of the algorithms presented in Section 3. The first set consists of randomly generated trees where every vertex has at most two children while the second set contains random trees where each vertex can have up to ten children. In both sets, the cost of each edge and the profit of each vertex is a random integer from the set $\{1, 2, \dots, 10,000\}$. Both sets contain 100 trees for each number of vertices from 1,000 to 10,000 in steps of 500 vertices. The fixed costs for all problem instances has been chosen as 1,000 times the number of vertices in the graph. This produces solutions containing around 50% of all vertices for the graphs where each vertex has at most 10 children. For the graphs where each vertex

has at most two children, the percentage is around 35%. To execute the three algorithms on the test sets as a documented and repeatable experiment and for analyzing the results, we used the tool set ExpLab [9].

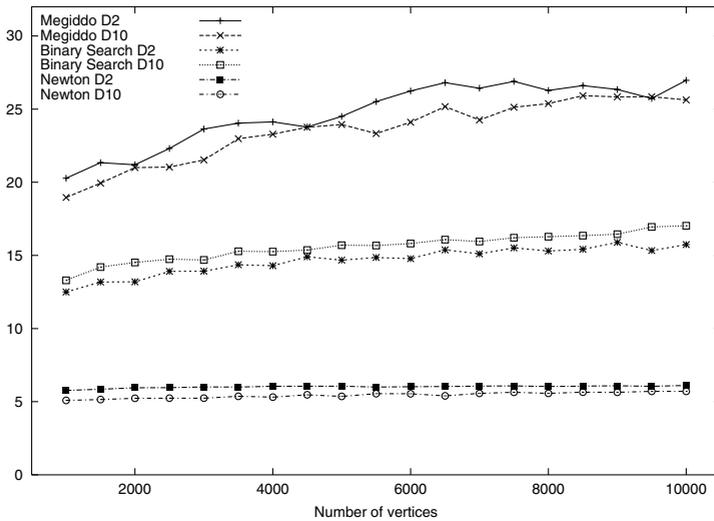


Fig. 2. The average number of calls to the linear algorithm executed by the three algorithms on the benchmark set with maximum degree 2 and maximum degree 10

Figure 2 shows the average number of calls over all trees with the same number of vertices for the three algorithms and the two benchmark sets. The number of calls grows very slowly with the size of the graphs for all three algorithms. In fact, the number of calls barely grows with the number of vertices in the graph for Newton’s method.

Our variant of Megiddo’s method needs more calls than the other two methods. For the leaves of the tree, the algorithm behaves just like binary search. The reason why the number of calls is higher than for binary search is that our new algorithm not only executes calls at the leaf level but also higher up in the tree. These are usually very few and not on every level. So on a level where additional calls have to be made, there are usually only one or two open decisions. Therefore, the binary search in our new algorithm can not effectively be used except at the leaf level. Because of this fact, the pure binary search algorithm can “jump” over some decisions that parametric search has to make on higher levels.

The reason why Newton’s method needs fewer calls than the binary search method is the random nature of our problem instances. Binary search starts with a provable upper bound for t^* which in our case is the sum of all vertex profits divided by the fixed costs. This upper bound is far away from the objective value of the optimal solution. After the first iteration of Newton’s method, the

value t is the objective function value of the whole tree. This value is a good lower bound for the optimal solution because the profits and costs are random and with the fixed costs we have chosen, the optimal tree contained 35-50% of all vertices. Therefore, Newton's method needs only a small number of steps to reach the optimal solution and the number of calls grows only very slowly with the size of the graphs.

Figure 3 shows that the number of calls to the linear algorithm determines the running time: our new algorithm is the slowest and Newton's method the fastest. The running times grow slightly faster than linear with the size of the graphs. Since each call to the algorithm for the linear problem needs linear time, the fact that the number of calls grows with the size of the graph (albeit very slowly) is the reason for this behavior. We executed the experiments on a PC with a 2.8 GHz Intel Processor with 2GB of memory running Linux. Even for the graphs with 10,000 vertices, the problems can be solved in less than 1.8 seconds.

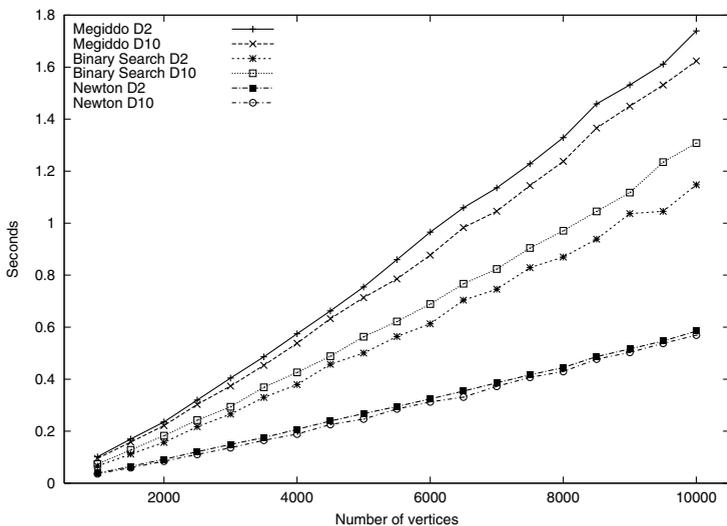


Fig. 3. The average time used by the three algorithms on the two benchmark sets

We also executed an experiment where we used only the 100 graphs of the test set with maximum degree 10 that have 10,000 vertices. We increased the fixed costs c_0 exponentially and ran all three algorithms on the 100 graphs for each value of c_0 . We started with $c_0 = 100$ (where the solution contained only a few vertices) and multiplied the fixed costs by 10 until we arrived at 10^{11} (where the optimal solution consisted almost always of the whole tree).

Figure 4 shows how the time needed by the three algorithms depends on fixed costs. It is remarkable that for small fixed costs, binary search is faster than Newton's method but for fixed costs of more than 10,000, Newton's method is

faster. The reason is the same we have already given for the better performance of Newton’s method in our first experiments. For large fixed costs, the percentage of the vertices contained in an optimal solution rises and so the value of the first solution that Newton’s method tests, which is the value of the whole graph, is already very close to the optimal value. Binary search has to approach the optimum solution from the provable upper bound for the objective function value which is far away from the optimal solution when this solution is large and therefore contains many edges.

Parametric search is not much slower than binary search for high fixed costs. As the plot shows, the reason is not that parametric search performs significantly better for higher fixed costs but that the performance of binary search deteriorates for the reasons given in the last paragraph.

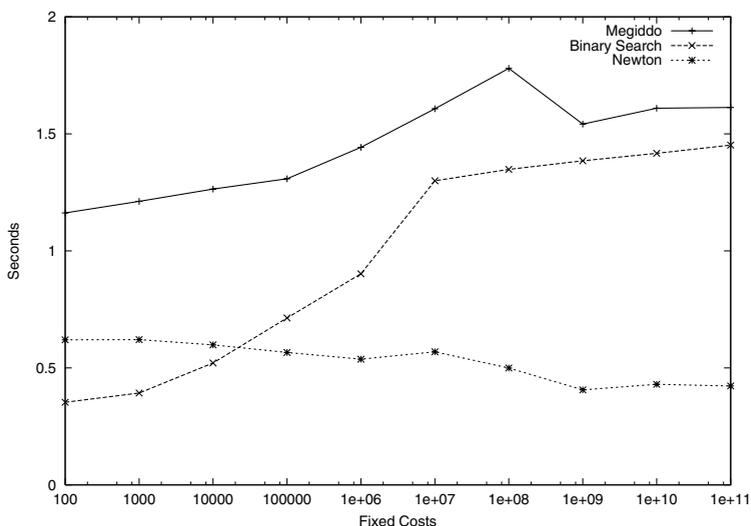


Fig. 4. Time used by the three algorithms for growing fixed costs (logarithmic x-axis)

5 Conclusions

In this paper, we have presented three algorithms for solving the fractional prize-collecting Steiner tree problem (PCST problem) on trees $G = (V, E)$. We have shown that Newton’s algorithm has a worst case running time of $O(|V|^2)$. We have also presented a variant of parametric search and proved that the worst case running time of this new algorithm is $O(|V| \log |V|)$. Our computational results show that Newton’s method performs best on randomly generated problems while a simple binary search approach and our new method are considerably slower. For all three algorithms, the running time grows slightly faster than linear with the size of our test instances.

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