Paper Report

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Abstract

This is a report on the paper "A fast and simple algorithm for the maximum flow problem" by James B. Orlin and R. K. Ahuja (1989). This paper was presented as a part of a course on Matching and Flow Algorithms at the Chennai Mathematical Institute.

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0.1 Overview and Notation

The paper's main result is a new algorithm for computing the maximum flow in a flow network, motivated by the celebrated *push-relabel* algorithm for the same problem. The authors first review the push-relabel algorithm for the maximum flow problem; then they identify the bottleneck in the algorithm and try to find a way around it. In this section, we will introduce the relevant definitions and notation.

0.1.1 A review of flows and preflows. Throughout this report, we will be working with a directed network G = (N, A), where N is the set of nodes, and A the set of arcs in the network. For every arc $(i, j) \in A$, we will use the symbol u_{ij} to denote it's (integral) capacity. Also, we define n := |N| and m := |A|. The source s and sink t will be two distinguished nodes in the network. We will assume that if $(i, j) \in A$, then $(j, i) \in A$ (with (j, i) possibly having zero capacity).

We will use the symbol U to denote the *maximum* capacity of any arc emanating from the source vertex s, i.e

$$U := \max_{(s,j) \in A} \left\{ u_{sj} \right\}$$

A flow is a function $f: A \to \mathbb{R}$ that satisfies the following.

• For all $i \in N - \{s, t\}$, we have

$$\sum_{\{j:(j,i)\in A\}} f_{ji} - \sum_{\{j:(i,j)\in A\}} f_{ij} = 0$$

This is also called the *flow conservation law*.

• For each arc $(i, j) \in A$, we must have

$$0 \le f_{ij} \le u_{ij}$$

This is called the *capacity constraint*.

The maximum flow problem tries to compute a flow f such that the following quantity, called the *size* of the flow f, is maximized.

$$\sum_{\{j:(j,t)\in A\}} f_{jt}$$

Next, we quickly discuss *preflows*. A *preflow* f is a function $f : A \to \mathbb{R}$ that satisfies all capacity constraints, and satisfies the following relaxation of the flow conservation law for all $i \in N - \{s, t\}$.

$$\sum_{\{j:(j,i)\in A\}} f_{ji} - \sum_{\{j:(i,j)\in A\}} f_{ij} \ge 0$$

For a node i, the above quantity is called the *excess* at node i, and is denoted by the symbol e_i . So, for a preflow, all the excesses must be non-negative.

A node $i \in N - \{s, t\}$ with positive excess is said to be *active*. The *residual capacity* of an arc $(i, j) \in A$, with respect to a preflow f, is defined as

$$r_{ij} = u_{ij} - f_{ij} + f_{ji}$$

The *residual network* for a preflow f is defined to be the network containing only those edges which have a positive residual capacity.

Finally, for each vertex i, we define the arc adjacency list A(i) as the set

$$A(i) := \{ (i, k) \in A : k \in N \}$$

0.1.2 Valid distance functions. A valid distance function is a function $d: N \to \mathbb{Z}^+$ for a preflow f that satisfies the following two conditions.

• d(t) = 0.

• $d(i) \le d(j) + 1$ for every arc $(i, j) \in A$ with $r_{ij} > 0$.

An arc $(i, j) \in A$ in the residual network is called *admissible* if d(i) = d(j) + 1; otherwise it's called an *inadmissible arc*.

0.2 The PUSH-RELABEL Algorithm

In this section, we will review the PUSH-RELABEL algorithm for the max flow problem. The algorithm consists of the following three subroutines.

- (1) PREPROCESS. In this subroutine, we initialize a preflow and a valid distance function. For each arc $(s, j) \in A(s)$, send u_{sj} units of flow. Also, initialize d(s) = n, d(t) = 0 and d(i) = 1 for all $i \neq s, t$. It can be easily verified that this is a valid distance function.
- (2) PUSH(i). Here, we select an admissible arc (i, j) in A(i), and send $\delta = \min \{e_i, r_{ij}\}$ units of flow from node i to j.

A push of flow on arc (i, j) is said to be *saturating* if $\delta = r_{ij}$, and non-saturating otherwise.

(3) RELABEL(i). Here, we just replace d(i) by min $\{d(j) + 1 : (i, j) \in A(i) \text{ and } r_{ij} > 0\}$. This is called a *relabel step*.

The PUSH-RELABEI algorithm can be described by the following pseudocode.

Algorithm 1 PUSH-RELABEL Algorithm

1: PREPROCESS.			
2: while there is an active node do			
3: Select an active node i .			
4: if there is an admissible arc in $A(i)$ then			
5: $PUSH(i)$.			
6: else			
7: $\operatorname{RELABEL}(i).$			
8: end if			
9: end while			

Proposition 0.1. The PUSH-RELABEL algorithm satisfies the following properties.

- (1) It maintains valid distance labels at each step. The distance labels only increase over the course of the algorithm, and at each step, the distance label of some node strictly increases.
- (2) For each node $i \in N$, d(i) < 2n.
- (3) The number of relabel steps is less than $2n^2$.
- (4) The number of saturating pushes is atmost nm.
- (5) The number of non-saturating pushes is at most $2n^2m$.

From this proposition, one can easily conclude that the time complexity of the PUSH-RELABEL algorithm is $O(nm + n^2m)$.

0.3 The bottleneck, and the Excess Scaling Algorithm

The authors remark that the bottleneck operation in many preflow-based algorithms is the number of non-saturating pushes. Intuitively, each saturating push changes the structure of the residual network (by removing an edge from the network). However, a non-saturating push doesn't change this structure, and hence it becomes difficult to bound the total number of non-saturating pushes. The main result of the paper, namely the *Excess Scaling Algorithm*, relies on a good upper bound on the number of non-saturating pushes.

The authors have shown that their algorithm reduces the number of non-saturating pushes from $O(n^2m)$ to $O(n^2 \log U)$. In the upcoming sections, we will describe the algorithm.

0.3.1 Description of the Excess Scaling Algorithm. The algorithm consists of a number of iterations, called *scaling iterations*. Each scaling iteration roughly does the following.

- (1) For each scaling iteration, we define the *excess dominator* to be the least integer Δ that is a power of 2 and which satisfies $e_i \leq \Delta$ for all $i \in V$.
- (2) After each scaling iteration, we ensure that Δ decreases by a factor of 2.
- (3) Note that when we initialize our preflow in the PREPROCESS subroutine, the maximum possible excess on any vertex is precisely $\max_{(s,j)\in A} \{u_{sj}\}$, which we denoted by U. So naturally, we will start off with $\Delta = 2^{\lceil \log U \rceil}$.
- (4) Since each scaling iteration decreases Δ by a factor of 2, we see that there will be $1 + \lceil \log U \rceil$ scaling iterations.

Suppose Δ is the excess dominator at start of a scaling iteration. We will somehow guarantee that each *non-saturating* push during this pushes at least $\Delta/2$ units of flow; this will help us in bounding the number of non-saturating pushes better. Moreover, we also have to ensure that during a scaling iteration, the excess dominator *never increases*.

To ensure that each non-saturating push has a value of at least $\Delta/2$, we do the following.

- (1) In a scaling iteration, we only push flow from vertices with an excess more than $\Delta/2$.
- (2) Moreover, among all vertices with excess more than $\Delta/2$, we push flow from the vertex with the minimum distance label. Since we push flow on only admissible edges, this choice will ensure that flow is being pushed to a vertex with excess at most $\Delta/2$.
- Next, we describe the data structures that are maintained throughout the algorithm.
 - (1) For each $r \in \{1, 2, ..., 2n 1\}$, we maintain a list denoted by LIST(r); this list will just be the set

$$\left\{i\in V: e_i > \frac{\delta}{2}, d(i) = r\right\}$$

Note that since d is a valid distance function, d(i) < 2n for all i at all times in the algorithm. Hence, we only need to maintain these lists for values of r atmost 2n-1. In practice, LIST(r) for each r will be implemented as a linked list, in which we can add and remove elements in O(1) time.

- (2) We will maintain a variable called *level*, which will represent the smallest index r for which LIST(r) is non-empty.
- (3) For each vertex i, an edge adjacency list A(i) will be maintained. Moreover, for each i, we will maintain a *current edge*, which will be an edge in A(i) which is a potential candidate for pushing flow out of i.

0.3.2 Pseudocode. The algorithm can be described precisely by the given pseudocode. It consists of the subroutine PUSH/RELABEL(i), for which we've also provided the pseudocode.

Algorithm 2 Excess S	Scaling A	Algorithm
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8				
1: PREPROCESS.				
$2: K := 1 + \lceil \log U \rceil$	\triangleright The number of scaling iterations			
3: for $k = 1$ to K do				
4: $\Delta = 2^{K-k}$	\triangleright The excess dominator			
5: for each vertex i do				
6: if $e_i > \Delta/2$ then				
7: Add i to $\text{LIST}(d(i))$				
8: end if				
9: end for				
10: level := 1.				
11: while $level < 2n$ do				
12: If $\text{LIST}(level) = \phi$, then $level := level + 1$.				
13: Otherwise, select a vertex i from LIST(le	vel), and do $PUSH/RELABEL(i)$.			
14: end while				
15: end for				

Algorithm 3 PUSH/RELABEL(i)

- Starting from the current edge of i, find an admissible edge (i, j) ∈ A(i) with r_{ij} > 0, incrementing the current edge pointer if necessary.
 if an admissible edge (i, j) has been found them.
- 2: if an admissible edge (i, j) has been found then
- 3: Push min $\{e_i, r_{ij}, \Delta e_j\}$ units of flow on arc (i, j).
- 4: Update residual capacity r_{ij} and excesses e_i and e_j .
- 5: If (updated) $e_i \leq \Delta/2$, delete *i* from LIST(*d*(*i*)).
- 6: If $j \neq s, t$ and if the (updated) $e_j > \Delta/2$, then add node j to LIST(d(j)); and set level := level 1.
- 7: **else**

 \triangleright No admissible edge found in the previous step

- 8: Delete *i* from LIST(d(i)).
- 9: Update label d(i) as usual.
- 10: Add i to LIST(d(i)), and set the *current edge* of i to the first edge of A(i).

11: **end if**

Lines 11-13 of the main algorithm do the following: among all vertices with excess more than $\Delta/2$, select the vertex with the minimum distance label, and try to push flow out of it; if not, relabel it. Line 3 of the PUSH/RELABEL(*i*) subroutine makes sure that, if the push is non-saturating, atleast $\Delta/2$ units of flow is pushed. This is quite easy to see: if the push is non-saturating, then the minimum of the three quantities is actually the minimum of $\{e_i, \Delta - e_j\}$; we know that $e_i > \Delta/2$ and $e_j \leq \Delta/2$ (since (i, j) is an admissible edge, and *i* is the vertex with minimum distance label among all vertices with excess more than $\Delta/2$); hence, the minimum of these two quantities is atleast $\Delta/2$.

Proposition 0.2. The Excess scaling algorithm satisfies the following.

- (1) Each non-saturating push from vertex i to vertex j sends at least $\Delta/2$ units of flow.
- (2) No excess ever increases above Δ withing a scaling iteration.
- (3) The number of non-saturating pushes per scaling iteration is at most $8n^2$.

The above three claims are not hard to prove. The third claim is proved by defining a potential function F for a scaling iteration by

$$F = \sum_{i \in V} \frac{e_i d(i)}{\Delta}$$

The idea is to track how much F increases/decreases over a scaling iteration. It can be shown that whenever a push is done, F decreases; using this along with the fact that each non-saturating push has a value of atleast $\Delta/2$, it can be shown that the number of non-saturating pushes is at most $8n^2$.

0.3.3 Complexity of the algorithm. Finally, with a little bit of additional work and using the facts mentioned in the previous section, one can show that the complexity of the Excess Scaling algorithm is $O(nm + n^2 \log U)$. Here, the term nm comes from an upper bound on the number of saturating pushes, and the term $n^2 \log U$ comes from the upper bound on the number of non-saturating pushes. As before, the number of relabel operations is again $O(n^2)$.