

Combinatorial Optimization

Jacob Denson

November 30, 2016

Table Of Contents

1	Matchings and Flows	2
1.1	Bipartite Matching	2
1.2	Vertex Covers	4
1.3	The Hopcroft-Karp Matching Algorithm	5
1.4	Network Flow	6
1.5	Goldberg-Tarjan Push Relabel Algorithm	10
1.6	Cuts in Undirected Graphs	11
1.7	Potentials	14
1.8	Minimum Cost Flows	15
1.9	Minimal Cost Bipartite Matching	16
1.10	Minimum Mean Cycle Cancelling	17
1.11	Minimum Spanning Trees and Arborescences	21
1.12	Arborescences	23
2	Matroids	26
2.1	Linear Programs for Matroids	28
3	Submodular Functions	32
4	Perfect Graphs	33

Chapter 1

Matchings and Flows

1.1 Bipartite Matching

Let $G = (V, E)$ be a graph. A **matching** is $M \subset E$ such that no vertex in V is the endpoint of more than one edge in M . M is maximal if $|M| \geq |M'|$ for any other matching M' . The maximal matching problem asks us to find a fast algorithm to find a maximum matching in any graph.

There is a polynomial time algorithm which can find matchings on any graph, but the problem is much more simple if G is bipartite – that is, if we may partition V into the disjoint union $W \sqcup U$ of two sets of vertices, such that every edge in E contains a point in W and a point in U . The bipartite matching asks us to find a maximal matching in a bipartite graph.

Suppose we are given a particular matching in a bipartite graph. Is there a reliable procedure to improve the matching? We could proceed by a guess and check method – we remove an edge in our matching, then try and add an edge using one of the vertices which has been freed up. If this edge cannot be taken because the end point is attached to an edge still in the matching, we remove that edge, freeing up more vertices. If we ever end up adding more edges than we started with (which occurs when we add an edge not attached to any points in the current matching), then we find a matching with an extra edge than before. This process is formalized by the ‘augmenting paths’ construction.

Given a particular matching M , construct a directed graph $G_M = (V \cup \{s, t\}, E_M)$, where s and t are new vertices. Let $w \in W, u \in U$. Construct the edges E_M such that

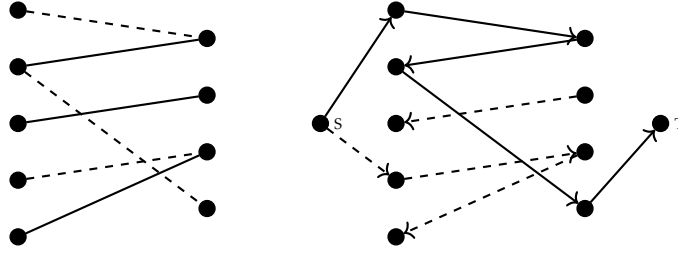


Figure 1.1: The left graph shows a matching in a graph of cardinality 3. The solid lines correspond to the edges chosen. On the right is the augmenting path graph corresponding to the matching. It contains a path from s to t , hence there is a way to improve the matching to a match of cardinality 4.

- $wu \in E_M$, if $wu \in M^c$.
- $uw \in E_M$, if $uw \in M$.
- $sw \in E_M$, if w is not the endpoint of any edge in M .
- $ut \in E_M$, if u is not the endpoint of any edge in M .

Let $(s, w_1, u_1, \dots, w_n, u_n, t)$ be a path in G_M from s to t , with $w_i \in W$, $u_i \in U$. Then $w_i u_i \in M^c$ and $u_i w_{i+1} \in M$, and w and u are both unused in M . Let M' be obtained from M by removing all edges of the form $u_i w_{i+1}$, and adding all edges of the form $w_i u_i$. It is easy to convince yourself that M' is a matching, with one more edge than M .

Conversely, suppose that there is no path from s to t in G_M . We claim M is then a maximal matching on G . Otherwise, we would have a matching M^* with $|M^*| > |M|$. Consider the multigraph $H = (V, M \sqcup M^*)$. Every vertex in V has at most degree two (for it can only be the endpoint of a single edge in M , and a single endpoint in M^*), and thus V breaks into cycles and paths. For any cycle C , we find $|M \cap C| = |M^* \cap C|$, for the edges in the cycle must alternate being in M and M^* . Since $|M^*| > |M|$, there must be a path P in H containing more edges in M^* than in M . P must therefore begin and end with a path in M^* , and we see that this path is exactly an augmenting path for M .

This argument justifies the correctness of the following matching algorithm, a variant of the ‘Ford Fulkerson Algorithm’. We take any initial

matching M . Then, we construct the augmenting path graph G_M , and use any of the standard algorithms to find a path from s to t . If we have $2n$ nodes, and m edges, then we can find a path in $O(m)$ time, and we can bound the number of augmentations we require to $O(n)$, so the algorithm runs in $O(nm)$ time.

1.2 Vertex Covers

A vertex cover $C \subset V$ is a set of points containing at least one endpoint of every edge in E . Clearly if M is a maximal matching in G , then $|M| \leq |C|$, because there is an injective function $f : M \rightarrow C$, obtained by mapping an edge to one of its endpoints which lies in C .

Theorem 1.1 (König - Egerváry). *Let $G = (W \sqcup U, E)$ be bipartite, and let M be a maximal matching in G . Let $Z \subset V \cup \{s, t\}$ be the nodes reachable from s in the augmenting path graph G_M . Then $C = (W - Z) \cup (U \cap Z)$ is a cover with $|C| = |M|$.*

Proof. First we prove C is a cover. Let $wu \in E$ be arbitrary.

- Suppose that $wu \in M$. If $u \in Z$, then $w \in Z$, because uw is an edge in G_M , so that if $w \notin C$, $u \in C$.
- Suppose that $wu \notin M$. Then wu is an edge in G_M , so that if $w \in Z$, $u \in Z$, and therefore if $w \notin C$, $u \in C$.

Now we show $|U| = |M|$. First, note that C contains *exactly one* of each of the endpoints of M , so $|U| \geq |M|$. Conversely, let $wu \in E - M$ be arbitrary. If $u \in C$, then u is reachable from s , and hence u is the endpoint of an edge in M , for otherwise t is reachable from s through u . Conversely, if $w \in C$, then w is not reachable from s , hence w is on an edge in M (for otherwise there is an edge directly to w from s in G_M). This shows that every vertex in C is on one of the edge in M , and hence $|U| = |M|$. \square

This gives us a method of finding a min cover in a graph. One equivalent form of this theorem is very useful. Given an undirected graph $G = (V, E)$, let N be the neighbour function $N(S) = \{v \in V : sv \in E, s \in S\}$.

Theorem 1.2 (Hall). *Let $G = (W \sqcup U, E)$ be a bipartite graph. Then G has a matching M which covers W if and only if $|N(S)| \geq |S|$ for all $S \subset W$.*

Proof. Certainly if a matching M covers W , then it induces an injective function from S to $N(S)$ for each $S \subset W$. Conversely, assume no matching of M covers W . Then König's theorem tells us there is a cover C of $W \sqcup U$ containing less than $|W|$ nodes. Then write $A = C \cap W$, $B = C \cap U$. We find $N(W - A) \subset B$, so $|N(W - A)| \leq |B| < |W| - |A|$. \square

1.3 The Hopcroft-Karp Matching Algorithm

The augmenting path approach to finding a maximal matching in a Bipartite graph improves the length of a candidate matching by one for each cycle of the algorithm. If we were able to consistently find augmentations which improved matchings by more than one edge, we could find a faster algorithm for maximal matching.

If M is a matching, and N_M be the length of a shortest alternating path in G_M , we shall let a collection of vertex disjoint alternating paths P_1, \dots, P_n be called **M-blocking** if $|P_i| = N_M$, and every alternating path of length N_M shares a vertex with one of the P_i . The next lemma is a simple generalization of the augmenting paths approach.

Theorem 1.3. *If P_1, \dots, P_n are vertex disjoint M -alternating paths, then*

$$M' = (M - \cup P_i) \cup (\cup P_i - M)$$

is a matching with $|M'| = |M| + n$.

The Hopcroft-Karp algorithm proceeds as in Ford-Fulkerson, albeit improving matchings by taking M -blocking paths as augmentations, improving matchings in chunks. We shall discover a method to find an M -blocking in linear time, and also show the chunk method yields an asymptotic speedup in the number of cycles to find a maximal matching, so that Hopcroft-Karp is objectively faster than Ford-Fulkerson.

Lemma 1.4. *If M is a matching, and P_1, \dots, P_n an M -blocking set of paths. Then $M' = (M - \cup P_i) \cup (\cup P_i - M)$, and $N_{M'} \geq N_M + 1$.*

Proof. Let $Q = (v_1, \dots, v_k)$ be an alternating path in $G_{M'}$ with $|Q| \leq |N_M|$, where v_1 and v_k are M' exposed. Then v_1 and v_k are also M exposed, because switching from M to M' only removes exposed vertices, not adds them.

Let H be the directed graph with edges taken from all the edge lying on the P_i , and the directed edges of Q , except that we remove both copies of duplicate edges. The reverse vu of any edge uv in Q that is not an edge on some P_i must be on some P_j \square

1.4 Network Flow

Let $G = (V, E)$ be a directed graph with two identified vertices $s \neq t$. Let $\mu : E \rightarrow \mathbf{R}^+$ be a function measuring the ‘capacity’ of each edge in the graph (Thinking of the edges as if they were ‘pipes’ which can only carry a certain throughput). Define, for $U \subset V$,

$$\delta_{\text{out}}(U) = \{w \in V - U : vw \in E\} \quad \delta_{\text{in}}(U) = \{w \in V - U : wv \in E\}$$

A flow is a mapping $f : E \rightarrow \mathbf{R}^+$ such that $0 \leq f \leq \mu$, and for any $v \neq s, t$,

$$f(\delta_{\text{out}}(v)) = f(\delta_{\text{in}}(v)) \quad *$$

a relation known as the flow conservation law. The aim of the maximum flow problem is to find f such that the value function

$$\text{val}(f) = f(\delta_{\text{out}}(s)) - f(\delta_{\text{in}}(s))$$

is maximized. This is essentially the amount of flow which is created at f . It is also the amount of flow which is ‘absorbed’ at t , because

$$\begin{aligned} f(\delta_{\text{out}}(t)) - f(\delta_{\text{in}}(t)) &= f(\delta_{\text{out}}(t)) - f(\delta_{\text{in}}(t)) + \sum_{v \neq s, t} f(\delta_{\text{out}}(v)) - f(\delta_{\text{in}}(v)) \\ &= \sum_{v \neq s} \left(\sum_{vw \in E} f(vw) - \sum_{wv \in E} f(wv) \right) \\ &= \left(\sum_{\substack{v \neq s \\ vs \in E}} f(vs) + f(ss) \right) - \left(\sum_{\substack{v \neq s \\ sv \in E}} f(sv) + f(ss) \right) \\ &= -[\delta_{\text{out}}(f, s) - \delta_{\text{in}}(f, s)] \end{aligned}$$

Thus the maximum flow problem tells us a method of maximizing the amount of flow which gets to t .

*We extend functions $g : X \rightarrow \mathbf{R}$ to $g : 2^X \rightarrow \mathbf{R}$ by defining $g(A) = \sum_{x \in A} g(x)$

There is an interesting relation between flows on graphs, and another structure on these graphs known as a cut, which we will take full advantage of in finding solutions to the max flow problem. Define a (s, t) cut on a directed graph $G = (V, E)$ with vertices s and t , to be a partition of V into two sets, one containing s , and one containing t . It is however more simple to consider a cut to be a subset C of V containing s , but not t . If we have a capacity function μ , then we define the value of the cut to be $\text{val}(C) = \mu(\delta_{\text{out}}(C))$. The min cut problem is to find a cut of smallest value.

Lemma 1.5. *If f is a flow, and C is an (s, t) cut, then $\text{val}(f) \leq \mu(\delta_{\text{out}}(C))$.*

Proof. Since $t \notin C$, similar manipulations to the ones above show that

$$\begin{aligned} \text{val}(f) &= \sum_{v \in C} f(\delta_{\text{out}}(v)) - f(\delta_{\text{in}}(v)) \\ &= f(\delta_{\text{out}}(C)) - f(\delta_{\text{in}}(C)) \\ &\leq \mu(\delta_{\text{out}}(C)) \end{aligned}$$

Note that we obtain equality if $f(\delta_{\text{out}}(C)) = \mu(\delta_{\text{out}}(C))$ and $f(\delta_{\text{in}}(C)) = 0$, in which case f must be a maximal flow, and C a min cut. \square

As with maximal matchings, we attempt to find a maximal flow by finding ways of augmenting suboptimal flows to a maximal solution. Given a flow f , we construct the residual graph G_f , which has the same vertices as G , but whose edges are defined to be the set

$$E_f = \{uv \in E : f(uv) < \mu(uv)\} \cup \{vu : uv \in E, f(uv) > 0\}$$

We shall denote the element vu in the second set which makes up E_f by \overleftarrow{uv} , since we obtained it from an edge uv in G . Also define a capacity function

$$\mu_f(uv) = \mu(uv) - f(uv) \quad \mu_f(\overleftarrow{uv}) = f(uv)$$

Suppose we can find a simple path P from s to t in G_f , and define $\alpha = \min_{e \in P} \mu_f(e)$. Consider a new flow

$$f'(uv) = \begin{cases} f(uv) + \alpha : uv \in P \\ f(uv) - \alpha : \overleftarrow{uv} \in P \end{cases}$$

First we show that f' is a flow. By the choice of α , $0 \leq f' \leq \mu$. For each vertex $v \neq s, t$, we must show that flow conservation still holds. Let $e_1, e'_1, \dots, e_n, e'_n$ be the edges in G containing v as an endpoint obtained from P , by reversing edges of the form \overleftarrow{uv} , considered in the order they appear in P . We may write pair up the edges in this manner because, for every edge that enters v in P , there must be an edge that leaves v in P . Let

$$S_i = \begin{cases} f(e_i) - f'(e_i) & e_i = uv \\ f'(e_i) - f(e_i) & e_i = vu \end{cases}$$

Define S'_i similarly for e'_i . Then, since no edges are repeated,

$$\begin{aligned} f'(\delta_{\text{out}}(v)) - f'(\delta_{\text{in}}(v)) &= f(\delta_{\text{out}}(v)) - f(\delta_{\text{in}}(v)) \\ &\quad + \sum_{i=1}^n (S_i + S'_i) \end{aligned}$$

We now show $S_i + S'_i = 0$ for each i . This breaks into several cases.

- If $e_i = uv$, $e'_i = vw$, in which case $f'(e_i) = f(e_i) + \alpha$, $f'(e'_i) = f(e'_i) + \alpha$, and $S_i + S'_i = \alpha - \alpha = 0$.
- If $e_i = uv$, $e'_i = wv$, then e'_i was obtained from an edge of the form \overleftarrow{vw} in P , in which case $f'(e'_i) = f(e'_i) - \alpha$, and so $S_i + S'_i = -\alpha - (-\alpha) = 0$.
- If $e_i = vu$, $e'_i = wv$, then e_i and e'_i were obtained from reversed edges in P , and so $S_i + S'_i = (-\alpha) - (-\alpha) = 0$.
- If $e_i = vu$, $e'_i = vw$, then e_i was obtained from reversing edges, and so $S_i + S'_i = (-\alpha) + \alpha = 0$.

And it follows, because flow conservation holds at v for f , that it holds at v for f' as well. Finally, we find that $\text{val}(f') = \text{val}(f) + \alpha$, which can be shown by performing an analysis, similar to v above, for s , noting that s will have an extra edge at the beginning of the path, which causes the extra α . All that remains is to show this augmentation method yields a maximum flow in all cases, after enough iterations.

Let f be a flow for which G_f has no augmenting path. Let U be the nodes reachable from s in G_f , easily verified to be a cut.

$$\text{val}(f) = \mu(\delta_{\text{out}}(U))$$

proving maximality of the flow, and minimality of the cut U . To see this, let $uv \in \delta_{\text{out}}(U)$. Then $f(uv) = \mu(uv)$, for otherwise v would be reachable from s in G_f . Similarly, we must have $f(\delta_{\text{in}}(U)) = 0$, for if $vu \in \delta_{\text{in}}(U)$, and $f(vu) > 0$, then $\bar{v}u \in G_f$, and so v is reachable from s , a contradiction. We conclude that

$$\text{val}(f) = f(\delta_{\text{out}}(U)) - f(\delta_{\text{in}}(U)) = \mu(\delta_{\text{out}}(U))$$

which shows that our augmenting paths approach works. Notice that if all edge weights are integral, then the augmenting flows are always integral, and thus there exists a maximal flow with integral weights.

The Ford Fulkerson algorithm solves the max flow problem by repeatedly augmenting an initial flow. In principle, this approach is correct, but if you do not find a residual path in a smart way, this algorithm does not terminate in polynomial time for all inputs. If all edge weights are integers n_1, \dots, n_m , then each augmentation increases the value of the flow by at least the greatest common factor of the n_i . Since there is a bound on the size of a maximal flow, obtained by taking the trivial cut $\{s\}$, the algorithm will eventually terminate in time proportional to

$$O\left(\frac{n_1 + \dots + n_m}{\text{lcm}(n_1, \dots, n_m)}\right)$$

which is exponential in the bit complexity of the representation. We cannot do any better than this, as the graph in the figure above provides an example. If the edge weights are rational, similar results can be obtained by multiplying out denominators.

We achieve much better estimates if we use breadth first search to find our residual path – that is, we always take the shortest path (in length, not in weight) from s to t . But this, of course, requires a careful analysis.

Theorem 1.6 (Mader). *Let G be an undirected graph, and s, t vertices not directly connected by an edge. Then the maximal number of vertex-disjoint (s, t) paths is equal to the minimum cardinality of a set $B \subset V - \{s, t\}$ which contains a vertex in any path from s to t .*

Proof. Form a directed bipartite graph W , which consists of two copies of the vertices in G . If v is a vertex, then we shall separately denote its copies by v^1 and v^2 . If uv is an edge in G , attach edges u^2v^1 and v^2u^1 to W , and attach an edge v^1v^2 for each vertex v . Define a capacity function on w by

$$\mu(u^2v^1) = \mu(v^2u^1) = \infty \quad \mu(v^1v^2) = 1$$

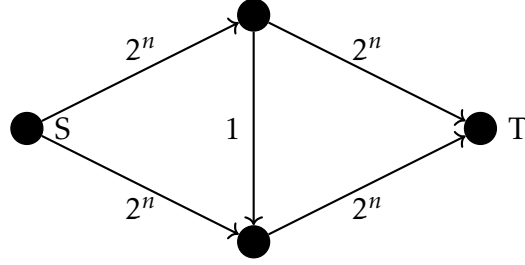


Figure 1.2: If our algorithm continuously switches between adding weights from left to right via the central edge, then we will need to compute $O(2^n)$ residual graphs before termination, even though the graph can be represented in $O(n)$ bits.

Note that any (s, t) path in G leads to an alternating path in W from s^2 to t^1 , which must cross at least one 1-capacity edge because there is no edge from s to t .

Let f be an integer-valued flow in W . Then $f \leq 1$, because an infinite capacity edge u^2v^1 satisfies $f(u^2v^1) \leq f(\delta_{\text{in}}(u^2)), f(\delta_{\text{out}}(v^1))$ \square

1.5 Goldberg-Tarjan Push Relabel Algorithm

Our previous algorithm found max flows in a graph by starting with a flow, then slowly augmenting it, improving the flow value until it is maximum. An alternative idea would be to take a flow-like function, which would be a maximum flow if it obeyed flow conservation, and slowly correcting the flow until we have a maximum flow. Define a **preflow** on a graph with capacity function μ to be a positive real-valued function f , defined on the edges, such that $f \leq \mu$, and $f(\delta_{\text{out}}(v)) - f(\delta_{\text{in}}(v)) \geq 0$ for all $v \neq s$. A preflow defines a residual graph G_f in the same way that a flow does. A non-negative integer valued function ψ on vertices is a **distance label** for a particular preflow f if $\psi(s) = |V|$, $\psi(t) = 0$, and $\psi(v) \leq \psi(w) + 1$ for $vw \in G_f$. Notice that if there is a path from s to t in G_f , then a distance label cannot exist, because if it has length k , then the inequality property gives

$$\psi(s) \leq (k-1) + \psi(t) \leq (|V|-1) + \psi(t) < |V|$$

Thus if we continuously keep track of a preflow and a distance label, and keep adjusting the preflow and distance label to obtain a flow with a distance label, we have obtained a maximal flow. In order to describe a method achieving this, we call a vertex $v \neq s, t$ **active** if $f(\delta_{\text{out}}(v)) - f(\delta_{\text{in}}(v)) > 0$, and an edge $vw \in G_f$ **admissible** if $\psi(v) = \psi(w) + 1$. The goal is to find a preflow with no active vertices, in which case we have a flow.

We now describe the push-relabel algorithm for finding the maximal flow in a graph. Begin by setting

$$f(e) = \begin{cases} \mu(e) & e = sv \\ 0 & \text{otherwise} \end{cases} \quad \psi(v) = \begin{cases} |V| & v = s \\ 0 & v \neq s \end{cases}$$

Then ψ is a distance label for the preflow f , since G_f has no edges leaving s .

Suppose there is an active vertex v relative to f . If there are no admissible edges leaving v , we may introduce an admissible edge by setting $\psi(v) = \min_{w \in G_f} \psi(w) + 1$. This leaves the f a preflow, and ψ a distance label. If there is an admissible edge vw leaving v , we set $f(vw)$ to be

$$\min[\mu_f(vw), f(\delta_{\text{out}}(v)) - f(\delta_{\text{in}}(v))]$$

Then $f \leq \mu$ still holds, and $f(\delta_{\text{out}}(v)) - f(\delta_{\text{in}}(v)) \geq 0$, since we can only have let at most $f(\delta_{\text{out}}(v)) - f(\delta_{\text{in}}(v))$ flow out. Also $f(\delta_{\text{out}}(w)) - f(\delta_{\text{in}}(w)) \geq 0$, since we have only let more flow in. The distance labelling is still a distance labelling, because we introduce an edge wv in G_f , which requires that $\psi(w) \leq \psi(v) + 1 = \psi(w) + 2$, and an edge from vw , so that $\psi(v) = \psi(w) + 1 \leq \psi(w) + 1$. Thus the fact that ψ is a distance labelling of the preflow f is an invariant of the algorithm.

Lemma 1.7. *If f is a preflow and ψ a distance label for f , then if $f(\delta_{\text{out}}(v)) - f(\delta_{\text{in}}(v)) > 0$, there is a (v, s) path in G_f .*

Proof. s □

1.6 Cuts in Undirected Graphs

Let G be an undirected graph with capacities $\mu \geq 0$. For a subset of vertices U let $\delta(U)$ denote all edges with a single endpoint in U . A **cut** in

such a graph is $\emptyset \subsetneq U \subsetneq V$, with value $\mu(\delta(U))$. The global minimum cut problem is to find the cut with $\mu(\delta(U))$ minimized, without regard to where any particular pair (s, t) resides. A naive way to calculate the global minimum cut is to calculate a minimum (s, t) cut, for all pairs (s, t) . This requires n^2 calls to the maximum flow problem.

And while we discuss this problem, we briefly sketch out how to compute a max flow in an undirected graph. Given such a graph G with capacity μ , we define H to be the directed graph with the same vertices as G , and for each edge $uv \in G$, two edges uv and vu in H , with a capacity function ψ such that $\psi(uv) = \psi(vu) = \mu(uv)$. Given any flow f in the directed graph, we obtain a flow f' in the undirected graph of the same value by subtracting $\min(f(uv), f(vu))$ from each edge. Thus we can compute a max flow in $O(V^2E)$ time in an undirected graph as well. The value of min cuts in the directed graph is the same as the value of min cuts in the undirected graph, so the max flow min cut theorem holds here as well.

More smartly, we notice that if s is fixed, s lies on either side of the cut, and since $\mu(\delta(U)) = \mu(\delta(U^c))$, we need only try each (s, t) cut as t ranges over all $t \neq s$, which only requires $n - 1$ calls to the max flow function. In a directed call, we need try all (s, t) cuts and (t, s) cuts, which requires $2(n - 1)$ calls to max flow.

We can do much better than just computing a min-cut in $O(n)$ times. It turns out that there is a way to compactly represent all minimum cuts via $O(n)$ calls to a min-cut algorithm. The first step to the representation is to find a useful relationship between the values of the numbers $\lambda_{u,v}$, the capacities of a min (u, v) cut.

Lemma 1.8. *Let v_1, \dots, v_k be vertices. Then $\lambda_{v_1, v_k} \geq \min(\lambda_{v_1, v_2}, \dots, \lambda_{v_{k-1}, v_k})$.*

Proof. Assume that $k \geq 3$, since the proof is trivial for $k = 2$. Let C be a (v_1, v_k) cut with $\mu(\delta(C)) < \min(\lambda_{v_1, v_2}, \dots, \lambda_{v_{k-1}, v_k})$. Inductively, we conclude that $v_2, \dots, v_k \in C$. But then by definition C cannot be a (v_1, v_k) cut. \square

Let $T = (V, E)$ be a tree, and $e \in E$ an edge in the tree. Then the set of nodes in some connected components of the graph $T' = (V, E - \{e\})$ is known as the **fundamental cut**. There are only two connected components, one for each vertex on the ends of e . A **Gomory-Hu Tree** for an undirected graph G is a tree T with the same vertices as G , such that the fundamental cut for every edge $uv \in T$ is a minimal uv cut. It turns out

that Gomory-Hu trees always exist, can be computed in $O(n)$ calls to a min (s, t) cut function, and compactly represent every min cut in a graph, for if (v_1, \dots, v_n) is the unique simple path in T , then

$$\lambda_{v_1, v_n} = \min(\lambda_{v_1, v_2}, \dots, \lambda_{v_{n-1}, v_n})$$

Certainly the left side is \geq to the right side. On the other hand, by construction of the tree, the fundamental cut of (v_i, v_{i+1}) is also a (v_1, v_n) , hence the left side is \leq to the right side.

Theorem 1.9 (Cut Submodularity). *For $A, B \subset V$, we have*

$$\mu(\delta(A)) + \mu(\delta(B)) \geq \mu(\delta(A \cap B)) + \mu(\delta(A \cup B))$$

Proof. Fix an edge uv . We shall show that every edge counted on the right side occurs at least as many times on the left side.

- Suppose $uv \in \delta(A \cap B)$, $uv \notin \delta(A \cup B)$. Then we may assume $u \in A \cap B$, $v \notin A \cap B$, $v \in A \cup B$. Then $\mu(uv)$ is counted once on the right hand side, and once on the left hand side.
- Suppose $uv \notin \delta(A \cap B)$, $uv \in \delta(A \cup B)$. Thus we may assume $u \in A \cup B$, $v \notin A \cup B$, $u \notin A \cap B$. Then $\mu(uv)$ is counted once on the right hand side, and once on the right hand side as well.
- Suppose $uv \in \delta(A \cap B)$, $uv \in \delta(A \cup B)$. Then we may assume $u \in A \cup B$, $v \notin A \cup B$, and so $u \in A \cap B$. Then $\mu(uv)$ is counted once on the right hand side, and twice on the left hand side.

Equality holds when $\delta(A \cap B)$ is disjoint from $\delta(A \cup B)$. □

Lemma 1.10. *Let $s \neq t$ be two vertices in a graph G , and let C be a min (s, t) cut. If $u \neq v$ are also distinct vertices, then there is a min (u, v) cut D with $C \subset D$ or $D \cap C = \emptyset$.*

Proof. Note that the statement is really redundant, because $C \subset D$ occurs exactly when $D^c \cap C = \emptyset$. Let B be any minimum (u, v) cut. We will modify B to another min cut with the desired properties. Note that

$$\mu(\delta(C)) + \mu(\delta(B)) \geq \mu(\delta(C \cap B)) + \mu(\delta(C \cup B))$$

We may assume $s, u \in B$. Then $t \notin B$, $C \cap B$ is an (s, t) cut, hence $\mu(\delta(C \cap B)) \geq \mu(\delta(C))$, and so

$$\mu(\delta(B)) \geq \mu(\delta(C \cup B))$$

Now $u \in C \cup B$, and $v \notin C \cup B$ (WHY?), so $D = C \cup B$ satisfies the properties of the lemma. □

1.7 Potentials

Fix a directed graph G with edge costs c . A **potential** for G is a mapping $\phi : V \rightarrow \mathbf{R}$ such that

$$c_\phi(uv) = c(uv) - (\phi(v) - \phi(u)) \geq 0$$

for all edges uv in G . The costs c_ϕ are known as the potential costs.

Lemma 1.11. *For any $s - t$ path P in G ,*

$$c_\phi(P) = c(P) - (\phi(t) - \phi(s))$$

Proof. If $P = (v_1, \dots, v_n)$, then we have an alternating sum

$$c_\phi(P) = \sum_{k=1}^{n-1} c(v_k v_{k+1}) + \phi(v_k) - \phi(v_{k+1}) = \sum_{k=1}^{n-1} c(v_k v_{k+1}) + \phi(v_1) - \phi(v_n)$$

and the right side is exactly the form considered in the statement of the theorem. \square

This is a discrete version of the fact that the energy change of a particle under a conservative force is invariant of the path of the particle, but only the change of the particles position in space. A simple corollary to the lemma is that the minimal cost paths relative to the metric c are the same as the minimal cost paths relative to c_ϕ .

If a graph has negative edge weights, but possesses some potential ϕ , then we can compute minimal paths under the cost function c_ϕ (which is now has positive edge weights everywhere) using Dijkstra's algorithm. There is a simple method to finding a potential, which we now detail.

If P is a cycle in a graph G with potential ϕ , we calculate

$$c_\phi(P) = c(P) \geq 0$$

so that G has no negative weight cycles. Conversely, if G has no negative weight cycles, then it turns out that G has a potential, and this potential can be computed in polyomial time.

Theorem 1.12. *A potential can be computed for a graph G can be computed in $O(nm)$ time.*

Proof. Given G , compute a graph H by adding a new vertex r to G , and adding edges from r to every vertex in G . Set $c(rv) = 0$ for all vertices $v \in V$, and let $\phi(v)$ be the cost of the cheapest walk from r to v . This is well defined since G has no negative weight cycles (there is a cheapest walk with no repeated vertices, for if a path has a repeated vertex, it contains a cycle, which must have non-negative weight, and this cycle can be removed with no penalty to the length of the walk – we then need only consider the finite set of paths with no repeated vertices to obtain an answer). Note that $c(uv) + \phi(u) \geq \phi(v)$, so ϕ is a potential for G . The Bellman-Ford can compute ϕ in $O(nm)$ time. \square

Let ω be a differential 1-form on a manifold M such that

$$\int_{\gamma} \omega \geq 0$$

for any cycle γ . Then if γ and δ are two paths from a point x to a point y , we find

$$\int_{\gamma} \omega - \int_{\delta} \omega, \int_{\delta} \omega - \int_{\gamma} \omega \geq 0$$

so $\int_{\gamma} \omega = \int_{\delta} \omega$, and given a fixed x , we can define

$$F(y) = \int_{\gamma} \omega$$

where γ is a path from x to y , and one verifies that $dF = \omega$, so that

$$\int_{\gamma} \omega - (F(y) - F(x)) = 0$$

a stronger fact which is analogous to the result above, although our result is weaker because in directed graphs we cannot perform the swapping trick to obtain equality of all paths.

1.8 Minimum Cost Flows

Throughout this section, we fix a directed graph with edge costs c , edge capacities μ , fixed vertices s, t , and a target flow value γ . The minimum cost

flow problem tries to find a flow f with value $\text{val}(f) \geq \gamma$, which minimizes the cost $\text{cost}(f) = \sum c(e)f(e)$, or to determine if no such flow exists.

We can certainly compute in polynomial time if a flow exists satisfying the constraints, because we can find a maximal flow in the graph. It is not so clear that we can find such a flow with minimal cost. It shall always be the case that the minimal flow satisfies $\text{val}(f) = \gamma$, because given any flow f , the value of the flow $g = \text{val}(f)/\gamma f$ satisfies $\text{val}(g) = \gamma$, and $\text{cost}(g) = \text{val}(f)/\gamma \sum c(e)f(e)$, which will be smaller than $\text{cost}(f)$ if $\text{val}(f) \neq \gamma$.

1.9 Minimal Cost Bipartite Matching

We now take a brief aside to discuss how minimal cost flows can be used to solve the analogous problem in bipartite graphs. Given the name, the minimal cost bipartite matching problem should be fairly self explanatory. We take a particular Bipartite graph $G = (L \cup R, E)$ with edge costs c , and we must find a minimum cost perfect matching on G in polynomial time. That is, our goal is a perfect matching M which minimizes

$$\text{cost}(M) = \sum_{e \in M} c(e)$$

The similarity to the minimal cost flow problem immediately suggests a connection between the two problems.

We can formulate the perfect matching problem as a minimal cost max flow problem, by modifying slightly the standard reformulation of the maximal matching problem as a max flow problem. The original process is to construct a directed bipartite graph H , whose vertices are obtained from G by adding two new vertices s and t , orienting all edges in E so that they point from L to R , and adding new edges sv , for $v \in L$, and wt , for $w \in R$. If we define a capacity function $\mu = 1$, then there is a one-to-one correspondence between integral valued $s - t$ flows and matchings in the graph. To see this, note that an integral valued flow f gives rise to a characteristic function χ_{A_f} on the edges of G , and the flow conservation law shows that if e and e' are adjoint, then $\chi_{A_f}(e) + \chi_{A_f}(e') \leq 1$, so that A_f forms a matching on G . Correspondingly, the reverse process takes a matching to a characteristic function on the edges of G , which can easily be completed into a flow on H .

Now we define a cost function c' on H in a way which mirrors the cost function c on G . We let

$$c'(sv) = c'(wt) = 0 \quad c'(vw) = c(vw)$$

Then we see that the correspondence described above preserves the cost function. That is, if a matching M corresponds to a flow f , then $\text{cost}(M) = \text{cost}(f)$, because f is obtained from M by adding edges of the form sv and wt , which add no value to the cost of the flow. Thus if we find a minimal cost maximal flow in H , we can find a minimal cost perfect matching in G , and we know we can find such a maximal flow using the successive shortest paths algorithm, and we need only compute $|L| = |R|$ successive paths, since each iteration of the algorithm improves the length of the candidate path, hence now have an algorithm for the minimal cost perfect matching problem which runs in $O(n(m + n \log n))$ time.

1.10 Minimum Mean Cycle Cancellation

We already know that a flow f has minimal cost if G_f contains no negative-cost cycles. This suggests that a strategy for finding minimal cost flows is obtained from successively pruning negative-cost cycles from G_f until there are none left. As with Ford Fulkerson, we must be careful which cycles we prune in order to guarantee a good asymptotic speed to our algorithm. It turns out that fast asymptotic results are available if we prune negative cycles with ‘minimum mean’.

Algorithm 1 The Minimum Mean Cycle Canceling Algorithm

- 1: $f \leftarrow$ any (s, t) flow of value γ (found in $O(n^3 \sqrt{m})$ time)
 - 2: **while** G_f contains negative weight cycles **do**
 - 3: $C \leftarrow$ a cycle in G which minimizes its mean cost $c(C)/|C|$.
 - 4: Augment f along C .
 - 5: **return** f
-

The algorithm certainly terminates, since the cost of f decreases for every iteration of the algorithm, at a rate bounded by the rational values of the edge costs. It remains to be seen that the algorithm terminates in a

polynomial amount of time (and this is not immediately obvious, and is not true if we do not take cycles which minimize the mean cost).

First, we argue why we can find a minimum ratio cycle C in polynomial time.

Lemma 1.13. *Any Eulerian multigraph $H = (W, F)$ with edge costs c contains a cycle C such that*

$$\frac{c(C)}{|C|} \leq \frac{c(F)}{|F|}$$

Proof. Perform a cycle decomposition on H , writing $F = C_1 \amalg \cdots \amalg C_n$, we find

$$\frac{c(F)}{|F|} = \frac{c(C_1) + \cdots + c(C_n)}{|C_1| + \cdots + |C_n|}$$

If $c(C_i)/|C_i| > c(F)/|F|$ for each i , then $c(C_i)|F| > c(F)|C_i|$ also holds for each i , and so, summing up we find

$$c(F)|F| = \sum_i c(C_i)|F| > \sum_i c(F)|C_i| = |F|c(F)$$

a clear contradiction which proves the existence of a C_i satisfying the constraints of the lemma. \square

Given any cycle C , the corresponding multigraph (which contains duplicates of repeated edges in the cycle) is Eulerian, and can thus be broken up into disjoint cycles. The lemma above shows that if C is not a simple cycle, we can always find a strictly smaller cycle C' without increasing the mean value of the cycle, so that we need only search simple cycles to find a minimum mean.

For $k \leq n$, define $\delta_k(v)$ to be the cheapest closed walk from v to itself using exactly k (not necessarily distinct) edges. If we let $\delta_k(v, w)$ be the cheapest path from v to w using k edges, then we find

$$\delta_k(v, w) = \min_{u \in V} c(v, u) + \delta_{k-1}(u, w)$$

Each value of $\delta_k(v, w)$ can be computed in $O(m)$ time, and so the set of all $\delta_k(v, w)$ can be computed in $O(mn^3)$ time. Let v^* and k^* minimize $\delta_{k^*}(v^*)/k^*$, and pick k^* to be smallest to break ties. Then the walk from v^* in k^* stops is verified to be a cycle, from the above remark, and thus is a minimum ratio cycle, since it is a minimum ratio walk in general.

Theorem 1.14. *The number of iterations of the minimum mean cycle cancelling algorithm is $O(m^2 n \log n)$.*

Proof. Consider the sequence of flows

$$f_1, f_2, \dots, f_N$$

which are found in the MMCC algorithm, together with the cycles

$$C_1, \dots, C_{N-1}$$

which are augmented to produce the sequence of flows. For $i < j$, define

$$\alpha(i, j) = \{e, \overleftarrow{e} : e, \overleftarrow{e} \in C_i \cup C_j\}$$

Lemma 1.15. *If $\alpha(i, k) = \emptyset$ for $i < k < j$, then we find*

$$\frac{c(C_i)}{|C_i|} \leq \frac{n}{n - |\alpha(i, j)|} \frac{c(C_j)}{|C_j|}$$

Proof. Let $H = C_i \sqcup C_j - \alpha(i, j)$ (essentially, H is obtained from the cycles C_i and C_j by keeping duplicate edges, and ‘cancelling out’ edges with their reverse). Then H is Eulerian, for $C_i \sqcup C_j$ is Eulerian, and removing edges along with their reverse changes both the in degree and out degree of a vertex by the same amount. Furthermore, each edge in H is in E_{f_i} because $\alpha(i, k) = \emptyset$, so

$$\begin{aligned} \frac{c(C_i)}{|C_i|} (|C_i| + |C_j|) &= \frac{c(C_i)}{|C_i|} (|H| + |\alpha(i, j)|) \\ &\leq c(H) + |\alpha(i, j)| \frac{c(C_i)}{|C_i|} \\ &= c(C_i) + c(C_j) + |\alpha(i, j)| \frac{c(C_i)}{|C_i|} \end{aligned}$$

where we have used the fact that $c(H) = c(C_i) + c(C_j)$ since the cost of reverse edges cancels out. This can be rephrased as

$$\frac{c(C_i)}{|C_i|} (|C_j| - |\alpha(i, j)|) \leq \frac{c(C_j)}{|C_j|} |C_j|$$

And dividing out, we find

$$\frac{c(C_i)}{|C_i|} \leq \frac{|C_j|}{|C_j| - |\alpha(i, j)|} \frac{c(C_j)}{|C_j|}$$

and this inequality can be weakened to the form above, because $x/(x - y)$ is a decreasing function of x for $y > 0$, and $n/(n - |\alpha(i, j)|) \geq 0$. \square

Returning to our original proof, we see that $j = i + 1$ satisfies the theorem, and proves that the minimum mean ratios are non-decreasing. If $i < j$ is the smallest value such that $\alpha(k, i) \neq \emptyset$, for some $i \leq k < j$, then we find

$$\frac{c(C_i)}{|C_i|} \leq \frac{n}{n-2} \frac{c(C_j)}{|C_j|} = \left(1 - \frac{n}{2}\right) \frac{c(C_j)}{|C_j|}$$

and we have a bound $j \leq i + m$. To see this, note that at least one edge is removed in the residual graph from each of the cycles $C_i, C_{i+1}, \dots, C_{i+m}$ during the algorithm. Thus we conclude

$$\frac{c(C_i)}{|C_i|} \leq \left(1 - \frac{n}{2}\right)^n \frac{c(C_{i+nm})}{|C_{i+nm}|} \leq e^{-1/2} \frac{c(C_{i+nm})}{|C_{i+nm}|} \leq 2 \frac{c(C_{i+nm})}{|C_{i+nm}|}$$

If the edge weights are integral (and all rational cost problems can be reduced to this form), this already gives us enough information to prove a polynomial time bound. Note that since the minimum mean cycles are simple, we must have

$$\frac{|c(C_i)|}{|C_i|} \geq \frac{1}{n}$$

and

$$\frac{|c(C_1)|}{|C_1|} \leq \min_{e \in E} |c(e)|$$

so if N is large enough that $2^{-N} \min_{e \in E} |c(e)| \leq n^{-1}$, then

$$\frac{c(C_{i+Nnm})}{|C_{i+Nnm}|} \geq \frac{1}{2^N} \frac{c(C_1)}{|C_1|} \geq -\frac{\min_{e \in E} |c(e)|}{2^N} \geq -\frac{1}{n}$$

which forces the mean cycle ratio to be positive, and hence the algorithm will terminate. We can choose any $\lg(n \min_{e \in E} |c(e)|) < N$, so the algorithm terminates in at most

$$nm \lceil \lg n \min_{e \in E} |c(e)| \rceil$$

iterations, which is polynomial in the bit complexity of the algorithm, since if the costs are given with l bits, then the total bit complexity is about $\Theta(n + ml)$, and $|c(E)| \leq 2^l$, so

$$nm \lceil \lg n \min_{e \in E} |c(E)| \rceil \leq nm \lg(n) + lnm$$

This finishes our initial analysis of the mean cycle canceling algorithm. \square

The polynomial bound for this algorithm is absolutely awful, and we can do much much better. In the next lecture we will begin the arguments that give us a much better bound on the runtime on this algorithm.

1.11 Minimum Spanning Trees and Arborescences

Let G be an undirected graph, together with a (possibly-negative) edge cost function c . A **spanning tree** T on G is a connected subset of edges containing no cycles (a tree) such that every vertex is covered by an edge in the tree. Note that this definition is equivalent to T having no cycles, and $n - 1$ edges (a fact proven most easily by induction, by removing vertexes and edges from T which form the leaves of T). There is another equivalent definition which is more important to the spanning tree problems we'll discuss. T is a spanning tree if it has $n - 1$ edges, and $|E(S)| \leq |S| - 1$ for all vertex sets S (where $E(S) = \{uv : u, v \in S, uv \in T\}$).

TODO: Draw a picture of S bounding the number of edges in the set S

The minimum spanning tree problem is self explanatory – given G , find a spanning tree containing the fewest edges. We can formulate the problem as an *LP* in exponentially many constraints in functions $x : E \rightarrow \mathbf{R}$,

$$\begin{aligned} \min \quad & \sum_e c(e)x(e) \\ \text{s.t.} \quad & x(E(S)) \geq |S| - 1 \quad \forall S \subsetneq V \\ & x(E(V)) = n - 1 \\ & 0 \leq x \end{aligned}$$

Note that the constraint $x \leq 1$ is already encoded in the problem by taking S to be a pair of vertices corresponding to an edge. Our discussion above entails that any integral solution to this algorithm corresponds to a spanning tree in G . It can be shown that all extreme points are integral, but we shall only show that we can choose an optimal extreme point which is integral.

In its current form, the minimal spanning tree linear program is unfeasible, having far too many constraints. We obtain a much better optimization problem if we switch to the dual of this program. This can be motivated by the fact that x behaves more like a linear functional than a vector in this linear program (the constraints of the program are expressed as x 's operations on the vector space $\mathbf{R} \cdot E$ generated by the edges E), so that the program is the dual of some other, more easily understood linear program. The dual program finds functions $y : 2^V - \{\emptyset\} \rightarrow \mathbf{R}$ which satisfy

$$\begin{aligned} \max \quad & -\sum (1 - |S|)y(S) \\ \text{s.t.} \quad & \sum_{e \in E(S)} y(S) \geq -c(e) \quad \forall e \in E(S) \\ & y(S) \geq 0 \quad \forall S \neq V \end{aligned}$$

Kruskal's algorithm, a method for finding the minimal spanning tree, can be viewed as a combinatorial method to solving the dual LP of the spanning tree problem. We first recall the simple, greedy method to form a spanning tree. We can verify this algorithm's correctness using the dual-

Algorithm 2 Kruskal's Algorithm

- 1: Set $T = \emptyset, K = E$.
 - 2: **while** T is not a spanning tree **do**
 - 3: Remove $e \in K$ with minimal weight.
 - 4: Append e to T if it connects two connected components of T .
 - 5: **return** T
-

ity of linear programming. Suppose that T contains the edges e_1, \dots, e_n , which are placed in the order they were added to T . Let S_k be the component containing e_k in the graph consisting only of the edges e_1, \dots, e_k (it is the component that was freshly merged together in the k 'th iteration of the algorithm). Let $x : E \rightarrow \{0, 1\}$ be the indicator function of T , and

$y : 2^V - \{\emptyset\} \rightarrow \mathbf{R}$ be defined by letting $y(S_i) = c(e_j) - c(e_i)$, where j is the smallest index greater than i such that one of the endpoints of e_j contains points in X_i , and define $y(S_n) = y(V) = -c(e_n)$. Define $y(S) = 0$ otherwise. For any edge e , by the telescoping sum property of our definition we have

$$\sum_{e \subset S} y(S) = -c(e_i)$$

where i is the smallest index such that X_i contains both endpoints of e . The way we selected edges guarantees that $c(e_i) \leq c(e)$, so that our constraints are satisfied, and our tight for e_1, \dots, e_n .

We now verify complementary slackness, so that x is verified optimal. If $x(e) > 0$, then our calculation above shows

$$-\sum_{e \subset S} y(S) = c(e)$$

If $y(S) > 0$, then $S = S_i$ for some i , hence S_i is a connected tree, and contains $n-1$ vertices, so $x(E(S)) = |S| - 1$. Complementary slackness guarantees that x and y are optimal solutions to their corresponding algorithms, verifying correctness of Kruskal's algorithm.

1.12 Arborescences

An (out) **arborescence** in a directed graph with root s is a subset T of $n-1$ edges, such that there is a unique directed path from s to any other vertex. Given a cost function $c \geq 0$ and vertex s , we want to try and find the min-cost arborescence at s . Like the minimal cost spanning tree problem, we can express this problem as an LP,

$$\begin{aligned} \min \quad & \sum_e c(e)x(e) \\ \text{s.t.} \quad & x(\delta^{\text{in}}(U)) \geq 1 \quad \forall U \subset V - \{s\} \\ & x \geq 0 \end{aligned}$$

If T is an arbitrary arborescence at v , then the corresponding characteristic function certainly is a feasible solution to the problem. The problem with this LP is that there are solutions (possibly even optimal ones) which don't look like arborescences. However, we shall not use the LP to solve

linear programs, but instead use properties of linear programs to guarantee optimality. We can form a simple argument to show the existence of optimal solutions which are arborescence. First, note that we only ever have upper bounds to the algorithm, and there is no need to have $x(e) > 1$ to satisfy any of the constraints, due to the positivity of x , so we may always assume that an optimal integral solution takes value in $\{0,1\}$, and thus correspond to characteristic functions of edges. Second, we see that the subgraph formed contains paths from v to any other vertex. Since $x(\delta^{\text{in}}(V - s)) \geq 1$, there is an edge connecting s to some vertex v_1 . Then $x(\delta^{\text{in}}(V - \{s, v_1\})) \geq 1$, so there is some edge from s to v_1 , or v_1 to v_2 . In either case, there is an edge to v_2 . Continuing this process gives you a path to any vertex in the graph. We may now assume our subgraph is a tree, because we can always remove edges to obtain an algorithm that is at least as optimal. Thus we have argued that there is an optimal solution which is an arborescence, like we are looking for.

Let us form the dual LP, so we can use complementary slackness to verify optimality.

$$\begin{aligned} \max \quad & \sum_S y(S) \\ \text{s.t.} \quad & \sum_{e \in \delta^{\text{in}}(S)} y(S) \leq c(e) \\ & y \geq 0 \end{aligned}$$

We now describe the algorithm which gives us minimal cost arborescences. Note that y is only included here for purposes of algorithmic correctness, and needn't be coded up in real algorithms.

* DRAW A PICTURE OF CONTRACTING CYCLES

Algorithm 3 Edmond's Algorithm

- 1: $y_{\{v\}} = \min c(uv)$, for $v \neq r$.
 - 2: Set $F = \{e_w : w \in V - \{r\}\}$, where e_w is the cheapest edge entering $w \neq r$.
 - 3: **while** F has a cycle **do**
 - 4: Let G' be the graph obtained by contracting each cycle in (G, F) .
 - 5: Define $c'(uv) = c(uv) - y(\{v\})$ if v is on a cycle, else $c'(uv) = c(uv)$.
 - 6: Let (T', y') be a min-cost arborescence in G' .
 - 7: Define $T = (F \cup T') - \{uv : u \text{ on a cycle}, uv \in T'\}$.
 - 8: Set $y(S) = 0$ if S cuts a cycle, and $y'(S)$ if $S = [S]$ after contraction.
 - 9: **return** F
-

Chapter 2

Matroids

If you have a problem, and a matroid is involved, you can probably use a greedy algorithm to solve the problem. This is very useful, because greedy solutions are normally far faster than conservative methods. A matroid is a set of sets which can be ‘greedily expanded’ without losing the properties defining a matroid. Formally, a matroid is a collection of subsets of some set X containing the emptyset, closed under the subset relation, and such that if two sets A and B are in the matroid with $|A| < |B|$, there is $z \in B - A$ such that $A \cup \{z\}$ is in the matroid.

Example. *If G is a graph, the set of edge subsets which contain no cycle forms a matroid. Certainly it contains the emptyset, and certainly a subset of edges with no cycle contains no cycle.*

Example. *If V is a vector space, the collection of linearly independent vector sets forms a matroid. Essentially, you can greedily expand vectors to form a basis.*

Example. *If X is a finite set, and L is a laminar family over X , together with a function $f \geq 0$ defined on the Laminar family, then the same of subsets A such that $|A \cap B| \leq f(B)$ for each $B \in L$, forms a matroid known as a laminar family.*

Example. *If G is a directed graph, and $A, B \subset V$ are given, the set of subsets of A which contains vertex disjoint paths into B is a matroid, known as a Gammoid.*

Example. *Given a connected graph G , the set of bond matroid.*

Example. Given a field extension $F \subset K$, the set of finite subsets Y of F with transcendence degree $|Y|$ forms a matroid, known as the algebraic matroid.

An independence oracle is an algorithm to decide whether an arbitrary subset is an element of a given matroid in polynomial time. This is key to having efficient algorithms for solving matroids.

A base of a matroid is a maximal set – an element B such that if $B \subset B'$, and B' is in the matroid, then $B = B'$. If our matroid consists of finite sets, then two bases have the same cardinality. Indeed, if $|B| < |B'|$, there is $z \in B' - B$ such that $B + z$ is in the matroid, which cannot occur since this is an expansion of B .

The first algorithm we consider is to efficiently find a min weight base for a matroid over a set X , for some function $c : X \rightarrow \mathbf{R}$. This is a generalization of the minimum spanning tree problem. The algorithm to do this is to just greedily add the minimum element of X to a set until we cannot anymore. The set B we end up with is a base by construction, because the order we add elements is immaterial. To prove that B is minimal, consider a minimal base B^* , and order the elements of both sets

$$B = \{x_1, \dots, x_n\} \quad B^* = \{x_1^*, \dots, x_n^*\}$$

We prove by induction that $w(x_i) \leq w(x_i^*)$. Indeed, consider the partial set $B_{i-1} \setminus \{x_1, \dots, x_{i-1}\}$ and $B_i^* \setminus \{x_1^*, \dots, x_i^*\}$. There is a minimal weight element x_k^* of B_i^* which is not in B_{i-1} , and can be safely added to B , and $w(x_i) \leq w(x_k^*)$. If $k \leq i$, this shows that $w(x_i) \leq w(x_i^*)$. (FIGURE THIS PROOF OUT LATER)

The **rank** $r(A)$ of a subset A of X , with respect to a matroid M is the cardinality of the largest subset of A in M . Then r is order preserving, and $r(A) \leq |A|$. More interestingly,

$$r(A) + r(B) \geq r(A \cup B) + r(A \cap B)$$

To prove this, let J be a maximal subset of $A \cap B$ in the matroid, $K \subset A - B$ a complementary set for J , so that $|K| + |J| = r(A)$. This is always possible because we can extend J to a base in A . Do the same for J in B , forming L such that $|K| + |J| + |L| = r(A \cup B)$. By construction, $J \cup K$ is independent in A , and L is independent in B . But then by maximality we obtain the bound above. It turns out that the rank function of a matroid uniquely determines the matroid, but we won't discuss that.

A subset C of X is a **circuit** if it is not independent, but removing any particular element of C makes it independent. If Y is any independent set, and $Y \cup \{x\}$ is independent, then there is a unique circuit contained in $Y \cup \{x\}$. We call this circuit $C(Y, x)$. To find this circuit, consider a minimal element of $Y \cup \{x\}$ which is not independent. The hard part is showing this is the only circuit. Let C_1 and C_2 be two circuits. If $C_1 \neq C_2$, let $x \in C_1 \cap C_2$, and $y \in C_1 - C_2$ (without loss of generality). We will show there is a circuit contained in $(C_1 \cup C_2) - \{x\}$ contains a circuit. Indeed, using submodularity we find

$$\begin{aligned}
|C_1| - 1 + r((C_1 \cup C_2) - \{x, y\}) + |C_2| - 1 &= r(C_1) + r((C_1 \cup C_2) - \{x, y\}) + r(C_2) \\
&\geq r(C_1) + r((C_1 \cup C_2) - \{y\}) + r(C_2 - \{x\}) \\
&\geq r(C_1 - \{y\}) + r(C_1 \cup C_2) + r(C_2 - \{y\}) \\
&= |C_1| - 1 + r(C_1 \cup C_2) + |C_2| - 1
\end{aligned}$$

The point of this calculation is that now we know $r(C_1 \cup C_2) = r((C_1 \cup C_2) - \{x, y\})$, so that we cannot add x or y to any maximal independent set in $(C_1 \cup C_2) - \{x, y\}$ while keeping it independent. In particular, $(C_1 \cup C_2) - \{x\}$ is not independent, so there must be a circuit contained in the set.

That there is always a unique circuit ‘screwing us over’ is very useful in analyzing the matroid intersection problem. Given two matroids M and N over a set X , the intersection problem asks us to find the maximal set which is contained in both matroids. Similarly, we may have a subset Y independent in both M and N , and we wish to know whether we can extend Y while preserving independence. Consider the following algorithm: given Y , we construct a directed graph with nodes in Y and $X - Y$, and directed edges ab (for $a \in Y, b \in X - Y$) if $Y - a + b \in M$, and ba if $Y - a + b \in N$. We find a shortest path from Y_1 to Y_2 where

$$Y_1 = \{x \in X - Y : Y + x \in M\} \quad Y_2 = \{x \in X - Y : Y + x \in N\}$$

and we then augment along that path to find a better matroid. If there is no path from Y_1 to Y_2 , we have a maximal matroid.

2.1 Linear Programs for Matroids

Let M be a matroid over a set X , together with a weight function $w : M \rightarrow \mathbf{R}$. We can formulate the maximum weight independent set problem over

X as a linear program. Specifically, we consider the space of functions in \mathbf{R}^X , and take the linear program

$$\begin{aligned} \max \quad & \sum_{e \in X} w(e)x(e) \\ \text{s.t.} \quad & x(A) \leq r(A) \quad \forall A \subset X \\ & x \geq 0 \end{aligned}$$

in particular, the constraints guarantee that $0 \leq x(e) \leq 1$ for all $e \in X$, so integral points correspond to indicator functions over X . What's more, integral points also correspond to independent sets in the matroid. If $x \in \mathcal{X}_B$, then $B = x(B) \leq r(B) \leq B$, so $r(B) = B$, and so B is independent.

It turns out that (surprise surprise!) all extreme points are integral, so we can solve the problem in polynomial time. Our proof will use an interesting technique known as the uncrossing method, which 'thins' the constraints of the LP so they correspond to a Laminar family (If we say A and B are crossed if $A \cap B \neq \emptyset$, then by breaking them into disjoint chains we have 'uncrossed them'). First, we embed 2^X in \mathbf{R} by mapping $A \subset X$ to the 0/1 indicator function χ_A . Then we let \mathcal{C} to be the largest linearly ordered subset of 2^X (a *chain* in 2^X) only containing subsets A for which we have a tight constraint, $x(A) = r(A)$, and such that the χ_A are independent. We claim that the χ_A span \mathbf{R}^X .

Next, note that the functionals $x \mapsto x(A)$ are essentially the duals of the characteristic function χ_A , and since 2^X spans \mathbf{R}^X , the rows of the constraint matrix which x satisfies span \mathbf{R}^X . In particular, since x is extreme, the set of tight constraints spans \mathbf{R}^X , and it suffices for us to prove that $\text{span}(\mathcal{C})$ contains all χ_A such that $x(A) = r(A)$. Suppose that χ_B was not in $\text{span}(\mathcal{C})$, where $r(B) = x(B)$. If such a B exists, we may choose B such that the number of elements is minimized

$$\tau(B) = \{A \in \mathcal{C} : A \not\subset B \text{ and } B \not\subset A\}$$

is minimized. Note that we must have $\tau(B)$ must contain at least one element, or else $\mathcal{C} \cup \{B\}$ would be a chain of linearly independent vectors which is larger than \mathcal{C} . So let $T \in \mathcal{C}$ be such that $B \not\subset T$ and $T \not\subset B$. Note that by using the inclusion/exclusion principle, optimality, and the submodu-

larity property,

$$\begin{aligned}
r(B) + r(T) &= x(B) + x(T) \\
&= x(B \cup T) + x(B \cap T) \\
&\leq r(B \cup T) + r(B \cap T) \\
&\leq r(B) + r(T)
\end{aligned}$$

So that $x(B \cup T) + x(B \cap T) = r(B \cup T) + r(B \cap T)$, and since $x(B \cup T) \leq r(B \cup T)$, $x(B \cap T) \leq r(B \cap T)$, we can actually split this inequality into the equalities $x(B \cup T) = r(B \cup T)$, $x(B \cap T) = r(B \cap T)$.

It turns out that $\tau(B \cup T), \tau(B \cap T) \subsetneq \tau(B)$, which by optimality implies $\chi_{B \cup T}, \chi_{B \cap T} \in \text{span}(\mathcal{C})$. This is clearly in contradiction to our assumption, since then

$$\chi_B = \chi_{B \cup T} + \chi_{B \cap T} - \chi_T \in \text{span}(\mathcal{C})$$

so all that remains is to prove properties of the τ function. This is most easily shown by drawing out Venn diagrams, which is often useful in matroid theory, but we'll also provide a textual description (and drawing your own diagram will probably be much more revealing than looking at someone else's).

Consider S in \mathcal{C} . Then either $S \subset T$, or $T \subset S$. If $S \subset T$, then $S \subset B \cup T$ and so $S \notin \tau(B \cup T)$. If $T \subset S$, then $B \cap T \subset S$, so $S \notin \tau(B \cap T)$. If $S \in \tau(B \cup T)$ but $S \notin \tau(B)$, then we would have to have $T \subset S$. This also shows $S \subset B$, because if $B \subset S$, then $B \cup T \subset S$. But then $T \subset B$, which is a contradiction. Similarly, if $S \in \tau(B \cap T)$ but $S \notin \tau(B)$, then $S \subset T$ and $B \subset S$, so $B \subset T$, another contradiction. We finish this conclusion by noticing that $T \notin \tau(B \cap T), \tau(B \cup T)$ by obvious relations.

Thus we conclude that \mathcal{C} forms a basis of \mathbf{R}^X (in fact, \mathcal{C} forms a basis for \mathbf{Z}^X – the same proof follows through since we've never used real numbers, but we won't need this in our proof. Now our homework comes in handy. Consider the laminar constrained matching problem, over the Bipartite graph whose left vertices are subsets of X , and with only one right vertex, with an edge between every left vertex. We consider $L_L = \mathcal{C}$ as a laminar family over the left vertices, and $L_R = \emptyset$, with $b_A = r(A)$ for $A \in \mathcal{C}$. Every solution to this problem can be identified with a characteristic function over the vertices, and by the identification x can be seen as a feasible solution to the Laminar constraints. Since x is tight for *all* constraints, it must be a feasible solution to this problem, and in our homework we proved that therefore x has integral coordinates.

The matroid intersection problem can be formulated as an LP with integral extreme points in a very similar manner. We consider two matroids M and N , take the same solution space as the maximal independent set LP, and find solutions such that

$$\begin{aligned} \max \quad & \sum_{e \in X} w(e)x(e) \\ \text{s.t.} \quad & x(A) \leq \min(r_M(A), r_N(A)) \quad \forall A \subset X \\ & x \geq 0 \end{aligned}$$

The extreme points of this set are integral, which can be verified from the uncrossing technique of the last problem. Here we will end up with a laminar family of tight constraints \mathcal{C}_M and \mathcal{C}_N over each matroid, rather than a chain, prove that $\text{span}(\mathcal{C}_M \cup \mathcal{C}_N)$ is full, and then use the laminar family $L_L = \mathcal{C}_M$, $L_R = \mathcal{C}_N$ to prove integrality. Our homework problem does not generalize to intersections of three or more matroids, and for good reason! The matroid intersection problem for more than 2 matroids is NP complete.

Chapter 3

Submodular Functions

A **submodular function** is a function $f : 2^\Omega \rightarrow \mathbf{R}$, such that

$$f(S) + f(T) \geq f(S \cup T) + f(S \cap T)$$

One algorithm that occurs across computing science is to find a maximum set for a submodular function. Consider the greedy approximation which constructs a set by taking the point which increases the current set by the largest amount. If Z is the set constructed, and Z^* is an optimum set, we claim that $f(Z) \geq (1 - 1/e)f(Z^*)$ – this is the best approximation one can do in polynomial time, unless $P = NP$.

Lemma 3.1.

$$f(Z^*) \leq k(f(z) - f(z))$$

Proof.

$$f(Z^*) \leq f(Z^* \cup Z_{i-1})$$

□

Chapter 4

Perfect Graphs

There are many numbers of interest associated to graphs. For instance, we have $\alpha(G)$, the size of the maximum independent set, $\omega(G)$, the maximum clique size, $\chi(G)$, the chromatic number, and $k(G)$, the clique cover number (the minimum number of cliques required to cover the vertices of G). A graph is **perfect** if $\chi(H) = \omega(H)$ for every subgraph of G . For each graph, we associate the polytope

$$P_{\text{clique}}(G) = \{x \in \mathbf{R}\langle V \rangle : x \geq 0, x(C) \leq 1, \forall \text{ cliques } C : x(C) \leq 1\}$$

We claim this polytope is integral if and only if G is perfect.

Perfect graphs include the families of bipartite graphs and chordal graphs. Given a graph G , define the complement of G to be the graph \bar{G} obtained by ‘flipping edges’. Then

$$\alpha(G) = \omega(\bar{G}) \quad \chi(G) = k(\bar{G})$$

It is an important theorem of Lovatz, proved in 1972, that a graph is perfect if and only if its complement is perfect. Beyond the scope of this course is the proof that G is perfect if and only if G and \bar{G} contain no cyclic graphs of size ≥ 5 (the strong perfect graph theorem).

To prove Lovatz’s theorem, consider the graph G^{+v} , for $v \in V$, obtained by adding a new vertex connected to every vertex but v .

Lemma 4.1 (Replication Lemma). *G is perfect if and only if G^{+v} is perfect.*

Proof. The lemma is trivial if G has one vertex. For $|V| \geq 2$, it suffices to check $\chi(G^{+v}) = \omega(G^{+v})$. It is always true that $\alpha(G^{+v}) = \alpha(G) + 0/1$. If $\alpha(G^{+v}) = \alpha(G)$ \square

So suppose G is perfect. Let $x \in P_{\text{clique}}(G)$ be an extreme point. From what we know in class, we know $x \in \mathbf{Q}$. Thus there is an integer $\eta \neq 0$ such that $\eta x \in \mathbf{Z}$. So consider H , obtained by copying G η times, and then connecting all edges (u_i, v_i) , if $u = v$ or (u, v) is an edge in the original graph. The lemma above implies H is perfect. Let C_H be a max clique in H , with $|C_H| = \omega(H)$.