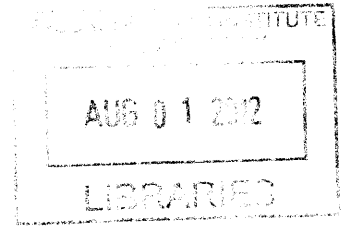


**A Functional Flow Framework for Cloud  
Computing**

ARCHIVES



by

Amy Zhang

B.S., EECS, Massachusetts Institute of Technology (2011)

B.S., Mathematics, Massachusetts Institute of Technology (2011)

Submitted to the Department of Electrical Engineering and Computer  
Science

in partial fulfillment of the requirements for the degree of

Master of Engineering in Electrical Engineering

at the

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## Abstract

This thesis covers a basic framework to calculate the maximum computation rate of a set of functions over a network. These functions are broken down into a series of computations, which are distributed among nodes of the network, with the output sent to the terminal node. We analyze two models with different types of computation costs, a linear computation cost model and a maximum computation cost model. We show how computation distribution through the given network changes with different types of computation and communication limitations. This framework can also be used in cloud design, where a network of given complexity is designed to maximize computation rate for a given set of functions. We provide a greedy algorithm that provides one solution to this problem, and create simulations for each framework, and analyze the results.

Thesis Supervisor: Muriel Médard

Title: Professor of EECS



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# Chapter 1

## Introduction

With the advent of cloud computing, many problems need to be addressed to keep the cloud secure, reliable, and low-cost. Much research has been conducted on possible protocols at a system level for cloud computing to minimize problems in privacy, reliability, and accountability, but research on a theoretical level has been more scarce. The goals of my project are to research and implement algorithms to both calculate and minimize access rates to the cloud under certain models, and to investigate the trade offs between communication and computation costs.

Our goal is to be able to design and analyze a network flow framework by modeling the amount of computation in each node of the network and using convex optimization methods to calculate the maximum computation rate and route flows through the network to achieve that rate. We consider an approach with given computation and communication limitations that affect the maximum computation rate. The two computation models we examine are a maximum computation cost (MCC) model and a linear computation cost (LCC) model. In the MCC model, the computation cost over the network is a function of the maximum computation, over all nodes in the network. In the LCC model, the computation cost in each node is a linear function of the amount of computation in that node.

We also propose an algorithm to design a cloud network under given communication and computation constraints. For a given set of functions, we can find a network topology that maximizes the computation rate under the given computation

and communication constraints. We explore a greedy algorithm that designs a cloud network with a given network complexity, which we measure as the number of edges in the network. We compare its performance with a random algorithm, which outputs a network topology of the given complexity by examining the trade-off between network complexity and computation rate for both algorithms.

This thesis is organized as follows.

Chapter 1 provides an introduction to the topic of cloud computing and outlining the problem we are trying to solve.

Chapter 2 gives a recap of both my previous work done in 6.UAP and related work in the field.

Chapter 3 provides the framework we use for the problem. It is an adaptation of the framework used by Shah *et al.* [8], where we assume we are given a network and a set of functions to be computed over that network, and analyze its performance.

Chapter 4 introduces the Maximum Computation Cost (MCC) model, where the computation cost over the network is proportional to the largest total self-flows over any node. In effect, we are trying to spread the load more evenly over the network if there tends to be clusters of high computation.

Chapter 5 introduces the Linear Computation Cost (LCC) model, where we place computation costs on specific nodes in the network, proportional to the amount of self-flow through each node.

Chapter 6 discusses a possible algorithm for network design, where we are given the set of functions to be computed and their probability distribution, and we find a network that maximizes its total computation rate.

Chapter 7 goes over the results found in Chapters 4, 5, and 6 and presents areas of future research.

# Chapter 2

## Previous Work

### 2.1 UAP work

My UAP work was to use estimated graph entropies to find theoretical upper and lower bounds on computation rates. As an introductory model, we created a simulation to test bounds on communication and computation costs, and find an algorithm that gives the optimal ratio between the two types of costs. We began by coding a cloud simulation with only a few nodes, where the source nodes and compute nodes are objects in C++. A function is decomposed into coloring functions which are assigned to nodes. We defined color functions for each object that can be used in a lookup table to find the desired function value for each node. There are also functions that compute over the subset of data received. We started off with a easily verifiable function, a sort. The first round is defined as when subsets of the source nodes are input into compute nodes.

I created a working simulation that took in an adjacency matrix of the graph and created a parallel set of the source nodes. It then recursively created the next parallel set until it encountered an empty parallel set, at which point it returned the solution. We also put in an on-off matrix that indicated which node data is output to, for cases where a node has two children it can potentially send data to and we only want to output to one to avoid repetition. A parallel set is a set of jobs to be done concurrently. The structure checks if there are threads available in the threadpool,

and if there are, it takes one and assigns one of the jobs within the set to that thread. Once all the jobs in that parallel set have been assigned and completed, the parallel set returns the solution.

In the most basic version of our model, we start off with three source nodes, each of which randomly generates a list of integers of size  $n$ . These nodes are put in the first parallel set. As each node from this set is picked up by a thread, it queries the on-off matrix for the node it is supposed to output to, and adds that node to the next parallel set. Once this has been done to every node in the first parallel set, it is empty and returns that the job is done. We then pick up the next parallel set and do the same, except now this is a set of compute nodes. We again query for the output nodes and put them in the next parallel set, but then, we also query for the input nodes, and grab the corresponding lists of data. If a node has two parent nodes putting out data to it, the data lists are combined by concatenation. The data is sorted and set as output for the next parallel set of nodes. We calculate the computation time by collecting the timestamp before and after each parallel set runs, and calculate the communication time by taking  $\sum \log(n)$ , where  $n$  is the size of the data set transmitted from a parent node to its child. Therefore, the first parallel set has a communication time of  $3\log(n)$ , because each of the three source nodes transmits  $n$  integers to a child node in the next parallel set.

## 2.2 Related Work

A major part of my UAP work was to investigate information theoretic rate bounds for computing a function over a network, where the input is found at the sources and the output at the receiver. Körner introduced graph entropy [6], which Feizi *et al.* [5] used to investigate graph coloring approaches to this problem. Distributed computing and in-network computation are both topics that have had significant interest, and have lead the way to network flow techniques, [2], [9]). Shah *et al.* [8] also used this framework for function computation considering only communication constraints. He simplifies cloud computing as a set of functions being computed over a network, with

a probability distribution for how often each function is called. There are jointly communication and computation limitations in a cloud computing network that need to be taken into account during design. The links between nodes have communication capacities, which are our communication constraints, and different computations require different computation power, which we represent as computation costs.

Shah provided a linear optimization method to find the maximum computation rate of a function over a network, and an algorithm to route the flows to achieve that rate [8]. The flow conservation constraint does not hold for functional flows because when a computation is done at a node, the flow coming out of that node is no longer the same as the flow coming in. However, [8] circumvented this by defining a self-flow on each node for each computation type. These self-flows make up the difference in flow when a computation is done at that node, allowing for flow conservation. These self-flows are proportional to the amount of computation done at each node, and therefore by putting a cost on these flows we can simulate computation costs. This framework we modify further as discussed in Chapter 3.

Reference [7] also considered the problem of minimum-cost multicast over coded networks with a decentralized approach, and used a max norm approach that we shall discuss in Chapter 4.





# Chapter 3

## Framework

### 3.1 Network

We are using the framework described in [8]. The network  $\mathcal{N} = (V, E)$  is a directed acyclic graph, where  $V$  is the set of nodes, and  $E$  is the set of edges. Each edge  $(u, v)$  has a capacity of  $c_{(u,v)}$ .  $\mathcal{N}$  has  $k$  source nodes,  $\{n_1, \dots, n_k\}$ , that have data values  $\{X_i(l)\}_{l \geq 0}$ , where  $X_i(l)$  belongs to a finite alphabet. For source node  $i$ , we use  $X_i$  to represent its data. In cloud computing, source nodes can be considered as datasets. In this case,  $X_i$  represents data in the dataset  $i$ .  $N(v)$  is the set of neighbors of node  $v$  in the network.

The terminal node,  $n_t$ , receives the output of a set of functions from the source data. These functions are given as computation trees.

### 3.2 Computation Trees

A computation tree is a graph  $\mathcal{G} = (\zeta, \Gamma)$ , where  $\zeta$  is the set of nodes, and  $\Gamma$  is the set of directed edges. Each directed edge  $\theta_i$  represents a separate computation to be mapped to a node in the network.  $\{\theta_1, \dots, \theta_k\}$  are outgoing edges from source nodes that represent the input data and  $\theta_t$  is the incoming edge to the terminal node  $\zeta_t$ , representing the output of the given function the computation tree represents.

$\{\zeta_1, \dots, \zeta_k\}$  represents the source nodes. There are many possible computation

trees (shown in Section 3.3 for a given function, and, for the scope of this problem, we assume that a computation tree is given. If the given computation tree excludes certain function mappings, those mappings will not be considered.

For a specific computation  $\theta \in \Gamma$ ,  $tail(\theta)$  and  $head(\theta)$  represent the tail and the head nodes of that edge in the computation tree, respectively. A computation  $\theta_i$  is a parent of a computation  $\theta$  if  $tail(\theta_i) = head(\theta)$ . Children of a computation  $\theta$  are defined similarly. We refer to the set of parents and children of computation  $\theta$  as  $\Phi_{\uparrow}(\theta)$  and  $\Phi_{\downarrow}(\theta)$ , respectively.

### 3.3 Mapping

A given computation tree can be mapped over a network by mapping its computations (edges)  $\theta$  to different nodes of the network. A mapping  $M$  is feasible when all computations  $\theta$  can be computed by the data received from their parents in the network. In other words, each computation must be mapped to a node that has a path leading from the nodes mapped to  $\Phi_{\uparrow}(\theta)$ .

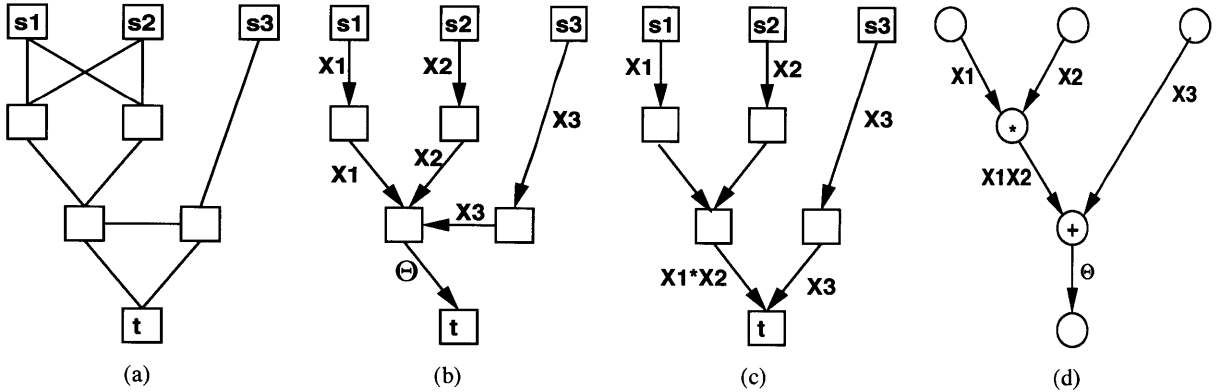


Figure 3-1: Mapping example: (a) Network  $\mathcal{N}$ . (b) A mapping of computation tree  $\mathcal{G}$  (d) on  $\mathcal{N}$ . (c) A different mapping of  $\mathcal{G}$  on  $\mathcal{N}$ . (d) A computation tree  $\mathcal{G}$  for the function  $\theta = X_1 * X_2 + X_3$ . [8]

Figure 3.3 shows a sample network  $\mathcal{N}$  and a possible computation tree for the function to be computed,  $\theta = X_1 * X_2 + X_3$ . Figure 3.3-(b) and (c) show possible

mappings of that computation tree on  $\mathcal{N}$ , one of many that can be found by our algorithm Embedding Edge LP, Figure 3-2.

---

**Algorithm 1:** Finding equivalent solution of the *Embedding-Edge LP* from a feasible solution of the *Node-Arc LP*.

---

**input :** Network graph  $\mathcal{N} = (V, E)$ , capacities  $c(e)$ , set of source nodes  $S$ , terminal node  $t$ , computation tree  $\mathcal{G} = (\Omega, \Gamma)$ , and a feasible solution to its *Node-Arc LP* that consists of the values of  $\lambda$ ,  $f_{uv}^\theta \forall \theta \in \Gamma, \forall uv \in E$ , and  $f_{uu}^\theta \forall \theta \in \Gamma, \forall u \in V$ .

**output:** Solution  $\{x(B) | B \in \mathcal{B}\}$  to the *Embedding-Edge LP* with  $\sum_{B \in \mathcal{B}} x(B) = \lambda$ .

Initialize  $x(B) := 0, B(\theta_i) = \emptyset$  (the null sequence),  $\forall B \in \mathcal{B}$  and  $\forall \theta_i \in \Gamma, \lambda' = 0$

**while**  $\lambda' \neq \lambda$  **do**

$z(t) := \lambda$  ;

$B(\theta_{|\Gamma|}) := t$  ;

**for**  $i := |\Gamma|$  **to** 1 **do**

$v := B(\theta_i)$  ; // valid, as  $B(\theta_i)$  has of only one node at this step

$u :=$  an element in  $N^+(v)$  such that  $f_{uv}^{\theta_i} > 0$  ;

**if**  $u \neq v$  and  $u \in B(\theta_i)$  **then**

// A cycle of redundant flow found: remove the flow from all the edges in the cycle

Let  $P$  be the path in  $B(\theta_i)$  upto the first appearance of  $u$  in it;

Delete  $P$  from  $B(\theta_i)$ . ;

$y := \min_{u'v' \in \{uv\} \cup P} (f_{u'v'}^{\theta_i})$  ;

$f_{u'v'}^\theta := f_{u'v'}^\theta - y \forall u'v' \in \{uv\} \cup P$

**end**

**else**

$z(u) := \min(z(v), f_{uv}^{\theta_i})$  ;

**end**

**if**  $u \neq v$  **then**

Prefix  $u$  in  $B(\theta_i)$  ;

$v := u$  ;

Jump to the second statement inside the **for** loop ;

**end**

**else**

$B(\eta) := u, \forall \eta \in \Phi_{\uparrow}(\theta_i)$  ;

**end**

**end**

$x(B) := z(s_1)$  ; // Flow extracted on  $B$

$\lambda' := \lambda' + x(B)$  ; // Total flow extracted

// Remove  $x(B)$  amount of flow from all the edges in  $B$ .

$f_{u'v'}^\theta := f_{u'v'}^\theta - x(B) \forall \theta \in \Gamma$  and  $\forall u'v' \in B(\theta)$  ;

// Remove  $x(B)$  amount of flow from all the relevant self-loops.

$f_{v'v'}^\theta := f_{v'v'}^\theta - x(B) \forall \theta \in \Gamma$  and  $v' = \text{start}(B(\theta))$  ;

**end**

---

Figure 3-2: Embedding Edge LP [8]

### 3.4 Algorithms

Note that, one computation tree can have several feasible mappings over a network. Define  $\mathcal{M}$  as the set of all mappings of a computation tree  $\mathcal{G}$  on our network  $\mathcal{N} = (V, E)$ . To obtain the maximum computation rate  $R$ , we need to find the set of mappings that obey capacity constraints of the network that sum to achieve  $R$ . We define  $R$  with a discrete model, where the sources put out discrete packets of data, and the terminal node receives discrete packets of the function output.

In Figure 3-4, Shah *et al.* gives a basic linear program that finds the maximum computation rate  $\lambda$  for a given computation tree over network  $\mathcal{N}$ . It also outputs the amount of flow of computation type  $\theta_1, \dots, \theta_{|\Gamma|}$ , where  $|\Gamma|$  is the number of computations, over every edge  $e \in E$  in the network. A computation flow is the amount of data being sent over an edge, which is easier to visualize when dealing with a discrete model.  $f_{uv}^\theta$  represents the flow from node  $u$  to  $v$  of computation type  $\theta$ .  $N(v)$  is defined as the set of nodes neighboring  $v$  and  $N'(v)$  is the set  $\{N(v) \cup v\}$ .

However, the scope of our problem is beyond this linear program. We want our algorithm to be able to compute multiple functions simultaneously on the network, and to give us the maximum total rate for both, weighted with their probability distribution. The probability distribution is the probability of each function being called by the network, and therefore we want to maximize the total average computation rate  $R_F = \sum_{i=1}^n p_i \lambda_i$ . Shah's algorithm also does not take into account computation and communication costs found in cloud computing. The multiple functions with probability distribution problem is solved as shown below for a set of functions  $\Theta = \{\Theta_1, \dots, \Theta_n\}$  given computation trees  $\mathcal{G}_i = (\Omega_i, \Gamma_i)$  for  $i = \{1, \dots, n\}$ , where  $\mathcal{G}_i$  is the given computation tree for function  $\Theta_i$ . The set of functions  $\Theta$  is given probability distribution  $\{p_1, \dots, p_n\}$ .

**Algorithm 1.**  $\max \sum_{i=1}^n p_i \lambda_i$   
*subject to*

$$f_{vv} + \sum_{u \in N(v)} f_{vu}^\theta - \sum_{u \in N'(v)} f_{uv}^\theta = 0, \forall \theta \in \Gamma_i \setminus \{\theta_{|\Gamma_i|}\}, \quad \forall \Gamma_i \in \Gamma \text{ and } \forall \eta \in \Phi_\downarrow(\theta). \quad (3.1)$$

$$\sum_{u \in N(v)} f_{vu}^{\theta_{|\Gamma_i|}} - \sum_{u \in N'(v)} f_{uv}^{\theta_{|\Gamma_i|}} = \begin{cases} -\lambda_i & v = t \\ 0 & \text{otherwise} \end{cases}, \quad \text{for } i \in \{1, \dots, n\} \quad (3.2)$$

$$f_{vv}^{\theta_i} = \begin{cases} \lambda_i & v = s_i \\ 0 & \text{otherwise} \end{cases}, \quad \text{for } i \in \{1, \dots, n\} \quad (3.3)$$

$$\sum_{\theta \in \Gamma} (f_{uv}^\theta + f_{vu}^\theta) \leq c(uv), \quad \forall uv \in E. \quad (3.4)$$

$$f_{uv}^\theta \geq 0, \quad \forall uv \in E, \forall \theta \in \Gamma_i \text{ and } \forall \Gamma_i \in \Gamma \quad (3.5)$$

$$f_{uu}^\theta \geq 0, \quad \forall u \in V, \forall \theta \in \Gamma_i \text{ and } \forall \Gamma_i \in \Gamma \quad (3.6)$$

$$\lambda_i \geq 0. \quad \text{for } i \in \{1, \dots, n\}. \quad (3.7)$$

Different ways of combining computation and communication costs will be discussed in Chapters 4 and 5. Computation costs are costs placed on certain, or all, nodes on a network for certain computations. These can represent inefficient machines or if certain computations are heavier than others. Communication costs are the capacities of each edge, a fixed ceiling that is provided.

We will explore the first constraint, the functional conservation of flows, further. The usual definition of flow conservation does not hold in the case of function flows, as shown in Figure 3-3. When a computation involving  $f_{(1,3)}^\theta$  and  $f_{(2,3)}^\theta$  occurs at node 3, then

$$f_{(1,3)}^\theta + f_{(2,3)}^\theta \neq f_{(3,4)}^\theta$$

To achieve flow conservation, we introduce self-flows,  $f_{(i,i)}^\eta$  at each node, so that our new conservation rule is,

$$f_{(1,3)}^\theta + f_{(2,3)}^\theta = f_{(3,4)}^\theta + f_{(3,3)}^\eta.$$

These self-flows represent the amount of computation done at each node, although different computations may have different amounts of self-flow. We shall use these self-flows when including computation constraints, as shown in Chapters 4 and 5.

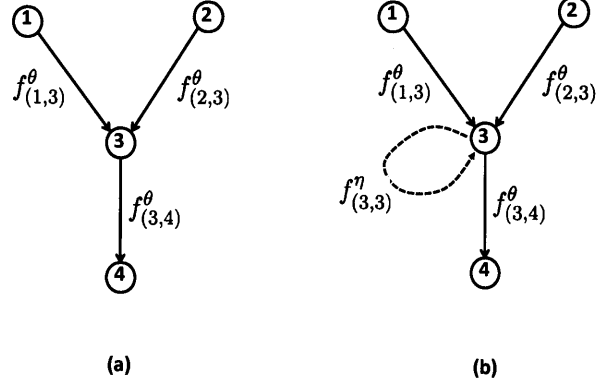


Figure 3-3: (a) An example that shows the flow conservation constraint does not hold in the function computation setup. (b) By adding a self-loop with infinite capacity to node 3, a modified flow conservation constraint holds in this case.

Furthermore, we know basic bounds for the total computation rate over a network given  $k$  sources and the min-cut rate of our network,  $\gamma$ .

**Theorem 2.** *Given a set of functions with  $k$  sources and a network  $\mathcal{N}$  with min-cut rate  $\gamma$ , the total computation rate  $R_F$  has bounds,*

$$\frac{\gamma}{k} \leq R_F \leq \gamma$$

*Proof.* The lower bound represents the simplest scenario, a centralized scheme where all computations are done at the terminal node. In this case, each source can send  $\frac{\gamma}{k}$  to the terminal node, following the min-cut max-flow theorem for multicast networks [1], giving us a total computation rate of  $\frac{\gamma}{k}$ .

The upper bound is achievable by envisioning a virtual node connected to all  $k$  sources with infinite capacity links. We also assume this virtual node has infinite computational capacity. In this case, the limiting factor is the min-cut rate of the network,  $\gamma$ , and again using the min-cut max-flow theorem, the upper bound of  $R_F$  is  $\gamma$ . □

---

**Node-Arc LP:** Maximize  $\lambda$  subject to following constraints any node  $v \in V$

1. Functional conservation of flows:

$$f_{vv}^\eta + \sum_{u \in N(v)} f_{vu}^\theta - \sum_{u \in N'(v)} f_{uv}^\theta = 0, \quad \forall \theta \in \Gamma \setminus \{\theta_{|\Gamma|}\} \text{ and } \forall \eta \in \Phi_\downarrow(\theta). \quad (3)$$

2. Conservation and termination of  $\theta_{|\Gamma|}$ :

$$\sum_{u \in N(v)} f_{vu}^{\theta_{|\Gamma|}} - \sum_{u \in N'(v)} f_{uv}^{\theta_{|\Gamma|}} = \begin{cases} -\lambda & v = t \\ 0 & \text{otherwise} \end{cases} \quad (4)$$

3. Generation of  $\theta_l \forall l \in \{1, 2, \dots, \kappa\}$ :

$$f_{vv}^{\theta_l} = \begin{cases} \lambda & v = s_l \\ 0 & \text{otherwise} \end{cases} \quad (5)$$

4. Capacity constraints

$$\sum_{\theta \in \Gamma} (f_{uv}^\theta + f_{vu}^\theta) \leq c(uv), \quad \forall uv \in E. \quad (6)$$

5. Non-negativity constraints

$$f_{uv}^\theta \geq 0, \quad \forall uv \in E \text{ and } \forall \theta \in \Gamma \quad (7)$$

$$f_{uu}^\theta \geq 0, \quad \forall u \in V \text{ and } \forall \theta \in \Gamma \quad (8)$$

$$\lambda \geq 0. \quad (9)$$


---

Figure 3-4: Node Arc LP [8]

Figure 3-2 details a linear program that finds the set of mappings that achieve rate  $\lambda$ . It does so by starting at terminal node  $t$  and exploring upwards to neighboring nodes while checking if these paths contain nonzero flow of type  $\theta_{|\Gamma|}$ . When the algorithm hits a node  $u$  where none of the edges to neighbors contain nonzero flow of type  $\theta_{|\Gamma|}$ , the algorithm assigns that computation to  $u$ , saves the path and the smallest flow through that path, and assigns  $u$  as the starting point for all computations in the set  $\Phi_\uparrow(\theta_{|\Gamma|})$ . The algorithm then repeats for computations  $\theta_{|\Gamma|-1}, \dots, \theta_1$ . Every run of the `for` loop, as shown in Figure 3-5, outputs an embedding and its rate, the smallest amount of flow on each path (the bottleneck), and the algorithm only stops when it achieves the maximum computation rate given by Node Arc LP.

```

for  $i := |\Gamma|$  to 1 do
   $v := B(\theta_i)$ ; // valid, as  $B(\theta_i)$  has of only one node at this step
   $u :=$  an element in  $N'(v)$  such that  $f_{uv}^{\theta_i} > 0$ ;
  if  $u \neq v$  and  $u \in B(\theta_i)$  then
    // A cycle of redundant flow found: remove the flow from all
    // the edges in the cycle
    Let  $P$  be the path in  $B(\theta_i)$  upto the first appearance of  $u$  in it.;
    Delete  $P$  from  $B(\theta_i)$ .;
     $y := \min_{u'v' \in \{uv\} \cup P} (f_{u'v'}^{\theta_i})$ ;
     $f_{u'v'}^{\theta} := f_{u'v'}^{\theta} - y \forall u'v' \in \{uv\} \cup P$ 
  end
  else
     $z(u) := \min(z(v), f_{uv}^{\theta_i})$ ;
  end
  if  $u \neq v$  then
    Prefix  $u$  in  $B(\theta_i)$ ;
     $v := u$ ;
    Jump to the second statement inside the for loop;
  end
  else
     $B(\eta) := u, \forall \eta \in \Phi_{\uparrow}(\theta_i)$ ;
  end
end

```

Figure 3-5: for loop in Embedding Edge LP [8]



# Chapter 4

## A Maximum Computation Cost Model

### 4.1 Description

In the Maximum Computation Cost Model, we are analyzing the behavior of computation flows over a network under certain limitations. The limitation is the maximum computation power constraint in a network. In this case, we show how our framework allows redistribution of computation flows over the network to achieve a better total computation rate. However, a max-norm is not everywhere differentiable, and therefore cannot be modeled as a convex program. Instead, we use the  $l^k$ -norm approximation [7] which converges to max-norm as  $k \rightarrow \infty$ . The new convex program constraints are below.

**Algorithm 3.**

$$\max \quad \sum_{i=1}^n p_i \lambda_i - \delta \left( \sum_{v \in V} \left( \sum_{\theta \in \Gamma} f_{vv}^{\theta} \right)^p \right)^{\frac{1}{p}}$$

*subject to equations 3.1-3.7.*

$\delta$  is the maximum computation cost parameter, a non-negative weight on the maximum self-flow that controls how evenly the computation is spread through the

network.

## 4.2 Simulations

In our simulations, we consider two functions to be computed at the receiver:

$$f_1 = X_1 * X_2 + X_3 * X_4 + X_5 * X_6$$

$$f_2 = X_1 * X_2 * X_3 + X_4 * X_5 * X_6.$$

The computation trees for  $f_1$  and  $f_2$  are shown in Figures 4-1 and 4-2. There are many possible computation trees for these functions, but the trees in Figures 4-1 and 4-2 encompass all possible mappings, because all possible computation orderings are achievable with these computation trees.

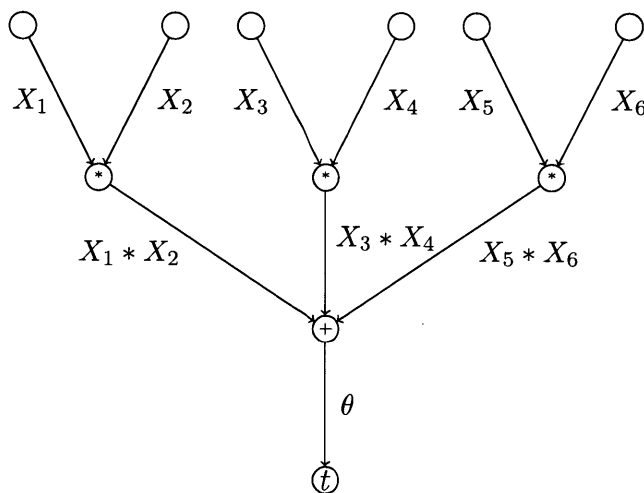


Figure 4-1: Function 1.

We demonstrate the effect of adding an approximation of Maximum Computation Cost (MCC) to our optimization problem, where we maximize the computation rate  $\lambda$  with an added  $l^k$ -norm cost,  $\delta * (\sum_{i,j} g_{i,j}^k)^{\frac{1}{k}}$  where  $g_{i,j}$  is the sum of all flow over edge  $(i, j)$  for all edges in the network and  $\delta$  is a control variable. As  $k \rightarrow \infty$  we are effectively putting a constraint on the maximum edge flow in the network [7].

The initial network for this simulation  $\mathcal{N}_1$  is shown in Figure 4-3. All edges have

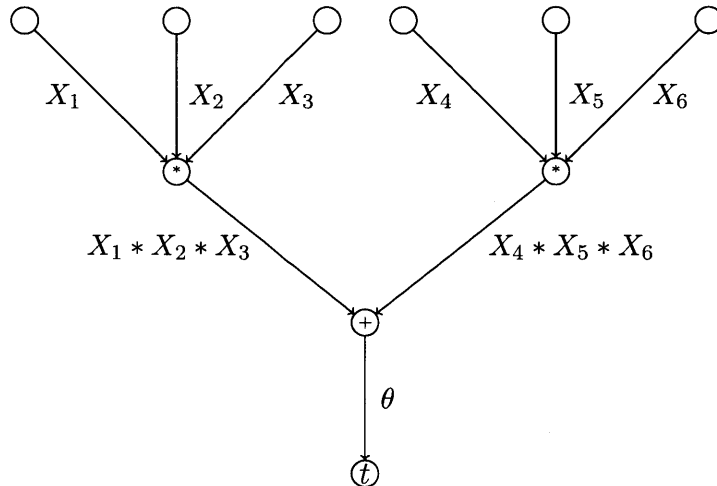


Figure 4-2: Function 2.

capacity 10 except the edge from  $n_{15}$  to the terminal node,  $n_t$ , has capacity 1. We chose edge capacities that cause flows to be denser in certain sections of the network than others. The min-cut rate of this network is therefore 11. We are computing  $f_1$  and  $f_2$  over the network simultaneously and with equal weights.

### 4.3 Results

Figure 4-4 shows the change in distribution of computation over the nodes in the network when we use the MCC constraint vs. when we do not use it. In this simulation we are using  $\delta = .1$  and  $k = 15$ . We achieved these values through trial and error. A  $\delta$  value too small did not affect flows in the network, and too large a value caused all flows to go to zero. The  $k$  is sufficiently large to provide us a reasonably accurate approximation of the max-norm. Using higher values did not change the computation costs, which means our  $k$  was able to find the maximum computation flow. We can see that the computation rate in nodes in the second layer,  $\{n_7, \dots, n_{12}\}$  increases in the model with the MCC parameter, spreading the computation load through the network more effectively than in the case without the MCC parameter.

Figure 4-5 shows the total computation rate over the network with varying  $\delta$  values. When  $\delta = 0$  we can achieve the maximum computation rate of 22, but as

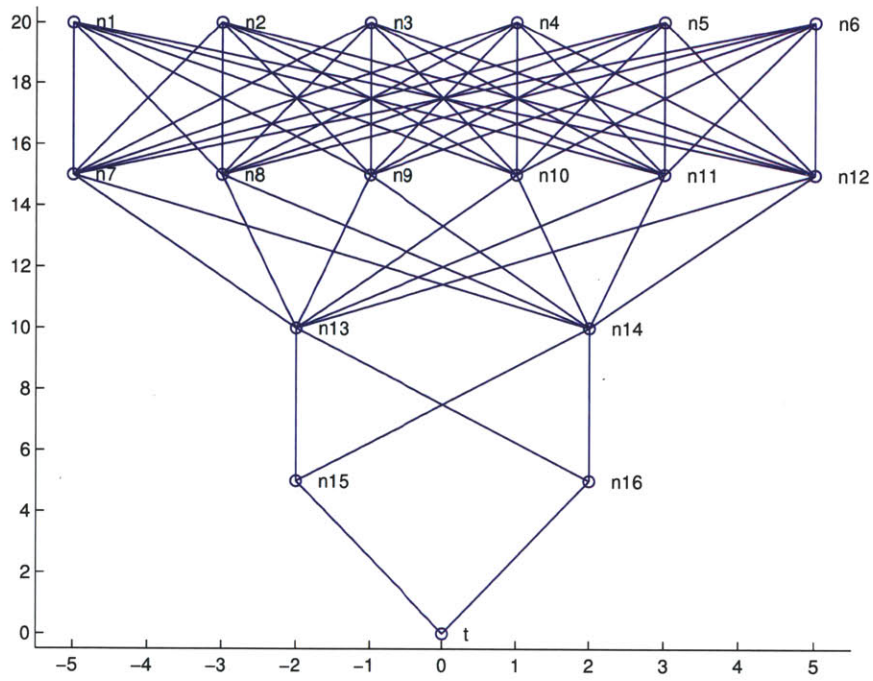


Figure 4-3: Initial communication network  $\mathcal{N}_1$ .

$\delta$  increases, the computation rate decreases, eventually to zero when  $\delta$  becomes too large.

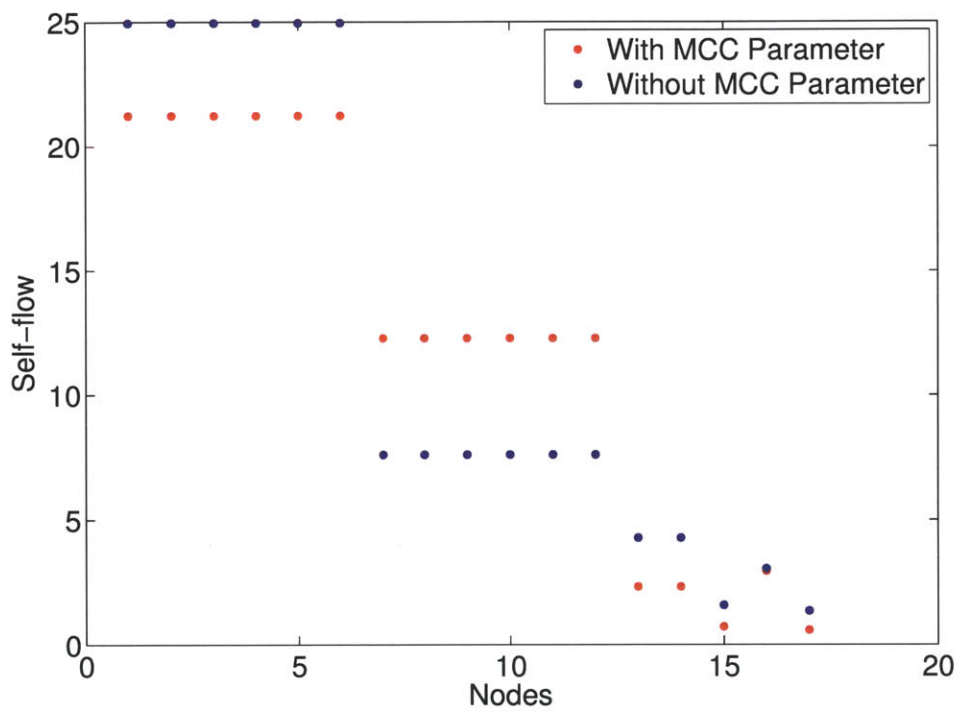


Figure 4-4: Self-Flow over nodes in the network  $\mathcal{G}_1$

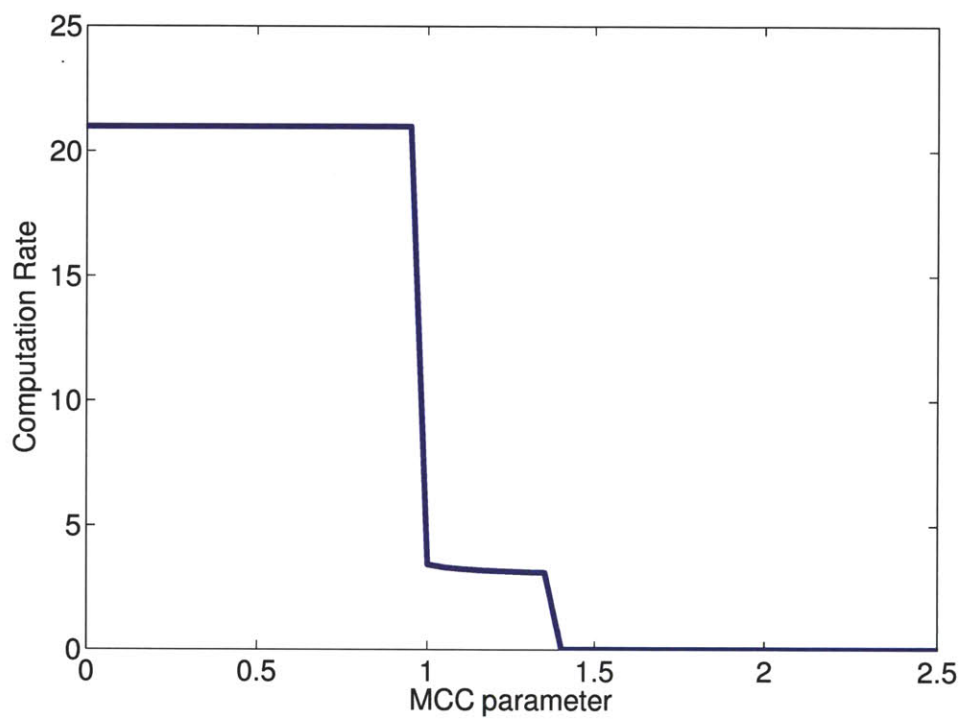


Figure 4-5: Total Computation Rate of  $f_1$  and  $f_2$  vs. MCC constraint.

# Chapter 5

## A Linear Computation Cost Model

### 5.1 Description

The Linear Computation Cost Model consists of modifying the original Node Arc LP in Figure 3-4 by adding costs to specific nodes proportional to the amount of self-flow through that node.

**Algorithm 4.**

$$\max \quad \sum_{i=1}^n p_i \lambda_i - \sum_{v \in V} \delta_v \left( \sum_{\theta \in \Gamma} f_{vv}^{\theta} \right)$$

*subject to equations 3.1-3.7.*

$\delta_v, v \in V$  are the linear computation cost parameters, non-negative weights on the total self-flow through each node  $v \in V$  that controls the computation cost. Unlike the MCC model, this is a linear function and therefore can be solved with a linear program.

## 5.2 Simulations

In this section, we consider the communication network  $\mathcal{N}_2(V_2, E_2)$  as depicted in Figure 5-1 and the receiver node only demands to compute  $f_1$  as defined in Figure 4-1. The network in Figure 5-1 has all edge capacities of 10, and we can see the bottleneck is at the minimum cut of the network, across the edges from  $n_9$  to  $t$  and  $n_{10}$  to  $t$ , and the min-cut rate of this network is  $\gamma = 20$ . We place costs on nodes  $n_7 - n_{10}$  to analyze the redistribution of flow and the effect on computation rate.

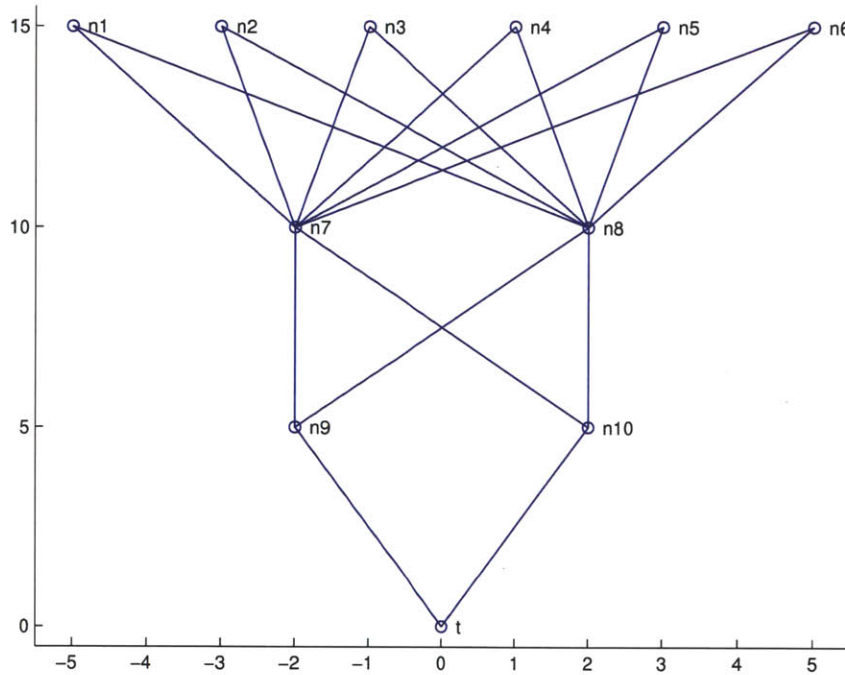


Figure 5-1: Initial communication network  $\mathcal{N}_2$ .

## 5.3 Results

Figure 5-2 shows the redistribution of flows occurring when all computations at  $n_8$  are given a Linear Computation Cost (LCC) constraint of 10. With no computation costs, flow is distributed equally among nodes  $n_7$ ,  $n_8$ ,  $n_9$ , and  $n_{10}$ . However, by



making  $n_8$  an expensive node at which to compute, flow is redistributed and getting passed on further to  $n_9$  and  $n_{10}$ .

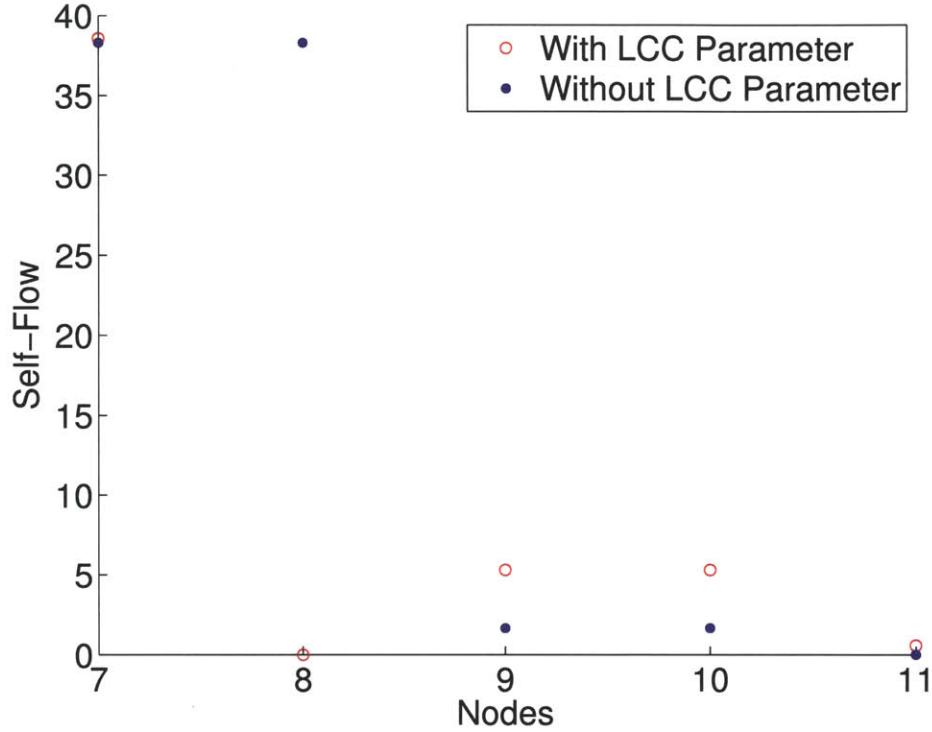


Figure 5-2: Self-Flow over nodes in the network affected by the bottleneck  $\mathcal{G}_2$ .

Figure 5-3 shows the effect on computation rate of varying LCC constraint  $\delta$  from 0 to 0.5 for all computations at  $n_7 - n_{10}$  (the bottleneck). We have placed costs on all nodes necessary for computation except the terminal node. We can see that, when  $\delta = 0$ , we achieve the upper bound for  $R_F$  as given by Theorem 2. As the LCC constraint increases, all of the computation gets redistributed to the terminal node, a more centralized scheme, which limits the computation rate to the bottleneck capacity of the network over the number of sources.

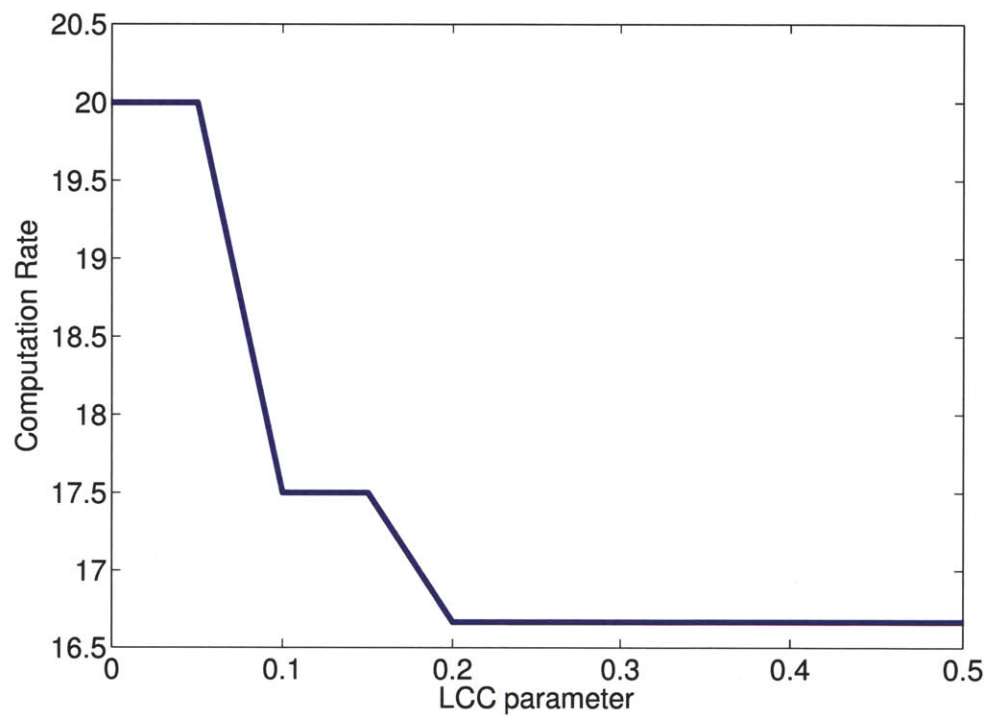


Figure 5-3: Computation Rate vs. LCC constraint on a subset of nodes.

# Chapter 6

## Network Design

In this chapter, we shall outline and simulate an algorithm that designs a network that maximizes computation rate for a specific set of functions and their probability distribution while minimizing network complexity.

**Definition 5.** *Given a network  $\mathcal{N} = (V, E)$ , where  $V$  is the set of nodes and  $E$  the set of all possible edges in  $\mathcal{N}$ , a sparse network  $\mathcal{N}_s = (V, E_s)$  is optimal for a set of functions  $\mathcal{F}$  if  $E_s \subseteq E$  and its computation rate  $R_s$  is maximized over all possible subsets of  $E$  of size  $|E_s|$ .*

In our algorithm, we are defining network complexity by the number of edges in our network. Therefore, an optimal network design of a given complexity is defined as the network with a given number of edges  $e$  with the highest computation rate of the set of all network topologies with a set of  $e$  edges that are a subset of all possible edges of our initial network.

### 6.1 Greedy Algorithm

In this section, we are sparsifying a dense network with a greedy approach. At each iteration, we run Node Arc LP and take away the edge with the smallest ratio of flow to capacity, which is the locally optimal solution to improving the ratio of computation rate and network complexity. The algorithm stops when we have reached the desired

number of edges. This algorithm can be combined with the models described in Chapters 4 and 5, but for the simulation described in 6.2 we stick with the basic model as described in the framework in Section 3.4.

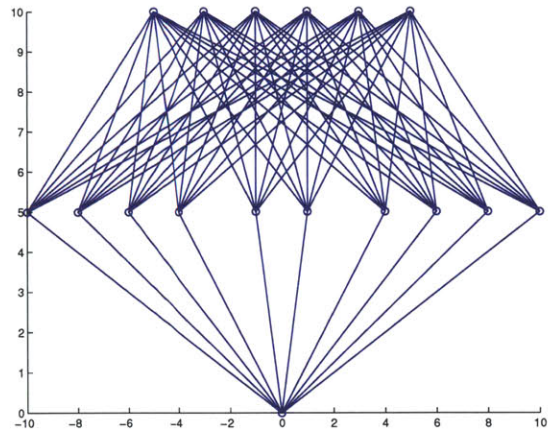
**Algorithm 6.** *The greedy algorithm is as follows,*

1. Use Node Arc LP to compute all flows  $f_{ij}^\theta$  for each edge for the network  $\mathcal{N}^r = (V, E^r)$ , where  $E^r$  is the set of possible edges in  $\mathcal{N}$ .
2. Eliminate the edge  $(u, v)$  with the smallest ratio of flow to capacity,  $(u, v) = \arg \min_{(u,v) \in E^r} \frac{g_{ij}^r}{c(i,j)}$  where  $g_{ij}^r = \sum_{\theta \in \Gamma} f_{ij}^\theta$  and  $c(i, j)$  is the capacity of edge  $(i, j)$ .
3. Update the set of edges:  $E^{r+1} = E^r \setminus \{(u_1, v_1)\}$ .
4. If  $|E^{r+1}|$  is the desired number of edges, terminate. Else, repeat.

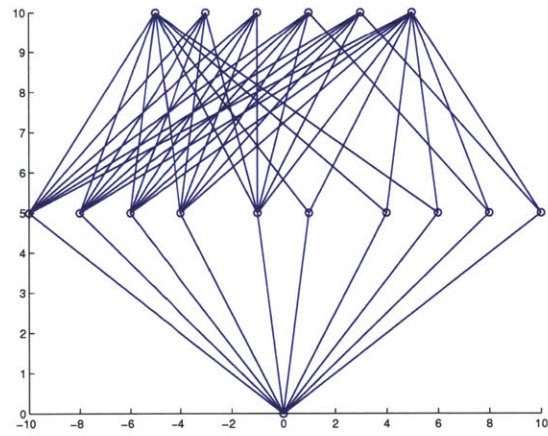
## 6.2 Simulations

We start out with an initial communication network  $\mathcal{N}_3 = (V, E)$  with 6 sources, one layer of 10 nodes, and a terminal node as shown in Figure 6-1(a). The set of possible edges are all possible pairings of nodes that are between layers. Therefore possible edges include and are limited to: all edges between  $\{n_1, \dots, n_6\}$  and  $\{n_7, \dots, n_{16}\}$  (but none within sets) and all edges between  $\{n_7, \dots, n_{16}\}$  and  $n_t$ .

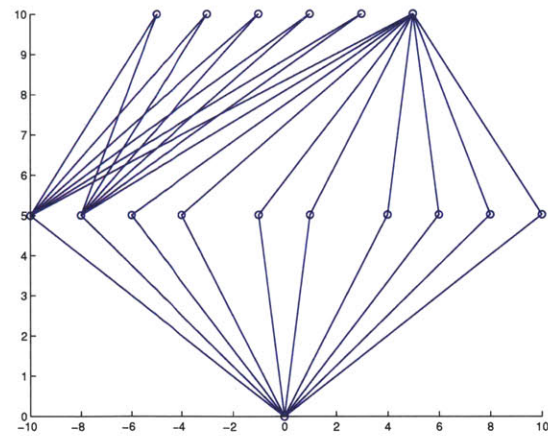
We calculate the ratio of flow to capacity along each edge, and remove the edge with the smallest ratio. No nodes on the same layer or non-adjacent layers are connected, as shown in Figure 6-1(a).



(a) Initial communication network  $\mathcal{N}_3$



(b) Iteration 20. Rate 1: 5.0411, Rate 2: 4.9589



(c) Iteration 40. Rate 1: 3.4555, Rate 2: 3.0445

Figure 6-1: Network Sparsification

## 6.3 Results

Figures 6-1(b) and 6-1(c) show the networks after iterations 20 and 40, where the original network in Figure 6-1(a) has been reduced by 20 and 40 edges respectively. Also in the description are the computation rates for  $f_1$  and  $f_2$ .

Figure 6-2 shows the total computation rate over  $\mathcal{N}_3$  as an edge is taken away in each iteration. We compare its performance with a random algorithm, where at each iteration a random edge is taken from the network. We run the random algorithm 5 times and average over the runs to account for scenarios where edges that are taken out cause the computation rate to drop to zero. The greedy curve does not drop at every iteration, because, in cases where an edge is removed, it is still possible to redistribute flows, which means the network still has an excess of edges that, when removed, do not diminish the min-cut of the network. If an edge is removed in a portion of the network that is not the min-cut, this does not affect rate because the rate is determined by the min-cut, by the min-cut max-flow theorem [1].

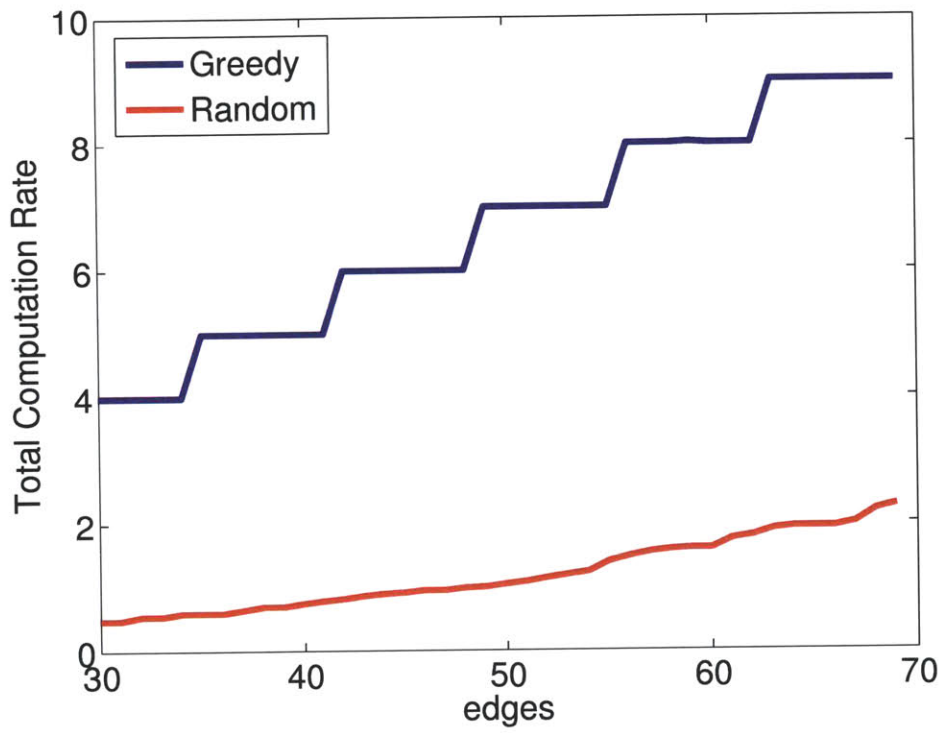


Figure 6-2: Computation rates summed over  $f_1$  and  $f_2$ .





# Chapter 7

## Conclusions

### 7.1 Project Accomplishments

In this thesis we have outlined a framework to find the maximum computation rate of a set of functions for a given network with communication limits (capacity constraints) in Chapter 3. We then discuss frameworks with different types of computation limitations, the Linear Computation Constraint Model (Chapter 5), and the Maximum Computation Constraint Model (Chapter 4). The LCC model uses a linear function at each node to represent computation cost, while the MCC model uses the maximum computation over all nodes in the network. Since this is not a convex problem, a relaxation of  $l^p$ -norm is used as a substitute for the max-norm [3]. While this is useful for achieving the best computation rate out of a given network, a better and more useful optimization problem to solve for cloud computing is how to design a network specialized for a set of functions. This problem is broached with a sub-optimal greedy algorithm, as described in Chapter 6. We found we were able to sparsify a given network while maintaining the same computation rate, and obtain a locally optimal sparsifying solution at each iteration. However, future work still needs to be done to find an optimal algorithm for designing a cloud network, as defined in Definition 5.

## 7.2 Future Work

We have tried various algorithms to achieve optimal sparse networks using the basic idea of our computation cost models, and adding a sparsity term  $\delta$  to minimize the number of edges in the network. We added it as a cost to all the flows over the network, excluding self flows. Unfortunately, the network is very sensitive to  $\delta$ , and flows go to zero too quickly to obtain a sparse network. More work will need to be done to find an optimal algorithm, but another direction for optimizing cloud networks is determining the physical location of source data. Given a set of possible data locations, we can create a cost model that represents their distances from each other to find the model with the highest computation rate.

In this thesis we have only analyzed two cost models, the LCC model and the MCC model. However, there are other types of costs in cloud computing to consider and model when routing flows. Common costs to model are network congestion and slower nodes with higher computation costs. A variation of the MCC model that can be used is a convex cost function, where the cost is proportional to the amount of self-flow through each node. Another aspect of to improve performance we did not address is functional compression. There is a certain amount of data distortion that occurs with compression and decompression of data, and therefore we do not always need an exact rate of computation when analyzing a network. We should investigate alternative methods, like graph coloring, to our convex programming solution and analyze the performance vs. accuracy of using information theory tools to approximate rates over the cloud.

Function flows can also be useful in dealing with privacy issues in the cloud. To protect private information and filter it from the results requested at the receiver of a network (e.g. pulling information from insurance records), we limit the areas of the network that Untrusted areas of the network can have limits on the type of information they can access, which can be represented using the LCC model with  $\delta = \infty$  on the computations requiring the sensitive information.

Future work can also be done to analyze different computation trees for functions.

A function has multiple computation trees, and while we have addressed that it is best to choose a computation tree that encompasses all mappings on a network, there also may be runtime advantages to having computation trees with similar topologies when dealing with multiple functions, even if it means excluding some mappings.



# Appendix A

## Code

### A.1 Node Arc LP code

```
1 function [ MaxFlow, SparseNet, flows, X, SelfFlows, edgeflows,
    FlowConsConstraintA, FlowConsConstraintB] = NodeArcLPa( CommNet,
    CompWeights, FuncWeights, sources, t, CompTrees, Capacity)
% Calculates flow for a set of functions with associated weights
3 % CommNet is a adjacency matrix representing the Communication Network.
% NetEdges is the list of directed edges in the network
5 % CompWeights is a set of arrays that are the penalties on each
    computation edge for computation cost.
% FuncWeights = weight associated with each function
7 % Sources is an set of arrays of integer values specifying the nodes
    that are sources for each function.
% t is an array of indices of the terminal nodes for each function.
9 % CompTrees is a set of matrices of the Computation Tree, where there is
    a +1 value if the row node is the predecessor of the col node.
% Capacity: capacity of each edge on the communication network
11 % MAXIMIZE: rate-sum(weight*rate of edge)-sparsity constant*sum(all edge
    flows)
    nodes = [];
13 selfflows = [];
    NumFunctions = length(FuncWeights);
15 [rows, cols]=size(CommNet);

17 flows = CommNet;
    for row = 1:rows,
19         for col = 1:cols,
                if row==col,
21                     flows(row, col)=1;
                end
23         end
    end
25 x=find(flows); %All nodes and edges in the network that have flow
```

```

27 %get the number of computation edges in each CompTree by counting
    %the number of parent nodes
29 CompEdges=zeros(1,NumFunctions);
    totalWeights=[];
31 sparseDesign=[];
    X=zeros(1,NumFunctions);
33 for index = 1:NumFunctions,
        CompEdges(index) = length(find(CompTrees{index}));
35     X(index)=CompEdges(index)*length(x);
        totalWeights = [totalWeights; zeros(X(index),1)];
37     selfflows = [selfflows; zeros(X(index),1)];
        sparseDesign = [sparseDesign; ones(X(index),1)];
39     for theta=1:CompEdges(index),
            for j=1:rows,
41                 sparseDesign(Matrix2Vector(j,j,flows,theta,index,X))=0;
                    selfflows(Matrix2Vector(j,j,flows,theta,index,X))=j;
43             end
        end
45     weights = CompWeights{index};
        [j,~] = size(weights);
47     for i=1:j,
            totalWeights(Matrix2Vector(weights(i,2),weights(i,2), ...
49                 flows,weights(i,3),index,X))=weights(i,1);
        end
51 end

53 totalX=sum(X);
    %minimizing over computation costs, and computation rates given function
55 %weights
    f=[totalWeights; -ones(NumFunctions,1).*FuncWeights'];
57 %FlowConservation
    FlowConsConstraintA = [];
59 for func=1:NumFunctions,
        for node = 1:rows,
61             for theta = 1:(CompEdges(func)-1),
                    newRow = zeros(1, totalX+NumFunctions);
63                 for eta = GetSuccessors(theta, CompTrees{func}),
                            newRow(Matrix2Vector(node, node, flows, eta, func,X)) =
                                1;
65                 end
                    for neighbor = GetNeighbors(node, CommNet,[]),
67                         newRow(Matrix2Vector(node, neighbor, flows, theta, func,
                                X)) = 1;
                            newRow(Matrix2Vector(neighbor, node, flows, theta, func,
                                X)) = -1;
69                         end
                    newRow(Matrix2Vector(node, node, flows, theta, func, X)) =
                        -1;
71                 FlowConsConstraintA = [FlowConsConstraintA; newRow];
            end
63         end
73     end
end
75 FlowConsConstraintB = zeros(length(FlowConsConstraintA(:,1)), 1);

```

```

77 %TConservation
   TConservationA = [];
79 for func = 1:NumFunctions,
   for node = 1:rows,
81     newRow = zeros(1,totalX+NumFunctions);
   for neighbor = GetNeighbors(node, CommNet, []),
83     newRow(Matrix2Vector(node, neighbor, flows, CompEdges(func),
        func, X)) = 1;
        newRow(Matrix2Vector(neighbor, node, flows, CompEdges(func),
        func, X)) = -1;
85     end
   newRow(Matrix2Vector(node, node, flows, CompEdges(func), func, X
        )) = -1;
87     if node == t(func),
        newRow(totalX+func) = 1;
89     end
   TConservationA = [TConservationA; newRow];
91 end
end
93 TConservationB = zeros(rows*NumFunctions, 1);

95 %Generation of theta_l
   totalSources = [];
97 for i=1:NumFunctions,
   totalSources = [totalSources, sources{i}];
99 end
   numSources = length(totalSources);
101 ThetaGenerationA = [];
   ThetaGenerationB = zeros(rows*numSources, 1);
103 row = 0;
   for node = 1:rows,
105     for func = 1:NumFunctions,
   for theta = 1:length(sources{func}),
107     newRow = zeros(1,totalX+NumFunctions);
   if node==sources{func}(theta),
109     newRow(totalX+func) = -1;
   end
111     newRow(Matrix2Vector(node, node, flows, theta, func, X)) =
        1;
   ThetaGenerationA = [ThetaGenerationA; newRow];
113     end
   end
115 end

117 EQConstraintA = [FlowConsConstraintA; TConservationA; ThetaGenerationA];
   EQConstraintB = [FlowConsConstraintB; TConservationB; ThetaGenerationB];
119

121 %Capacity Constraints (LEQ)
   CapConstraintA = zeros(length(Capacity), totalX+NumFunctions);
   rowcount = 0;
123 for row = 1:rows,
   for col = 1:row,
125     if CommNet(row,col)~= 0 && row>col,
        rowcount = rowcount + 1;

```



```

127         for func = 1:NumFunctions,
129             for theta = (1:CompEdges(func)),
                CapConstraintA(rowcount, Matrix2Vector(row, col,
                    flows, theta, func, X))=1;
                CapConstraintA(rowcount, Matrix2Vector(col, row,
                    flows, theta, func, X))=1;
131             end
            end
133         end
        end
135     end
    CapConstraintB = Capacity;
137
138     %Non-negativity Constraints (GEQ)
139     NonNegConstraintA = -1*eye(totalX+NumFunctions, totalX+NumFunctions);
    NonNegConstraintB = zeros(totalX+NumFunctions, 1);
141
142     size(CapConstraintA);
143     size(CapConstraintB);
144     size(NonNegConstraintA);
145     size(NonNegConstraintB);
    LEQConstraintA = [CapConstraintA; NonNegConstraintA];
147     LEQConstraintB = [CapConstraintB; NonNegConstraintB];
149
150     %Run linear program
    MaxFlow = linprog(double(f), double(LEQConstraintA), double(
        LEQConstraintB), double(EQConstraintA), double(EQConstraintB));
151
152     %Generate sparse network
153     SparseNet = zeros(size(CommNet));
155
156     for row = 1:rows,
        for col = 1:rows,
157             if CommNet(row, col)~= 0,
                for func = 1:NumFunctions,
159                     for theta = 1:CompEdges(func),
                        SparseNet(row, col)=SparseNet(row, col)+ MaxFlow(
                            Matrix2Vector(row, col, flows, theta, func, X));
161                     end
                end
163             end
        end
165     end
167
168     SparseNet = SparseNet+SparseNet';
    edgeflows = SparseNet;
169     SparseNet = SparseNet>1e-5;
171
172     SelfFlows = [];
    for i=1:rows,
173         SelfFlows = [SelfFlows, sum(MaxFlow(find(selfflows==i)))]];
    end
175 end

```



---

code/NodeArcLPa.m

## A.2 Embedding Edge LP code

```

function [ embeddings , weights ] = EmbeddingEdgeLP( CommNet, Capacity ,
    FuncWeights, sources , t, CompTrees, MaxFlow )
2 %EmbeddingEdgeLP Given the Network graph, capacities , set of arrays
    source nodes
    %s, set of terminal nodes t, computation trees , and the linprog solution
    , we generate
4 %the set of embeddings where sum(weight_i*x(B_i))=lambda
    %in each iteration of the while loop, we find an embedding with a
    nonzero
6 %flow and remove the corresponding edge flows to obtain another feasible
    %solution with a reduced rate. We start by finding a mapping of
8 %theta_terminal.
    [rows , cols]=size (CommNet);
10 weights={};
    embeddings={};
12 flows = CommNet;
    for row = 1:rows ,
14         for col = 1:cols ,
            if row==col ,
16                 flows (row , col)=1;
            end
18         end
    end
20
    x=find (flows ) ;
22 NumFunctions = length (t) ;
    lambda = MaxFlow (length (MaxFlow)-NumFunctions+1:length (MaxFlow)) ;
24 totalLambda=0;
    X=zeros (1 , NumFunctions) ;
26 for index = 1:NumFunctions ,
        CompEdges (index) = length (find (CompTrees {index}));
28        X (index)=CompEdges (index)*length (x) ;
    end
30
    for func=1:NumFunctions ,
32        newLambda = 0;
        weights {func} = [];
34        embeddings {func} = {};
        while abs (lambda (func) - newLambda)>1 && newLambda<lambda (func) ,
36            stopflag=0;
                thetaT = length (find (CompTrees {func}));
38            B {thetaT}=t (func) ;
                Embeddingflow = [];
40            trackpath = {};
                for i = thetaT:-1:1 ,
42                    trackpath {i} = [];
                        v = B {i} (length (B {i}));
44                    vtried = [];
                        %goes into a loop , keeps trying the same nodes over and over
                        again .
46                    while (stopflag==0),

```

```

48     zeroflows=0;
         possible = [v, GetNeighbors(v, CommNet, vtried)];
         maxf = 0;
50     u = 0;
         for each=1:length(possible)
52             if maxf < MaxFlow(Matrix2Vector(possible(each),v,
                 flows,i,func,X)),
                 u = possible(each);
54                 maxf = MaxFlow(Matrix2Vector(possible(each),v,
                 flows,i,func,X));
                 end
56     end
         if u == 0,
58             stopflag = 1;
         end
60     if MaxFlow(Matrix2Vector(u,v,flows,i,func,X))<=0,
         index = Matrix2Vector(u,v,flows,i,func,X);
62         MaxFlow(index)=MaxFlow(index)-trackmin;
         stopflag=1;
64     end
         vtried = [vtried, u];
66     if u~v && ~isempty(find(B{i}==u,1)),
         P=B{i}(1:find(B{i}==u));
68         if length(P)~=length(B{i}),
             B{i}=B{i}(length(P)+1:length(B{i}));
70         else
             B{i}=[];
72         end
         temp = [];
74         for nodei=1:length(P)-1,
             if P(nodei)~=P(nodei+1) && CommNet(P(nodei),P(
                 nodei+1))~=0,
76                 temp = [temp, MaxFlow(Matrix2Vector(P(nodei),P
                 (nodei+1),flows,i,func,X))];
             end
78         end
         temp = [temp, MaxFlow(Matrix2Vector(u,v,flows,i,func
             ,X))];
80         y = min(temp);
         %update the flows
82         for nodei=1:length(P)-1
             current = Matrix2Vector(P(nodei),P(nodei+1),
                 flows,i,func,X);
84             MaxFlow(current)=MaxFlow(current)-FuncWeights(
                 func)*y;
         end
86         MaxFlow(Matrix2Vector(u,v,flows,i,func,X))=MaxFlow(
             Matrix2Vector(u,v,flows,i,func,X))-FuncWeights(
             func)*y;
         else
88         Embeddingflow = [MaxFlow(Matrix2Vector(u,v,flows,i,
             func,X)), Embeddingflow];
         trackpath{i} = [u,v; trackpath{i}];
90         trackmin = min(Embeddingflow);

```

```

92         end
93         if v~=u,
94             B{i}=[u,B{i}];
95             v=u;
96         else
97             for eta = GetPredecessors(i, CompTrees{func}),
98                 B{eta} = u;
99             end
100            break;
101        end
102    end
103
104    if stopflag==0,
105        xB= min(Embeddingflow);
106        FuncWeights(func)*xB
107        newLambda = newLambda + FuncWeights(func)*xB
108        %Remove x(B) amount of flow from all edges in B
109        Path=[];
110        for theta=1:thetaT,
111            for i=1:length(trackpath{theta}(:,1)),
112                index=Matrix2Vector(trackpath{theta}(i,1),trackpath{
113                    theta}(i,2),flows, theta, func, X);
114                MaxFlow(index)=MaxFlow(index)-FuncWeights(func)*xB;
115            end
116        end
117    else
118        xB=0;
119    end
120    check=0;
121    for e=1:length(embeddings{func}),
122        check=0;
123        for b = 1:thetaT,
124            if length(embeddings{func}{e}{b})==length(B{b}),
125                if embeddings{func}{e}{b}==B{b},
126                    check= check+1;
127                end
128            end
129        end
130        if check==thetaT,
131            break;
132        end
133    end
134    if check==thetaT,
135        weights{func}(e) = weights{func}(e)+FuncWeights(func)*xB;
136    else
137        embeddings{func}{length(embeddings{func})+1} = B;
138        weights{func} = [weights{func},FuncWeights(func)*xB];
139    end
140 end

```

code/EmbeddingEdgeLP1.m

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