Parallel breadth First Search for Scale-Free Social Network Graphs

James Ihrig
University of Central Florida

Josh Eberst
University of Central Florida

Johnathan Pecoraro
University of Central Florida

Abstract—**Graphical representations are one of the most commonly used abstractions to represent large networks. Social media networks like Facebook and Twitter contain massive scale free networks with billions of nodes. The scale free nature of the connections between nodes presents several challenges to the performance of traditional searching algorithms. In this paper, we present two implementations of Parallel Breadth First Search algorithms on scale free networks without the use of specialized hardware.**

I. INTRODUCTION

Graph abstractions are commonly used in large scale network analysis [4]. With the proliferation of "Big Data" applications, parallelized versions of common graph analysis algorithms have a very high demand.

Many graph analysis problems make use of Breadth First Search (BFS), as the efficiency of the algorithm scales linearly with the number of nodes and edges.

Parallelizing Breadth First Search algorithms offers an excellent opportunity to increase the efficiency of this common graph analysis technique. However, scale free networks such as those created by social media connections are challenging to parallelize efficiently as the high distribution of edges to a few nodes is difficult to balance in memory [1]. Our research is targeted at the implementation of two

Parallel Breadth First Search algorithms without the use of specialized hardware on scale free networks. The first algorithm Parallel Breadth First Search with Partitioning (PPBFS) was originally designed to improve BFS performance in distributed memory architectures used by many modern supercomputers. Our implementation will specifically apply this solution to scale free networks, which was left as future work by the original authors. However, as we are not implementing specialized hardware, our research will be performed without the use of a distributed memory architecture.

The second Parallel Breadth First Search algorithm (MTBFS) we will implement is designed to support fine-grained, low overhead synchronization in a massively multithreaded system [1]. This algorithm will be used for gauging the performance of PPBFS, as MTBFS was intended to support scale free graphs. Since the focus of MTBFS is on increased parallelism through multi threading, the differences between the Cray MTA-2 system MTBFS was designed for and a standard multi-core CPU should be less drastic than the shift from the distributed memory architecture used by PPBFS. Memory access by BFS and other graph algorithms

is typically fine-grained and irregular. This leads to poor cache performance, especially in parallel versions of BFS as parallelization relies heavily on the cache performance. Some performance improvements can be made, but they can not be sufficiently generalized since cache performance depends largely on the structure of the graph [4].

We will compare both of these algorithms with the sequential version of BFS across two datasets. One dataset will be a sufficiently large import of a Facebook social network, and the other a generic scale free network generator GDBench [2].

II. RELATED WORK

A Scalable Distributed Parallel Breadth-First Search Algorithm on BlueGene/L

This paper talks about searching networks that are too large to fit into memory of a single machine. To handle this, divides the graph into partition where each node processes a set of vertices assigned to it. If it finds a vertex that does not belong to it, the node that owns it is notified. One major drawback this approach seems to have is that it waits for all processes to reach the same level before moving on. While this is important to truly follow a breadth first pattern, it may not be necessary to limit it in this way.

Designing Multithreaded Algorithms for Breath_First search and st-connectivity on the Cray MTA-2

The Multithreaded BFS algorithm implemented by Bader and Madduri will serve as our baseline parallel BFS algorithm. Although initially implemented on a massively multithreaded shared memory system without a data cache, we believe the fine-grained parallelization of this algorithm will still perform efficiently in a single CPU multi-core system. We have chosen this algorithm to serve as baseline parallelized implementation because it was specifically designed for the traversal of large scale-free graphs similar to those used in our benchmarks.

Parallel Breadth-First Search on Distributed Memory Systems

The partitioning Parallel Breadth-First Search algorithm for distributed memory systems implemented by Buluc and Madduri extends the research performed by Yoo on the Scalable Distributed Parallel Breadth-First Search, and highlights inefficiencies in the supporting structure of the other research. This research provided prior work summaries and criticism on the two algorithms we will adopt, but implementing the improvements made here is outside the scope of our research.

III. Technical Approach

Our projects technical approach will focus on comparing three versions of the BFS algorithm. The original implementations of our two parallel BFS algorithms were intended for use on super-computing hardware. The reason we have implemented two versions of parallel BFS algorithms is to in the performance of the algorithms when they are abstracted away from their specialized hardware. Accordingly we believe that the PPBFS algorithm's performance will be the most dependent on the hardware, and the MTBFS algorithm will be more hardware agnostic. These parallel algorithms will both have their performance compared with a traditional sequential implementation of BFS.

The performance comparison of the algorithms will be determined by the amount of time each algorithm takes to fully traverse every node in two scale-free graphing benchmarks. We have implemented a third party application, NodeXL, to generate large scale-free graphs within our application. The NodeXL program is an open source application written in the C# language. In general, it is used as a network visualization tool which gives users the ability to view a graph's connectivity on a graphical interface. Additionally, there is a social network plugin for the NodeXL application which provides the capability of importing a social network graph from Facebook. We have utilized the class libraries for NodeXL and the associated social network plugin within our application to assist with the generation of large graphs.

Our first benchmark will be based on a graph of a sufficiently large social network imported through the use of the Facebook social network plugin for NodeXL. Our second benchmark is a generic scale-free graph generated by the social network data generator GDBench. The GDBench tool is capable of generating scale-free graphs with millions of nodes, and saving it in a common GraphML format. The GraphML file generated by the benchmark will be imported using NodeXL into a graph consumable by the algorithms. These social networking benchmarks should both be sufficiently large to incur the cache performance challenges associated with many parallel BFS solutions.

IV. Technical Details

Our research is based on a distributed parallel approach to BFS, originally implemented on the IBM BlueGene/L supercomputer. We hope to attain performance increases using the same methods on a single CPU using multiple cores. Additionally, we will attempt to find optimizations to the algorithm such as allowing the processor to continue without waiting for send/receive messages before continuing the search for a given level/depth.

Parallel With Partitioning BFS (PPBFS)

The original algorithm used as a reference for our project was implemented as a distributed BFS on the BlueGene/L architecture. The algorithm takes advantage of the architecture's structure to develop efficient inter-processor communication, which is generally the bottleneck in distributed systems. Our

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Algorithm 1 PPBFS from [4]

1) Initialize $L_{v_s}(v) = \begin{cases} 0, & v = v_s, \text{where } v_s \text{ is a source} \\ \infty & \text{otherwise} \end{cases}$
2) **for** $l = 0$ to ∞ **do**
3) $F \leftarrow \{v \mid L_{v_s}(v) = l\}$, the set of local vertices with level l
4)   **if** $F = \emptyset$ for all processors **then**
5)   Terminate main loop
6)   **end if**
7) $N \leftarrow \{$neighbors of vertices in F (not necessarily local)$\}$
8)   **for all** processors q **do**
9) $N_q \leftarrow \{$vertices in N owned by processor $q\}$
10)   **Send** N_q to processor q
11)   **Receive** $\bar{N}_q$ from processor q
12)   **end for**
13) $\bar{N} \leftarrow \bigcup_q \bar{N}_q$ $\left(\text{The } \bar{N}_q \text{ may overlap}\right)$
14)   **for** $v \in \bar{N}$ and $L_{v_s}(v) = \infty$ **do**
15) $L_{v_s}(v) \leftarrow l + 1$
16)   **end for**
17) **end for**

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implementation uses the techniques described in the paper to develop our parallel version of the algorithm.

The implementation is slightly different than the general approach to performing a BFS. It starts by partitioning the graph equally between the available threads. Each thread will has it's own source node which it uses to start the BFS computation. While a thread is executing the breadth first search, it may come across nodes that it does not own. To solve this problem, a list of neighboring nodes is created. Each neighboring node that is not owned by the current thread is sent to the thread who owns it, which instructs the owning thread that the node has been visited by another thread and it needs to be processed accordingly. Likewise, each thread A also receives nodes that were visited by other threads which belong to A. Each vertex is given a level property. The level denotes the distance the vertex is from the source. During the breadth first search the level property is used to determine which vertices to expand upon. For example, if we are currently examining level one and we see a vertex with level three, we will not expand upon this vertex.

The BFS algorithm takes the partitioned graph as input where each node is initialized to infinity except for the source node, which is initialized to zero. It starts by looping through each level until there are no nodes at the current level (the source node is at level zero). For each iteration, a set of local nodes F is obtained that have level L. If there are no nodes in the set F, the algorithm terminates. Otherwise, a set is created of neighboring nodes N for each node in F. The nodes contained in N are not necessary all owned by the current thread. Then, in parallel, we loop through node in N and mark them as visited in an object shared by all threads. Finally, each remaining unvisited node is marked with the appropriate level index and the loop continues onto the next iteration. In detail, the algorithm works as follows: Loop from level zero to infinity Get set F of nodes owned by current thread with level L End loop if F is empty Get set N of neighbors of verticies in F Loop through all nodes in N and mark the

visited flag Set the level of each node Since the original algorithm was created for a distributed system, it utilized non-traditional methods for conducting the BFS. We expect to see some interesting performance gains from implementing this algorithm in a parallel fashion.

Multi-threaded BFS (MTBFS)

The MTBFS implementation used for reference in this paper was implemented on a Cray MTA-2 multithreaded architecture (Reference paper 2). Similar to our approach, the algorithm was tested against scale-free graphs. When tested with a graph containing 400 million nodes, a 40 processor system showed the multi-threaded algorithm to have a system speedup of about 30 time over the sequential implementation.

The algorithm takes in a graph with source node s, and returns a shortest-path array d such that d[v] contains the length of the shortest path from source node s to destination node v, where v is a node in the graph. The algorithm uses the standard approach for implementing a sequential BFS where you start by adding the source node to a queue, then continue looping until the queue is empty. Each iteration dequeues the first node, finds it's neighboring nodes, and adds all unvisited neighbors to the queue. The main difference is that looping through the queue and through the neighboring nodes is done in parallel. In detail, the algorithm works as follows:

- Initialize all nodes in the graph to negative 1
- Set source node to zero, clear the queue, and enqueue source node
- In parallel, loop through all nodes in the queue
  - Dequeue head node u
  - In parallel, loop through all neighbors v of u
  - Enqueue all unvisited neighbors, and set the return array's length
    * d[v] = d[u] + 1

This algorithm takes full advantage of the Cray MTA-2 system's architecture by using its fine-grained parallelism and zero-overhead synchronization while looping through queue and looping over each node's set of neighbors. This multi-threaded BFS technique along with the hardware architecture design offer a considerable speedup advantage over any sequential implementation of BFS on graphs of similar size.

V. CHALLENGES

We have noticed that the PPBFS algorithm [3] waits for all other threads to complete a search level before moving on to the next level. This places some sequential constraint on the algorithm and can potentially limit performance.

VI. REMAINING WORK

Thus far we have successfully implemented and modified the NodeXL class libraries to support additional properties for BFS, and implemented a sequential BFS algorithm utilizing a graph imported from Facebook with NodeXL. We have also implemented a sequential version of the PPBFS Algorithm.

  1) Finish the parallelization modifications to the PPBFS Algorithm

  2) Implement the MTBFS Algorithm
  3) Execute the algorithms

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