Parallelisation of Graphing Algorithms in Julia

Group 4

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Agenda

- Introduction to Julia
- Graphing algorithms
- Parallelization of Graphing algorithms
- Our implementation/benchmarking

What is Julia



Julia is

- Open source
- High-level
- High-performance
- Dynamic programming language

Designed for numerical computing

"Looks like python, feels like lisp, runs like C"

Syntax

Python

def sum(a):

$$s = 0.0$$

for x in a:

$$s += x$$

return s

<u>Julia</u>

function sum(a)

$$s = 0.0$$

for x in a

$$s += x$$

end

return s

end

- 🔹 1-based indexing 😠
- Macro support
- Homoiconicity
- Structs

Generality

Julia has metaprogramming support similar to Lisp

```
macro name (expr, expr)
```

...modify evaluation...

End

@name (expr, expr)

Compilation in Julia

Julia "runs like c" mostly due to it's compilation

Julia uses a JIT (just-in-time) compiler based on LLVM to generate native machine code

optimize unnecessary static branches out at runtime

Julia Tasks

Used to execute of multiple functions co-operativly

Declare a channel

```
c1 = Channel(32)
c2 = Channel(32)
```

!take and !fetch data from and !put data into channels

```
data = take!(c1)
put!(c2, result)
```

close () channels when done with them

```
close(c1);
close(c2);
```

Native Threads

Currently Experimental (only supports for loops)

@distributed

@Threads

```
Threads.@threads for var = range body

End
```

(does not support optional reduction parameter)

Processes

Future

- Remotecall the function
- Fetch() the result

RemoteChannel

- Create remote channel
- !put() !take() to/from remote channel

```
./julia -p 2
Create workers
```

RemoteChannel(pid::Integer=myid())

Create remote channel

Parallel computing constructs in Julia

Julia Tasks

- Useful for concurrency
- "Appear" as multiple threads
- In Julia all executed on one system thread
- Not hugely useful for parallel speedup
- Uses Tasks (Coroutines) with Channels to communicate

Native Threads

- Somewhat minimal support in Julia (still experimental)
- @thread and @distributed
- Fork-join approach
- OpenMP style loop parallelisation

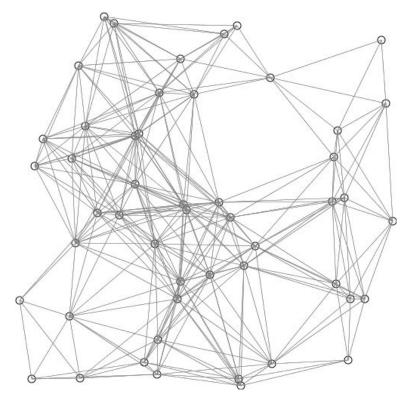
Processes

- Expensive and heavyweight
- Message passing
- Coarse grain granularity

Graph algorithms

Algorithms that solve a variety of problems that exist within graph theory. Categories include:

- Minimal Spanning Tree
- Graph Traversal/Searching
- Shortest Path
- Maximal Independent Subset
- Transitive Closure



https://en.wikipedia.org/wiki/Dijkstra%27s_algorithm#/media/File:DijkstraDemo.gif

Minimum Spanning Tree

A subset of the edges

- In a graph that is connected, edge-weighted and undirected.
- No cycles
- Minimal possible total edge weight

Inherently Sequential

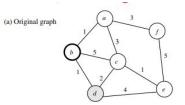
- Most MST algorithms "grow" the spanning tree
- Greedy algorithms

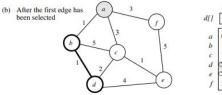
Somewhat parallelizable

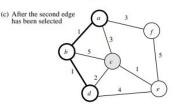
- The selection process of candidate nodes can be parallelized
- Distance matrix can be partitioned between processors

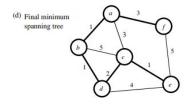
Overhead considerations

- Alternative methods of speedup may be more optimal, (Binary heap adjacency list etc.)



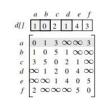












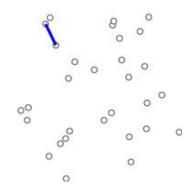


Prims Algorithm

Prim's Algorithm

A greedy algorithm for finding a minimum spanning tree in an undirected graph.

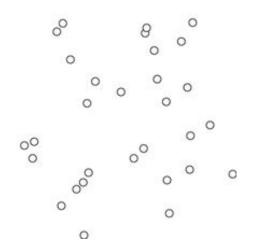
- Choose a vertex and build tree around it
- Building of tree is mostly sequential
- We can parallelise the process of selecting the shortest edge.
- Time complexity depends on data structure used



Kruskal's Algorithm

Another greedy algorithm to find a minimum spanning tree.

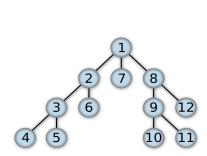
- Similar to Prim's but selects an edge instead of a vertex
- Better for sparse graphs
- Next selected edge not necessarily connected to previous edges.
- Also inherently a sequential algorithm



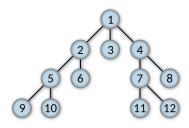
Traversing / Path finding

Search throughout a graph, starting from a source vertex. Includes finding shortest paths.

- Each node visited exactly once
- Examples include:
 - o **A***
 - o Dijkstra's
 - o BFS
 - o DFS



Order nodes are visited in DFS



Finding a shortest path with Dijkstra's

Order nodes are visited in BFS

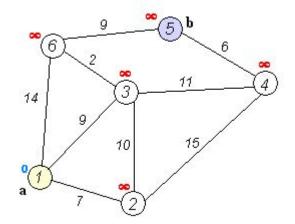
Finding a shortest path with A*

tps://upload.wikimedia.org/wikipedia/commons/thumb/3/33/Breadth-first-tree.svg/300px-Breadth-first-tree.svg, Ibs://upload.wikimedia.org/wikipedia/commons/thumb/1/1/Depth-first-tree.svg/300px-Depth-first-tree.svg ng Ibs://upload.wikimedia.org/wikipedia/commons/2/23/Diikstras_progress_animation.gif

Dijkstra's Algorithm

One of the most well known graph algorithms.

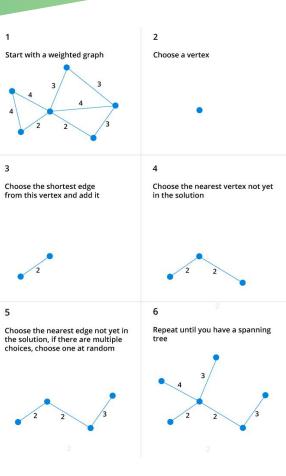
- Published in 1959
- Finds the shortest path to nodes in a graph from a given source node, producing a shortest path tree.
- Parallelizable in a way similar to Prims and Kruskal:
 - After a node is visited the adjacency matrix can be partitioned amongst the available processors n.
 - Each processor seeks to find a local minimum next node
 - Reduction is then done to find global minimum for next iteration
- Time complexity of O(E log V) with help of a binary heap



Prims in parallel

While the process of building the tree is sequential, we can parallelize the process of choosing the closest node.

- The distance array is partitioned amongst the available threads in a distributed for loop
- Each processor finds minimum distance in their array partition
- A reduction then occurs to find the globally closest node as the next node for the tree
- Updating the distance array after a node is also be made parallel



Dijkstra's all-sources in parallel

Source Partitioning

- Use p processors, distribute the vertices between the processors.
- Each processor sequentially executes a single source Dijkstra algorithm on its allocated vertices
- Can only use as many processors as vertices in the graph
- Lower overhead

Source Parallel

- Used if we have more processors than vertices
- P processors split into n nodes
- n/p processors working on each node
- Greater exploitation of parallelism

Initial Benchmarks

Run Time

- Is a parallel solution faster than a sequential one?
- What types of graphs are better suited for a parallel approach?

Efficiency

 How effectively does our solution use additional resources?

Dataset

Randomly generated graphs, both sparse and dense

Machine 1 (Toshiba Portégé A600):

- Intel Core 2 Duo SU9300 / 1.2 GHz
- 2 Cores
- 2GB RAM
- Lubuntu 18.10 64bit

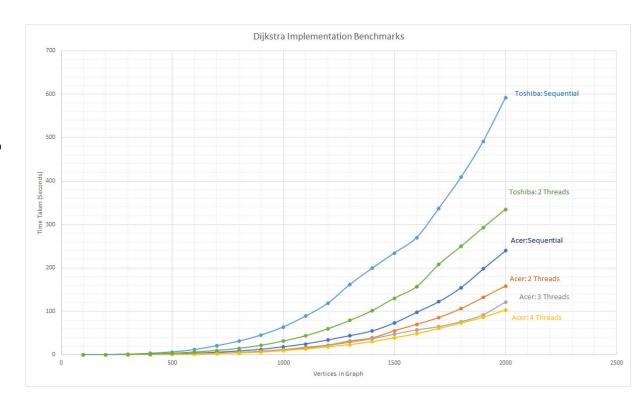
Machine 2 (Acer Aspire s7)

- Intel Core i7-5500U / 2.4Ghz up to 3Ghz
- 2 Cores with HyperThreading (4 Threads)
- 8GB RAM
- Ubuntu 18.04 64bit

Dijkstra all-sources

Source Partitioned all-sources Dijkstra implementation

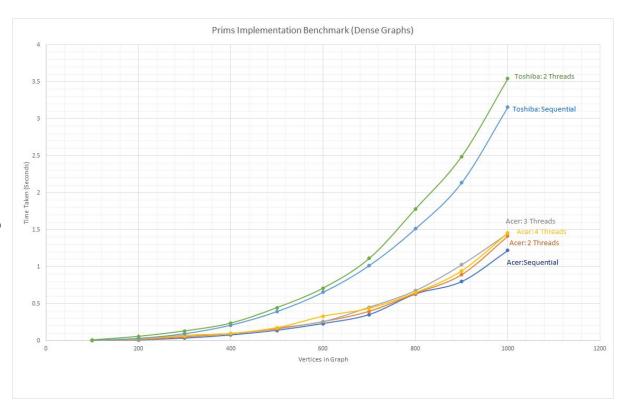
- Starting vertices are distributed between available threads. Used @threads macro
- Minimal inter-process communication
- Decent speedup across the board, (relatively efficient)
- Next step: source parallel implementation



Prims

Parallelisation becomes much more viable when the problem each thread has to solve is bigger

- Sparse graphs tended to suffer very heavily from overhead
- With more neighbouring nodes to process, the proportion of useful work to overhead increased



Next Steps

- More extensive benchmarking suite
 - SNAP Dataset
 - Benchmark.jl
- Algorithm improvements
 - Further Speedups
 - Data structure improvements
 - More complex algorithms (Maximal Independent set, Luby's)

