# Using MVAPICH2-GDR for multi-GPU data parallel graph analytics

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# Overview

#### 1. Review of previous results

• Slides from MUG 2014

#### 2. Recent issues

- Current CUDA driver/runtime reduced performance
- Single-node system performs worse than cluster

#### 3. Future plans

- GPU accelerated graph database
- Translation from Scala DSL



# Slides from MUG 2014



#### Parallel Breadth First Search on GPU Clusters using MPI and GPUDirect

Speaker: Harish Kumar Dasari, Scientific Computing and Imaging Institute, University of Utah

Advisor: Dr. Martin Berzins, SCI, University of Utah

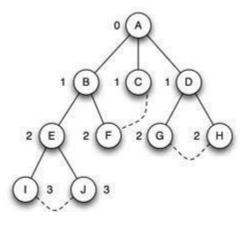
In collaboration with Dr. Zhisong Fu, Bryan Thompson, Systap, LLC.

http://sourceforge.net/projects/mpgraph/



#### Introduction

- Breadth First Search: It is a graph search algorithm that begins at the root vertex and explores all the connected vertices, traversing all vertices of a particular level before traversing the vertices of the next level
- At the end of the BFS we can find out the level of a vertex if it is connected to the root element and also its predecessor
- Useful in social media, logistics and supply chains, e-commerce, counter-terrorism, fraud detection etc.

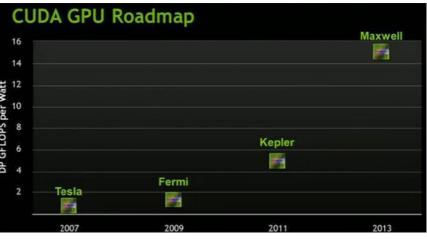




#### Introduction

- Why BFS?
  - Least work/byte of the graph algorithms
  - Building blocks for many other graph problems
- Why GPUs?
  - High Performance: NVIDIA K40 peak performance: 1.43 Tflops
  - High Energy Efficiency
  - Central for next generation of architectures



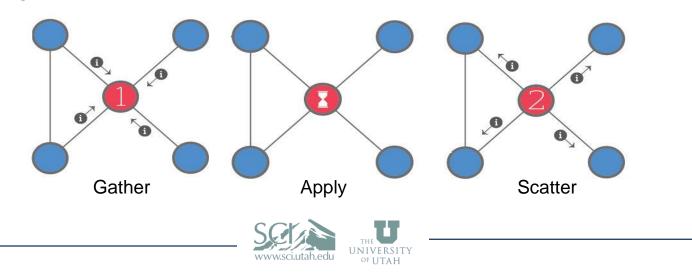






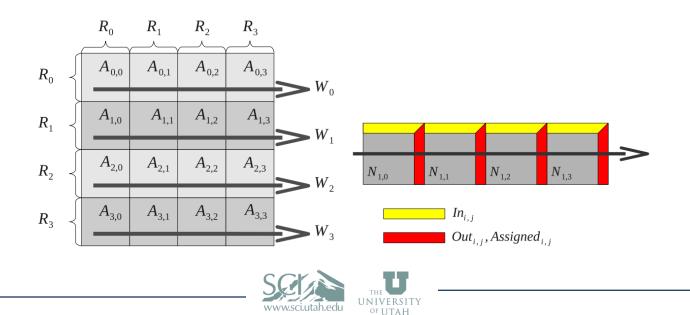
#### **Related Work**

- Scalable GPU Graph Traversal Single node multi-GPU, Merrill, Garland et al.
  - Around 12x speedup over idealized multi-core CPU
  - 3 GTEPS on single node
- MapGraph, Fu, Thompson et al.
  - Generalized for many graph algorithms using Gather Apply Scatter (GAS) abstraction
  - Provides an easy framework for the developer to develop solutions to other graph problems like SSSP(Single Source Shortest Path), PageRank etc.



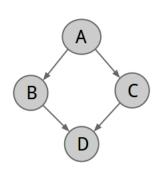
#### **Related Work**

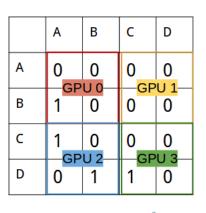
- Breaking the Speed and Scalability barriers for graph exploration on distributed-memory machines by Checconi, Petrini et al from IBM
  - BFS on Bluegene supercomputers, uses CPUs
  - On Graph500 data sets, on the order of 2<sup>40</sup> edges
  - 254 billion edges/sec with 64k cores
  - Uses 2D partitioning and waves for communication

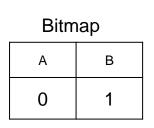


### Partitioning of the Graph

- RMAT graph generated using the Graph500 generator
  - Scale Free
  - Follows power law, at least asymptotically
  - undirected edges are converted to directed edges
- 2-D Partitioning of directed edges with a square layout
- Each subgraph resides in GPU memory
- Bitmaps used to represent the frontiers
  - Bit is set to 1 to represent active vertex





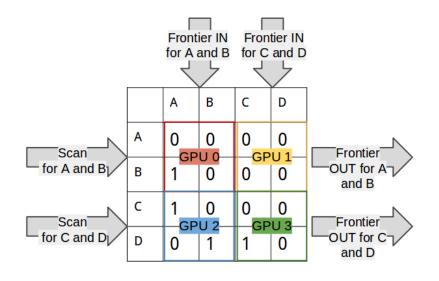


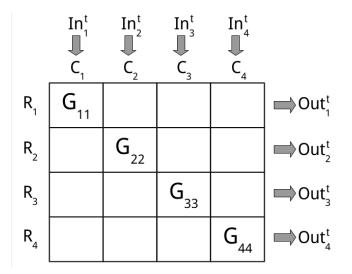




### The Algorithm and Communication

- Each GPU G<sub>ij</sub> takes in its input frontier bitmap In<sup>t</sup><sub>i</sub> and perform BFS on its subgraph to produce Out<sup>t</sup><sub>ij</sub>
- Parallel Scan for bitmaps along the row R<sub>i</sub> to produce prefix sum *Prefix<sub>ii</sub>* in Bitwise-OR



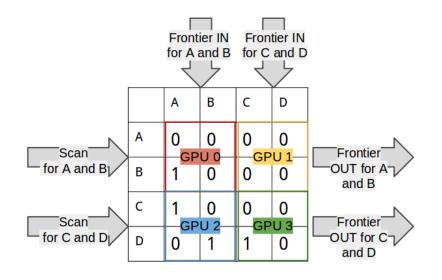


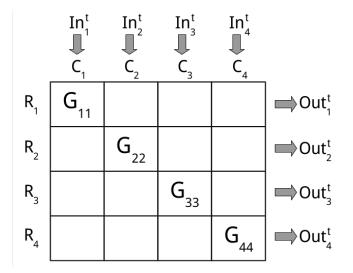


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### The Algorithm and Communication

- The Prefix is used to determine the vertices the GPU is assigned for predecessor updates
- $Out_i^t$  is broadcast across row  $R_i$  and also as  $In_i^{t+1}$  across column  $C_i$



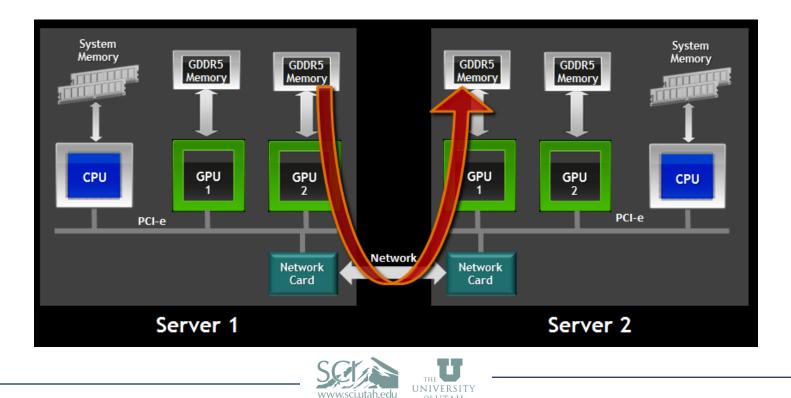






#### **Experimental Setup**

- 32 nodes and 64 NVIDIA K20c GPUs with 5GB DDR5 memory
- Two Mellanox InfiniBand SX6025 cards per node
- CUDA 5.5 used for these results
- Used GPUDirect support in MVAPICH2-GDR to avoid explicit copy of messages to host memory



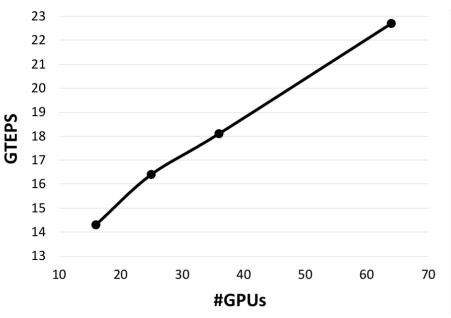
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### **Results - Strong Scaling**

- The scale of the problem remains the same as we increase the computational resources (GPUs)
- GTEPS= Giga(Billion) Traversed Edges Per Second = 10<sup>9</sup> edges per second

GPUs	Scale	Time	GTEPS
16	25	0.075	2.5
25	25	0.066	6.3
36	25	0.059	15.0
64	25	0.047	29.1

Number of Vertices in graph =  $2^{SCALE}$ Number of Directed Edges in graph =  $32^{SCALE}$ 





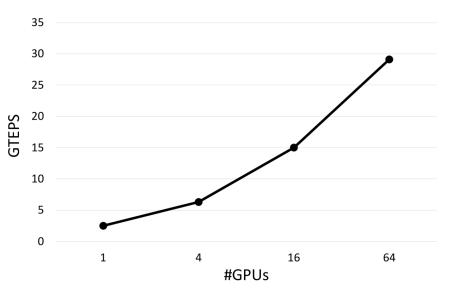
### **Results - Weak Scaling**

- Problem size grows proportional to the growth in computational resources (GPUs)
- Each GPU has same amount of work?

GPUs	Scale	Time	GTEPS
1	21	0.0254	14.3
4	23	0.0429	16.4
16	25	0.0715	18.1
64	27	0.1478	22.7

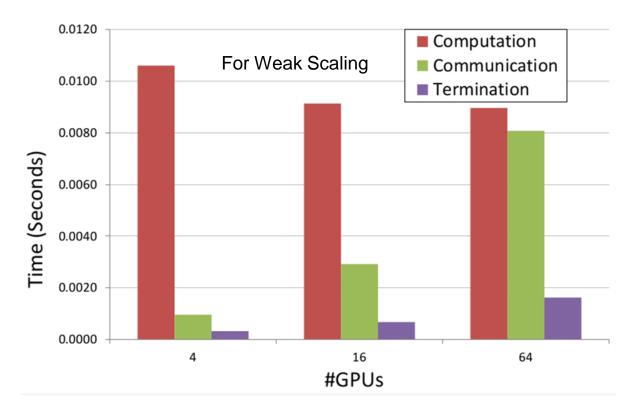
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#### **Communication vs Computation**

- Even if the work per GPU remains the same, the communication costs grow
- Impacts weak Scalability

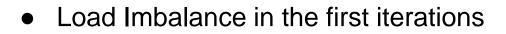


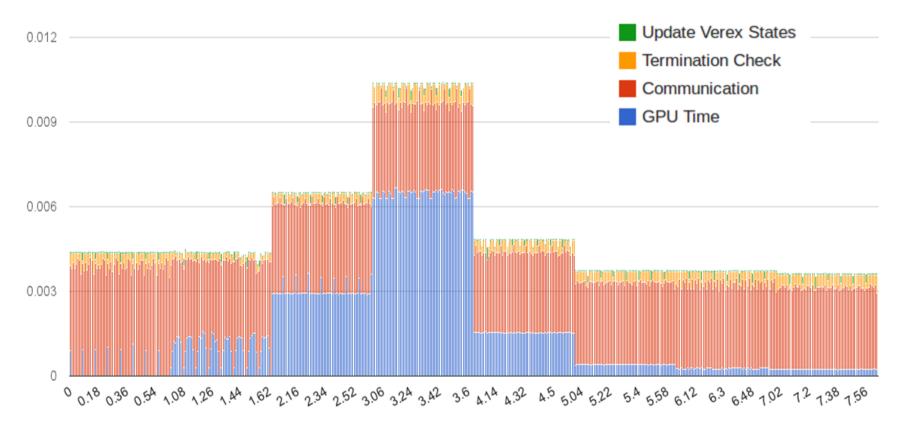


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### **Breakdown of Timings**

Near constant communication times across iterations





iteration . MPI rank





### **Recent Issues**

Reduced performance with version updates Nvidia PSG cluster: 16 K40s across 4 nodes

- Cuda 5.5
- MVAPICH2-GDR 2.0b
- 18.74 GTeps

- Cuda 7.0
- MVAPICH2-GDR 2.1 rc2
- 9.13 GTeps

Only 48% of previous performance



### **Recent Issues**

#### **Disappointing Single Node Performance**

16 MPI processes Scale 25 graph

Cirrascale	<ul> <li>Nvidia PSG Cluster</li> </ul>
• 8x K80 (16 GPU)	• 4 Nodes, 16 K40
Full PCIe-3 16x	• Cuda 7.0
• Cuda 7.0	• MVAPICH2-GDR 2.1 rc2
• MVAPICH2-GDR 2.1 rc2	• 9.13 GTeps
• 2.20 GTeps	



## Future Plans

#### GPU accelerated graph database

- BlazeGraph graph database (Java)
- Accelerate SPARQL queries with GPU

#### Translation from Scala DSL

- DSL defined operators
- Graph algorithms written in Scala
- Scala translated to native GPU code
- High performance without GPU experts



### Questions

