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Model Evaluation

A Parallel Model for Heterogeneous Cluster

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Motivation

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Clusters are becoming more heterogeneous

Distinct processors, accelerators, and network connections

To explore all the resources available in such a heterogeneous platform, a data-parallel application must divide its data across multiple devices

Distinct processing power of devices and the distinct latencies of the networks

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Which configuration leads to the best speedup?

Contribution

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- Present a parallel model that estimates the execution time of applications running on heterogeneous clusters
 - Extends some characteristics of the LogP model
 - Considers that processing units may have distinct computational power as well as they are interconnected by connections with distinct latencies
- The idea is to use the results of this estimation, in future works, to predict the best data division to be used in a heterogeneous cluster
 - Taking into account not only the processing power of each processor and accelerator, but also the communication and synchronization costs.

LogP Model

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LogP Model

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- Measures the effects of latency, occupancy and bandwidth on distributed memory multiprocessors
- Main parameters used in the LogP model
 - L represents an upper bound on the communication latency due to the use of point-to-point messages
 - o represents the overhead
 - **g** represents the minimum time interval between consecutive message transmissions/receptions by a processor (gap)
 - \blacksquare The reciprocal of the ${\bf g}$ parameter represents the communication bandwidth

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P represents the number of processor/memory modules

Related works

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- Heterogeneous processors interconnected by an Ethernet-based network
 - Homogeneous network

HLoGP model

- Takes into account the heterogenity of both computation and communication resources
- Large number of parameters is an issue
- This work proposes a simpler model that predicts the execution time of parallel applications
 - Regardless of the computational environment used, homogeneous or heterogeneous one.

The new model

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Model Evaluation

- Deal with modern heterogeneous environments, composed by distinct processors, accelerators and networks
 - **L**_d represents an upper bound on the communication latency of a device d;

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- \mathbf{o}_d represents the overhead in device d
- **g**_d represents the minimum time interval between consecutive message transmissions/receptions by a processor in a device d (gap)
- *R_P* represents the relative computing power of a processing unit
- Parameters and variables are used to describe mathematically the total execution time of an application

The new model

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- How to measure the relative computing power (R_P) ?
 - Running a benchmark on each processing unit to collect a metric, such as the processing units per time step
 - Using the average computation time that a processing unit takes to run some iterations of an application
- How to measure the values of the latency (L_d) and the gap (g_d) ?
 - Network benchmark is used for this purpose
 - Benchmark is executed for each type d of network that is available
 - Collects the values of L_d and g_d for distinct message sizes, ranging from 0 to 4MB

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- How to measure the overhead (**o**_d)?
 - Also measured with a specific benchmark
 - It considers that the overhead varies with the message size
- Use of benchmarks to collect the communication costs, overheads, as well as the relative performance of the processors and accelerators, can be executed only once

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Each time a new hardware or network is included in the system

Model Evaluation

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- Two kernels (EP and FT) and one application (SP) from the NAS benchmark were used in the initial validation of the model
 - Benchmarks were developed to execute in a CPU environment
- HIS (human immune system) simulator was chosen to evaluate the model on a hybrid environment

Uses GPUs and CPUs simultaneously

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- Embarrassingly Parallel kernel solves a typical problem of many Monte Carlo based applications
 - Generate pairs of Gaussian pseudorandom deviates
- Communication occurs only at the end of the computation
 - Collective MPI routine is used to combine the sums generate from all processors

Class C used in the evaluation

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Algorithm 1 EP

- 1: **main**
- 2: ... generate the seed for each process ...
- 3: ... calculate counts and sums in each process ...
- 4: ... Use MPI_Allreduce to send parameter to all processes ...

5: end-main

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- 3-D Fast-Fourier transform kernel
 - Used to numerically solve partial differential equation (PDE)

- All-to-all communication used to exchange the transpose results
- Class B used in the evaluation

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Algorithm 2 FT

1: main

3.

4:

- 2: for t from 1 to number of iterations do
 - ... evolve u0 to u1 (t time steps) in fourier space ...
 - ... calls the fft subroutine ...
- 5: ... transpose operations in each process ...
- 6: ... use MPI_Alltoall to exchange the transpose results ...

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- 7: ... call checksum ...
- 8: end-for
- 9: end-main

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Scalar Penta-Diagonal solver

Solves multiple, independent systems of nondiagonally-dominant, scalar pentadiagonal equations

- Coarse grained communication
- Class B used in the evaluation

SP

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Model Evaluation

Algorithm 3 SP						
1:	main					
2:	for t from 1 to number of iterations do					
3:	performs the block-diagonal matrix vector multiplicator					
4:	use MPI_Isend to send the buffer					
5:	use MPI_Ireceive to receive the buffer					
6:	performs aproximate factorization in the x-plane					
7:	use MPI_Isend to send the buffer					
8:	use MPI_Ireceive to receive the buffer					
9:	performs aproximate factorization in the y-plane					
10:	use MPI_Isend to send the buffer					
11:	use MPI_Ireceive to receive the buffer					
12:	performs aproximate factorization in the z-plane					
13:	use MPI_Isend to send the buffer					
14:	use MPI_Ireceive to receive the buffer					
15:	add the u vector					
16:	end-for					

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- Three dimensional simulator of the Human Immune System
 - Set of eight Partial Differential Equations (PDEs) used to describe how some cells and molecules involved in the innate immune response react to a pathogen.

- 200 × 200 × 200
- Border exchange occurs at the end of each time iteration

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Algorithm 4 HIS

1: main

2: ... define the mesh slice to be computed by each GPU/CPU ...

3: ... initialize submeshes according to their initial conditions ...

4: for t from 0 to final time do

5: ... call the functions/*kernels* in order to compute the PDEs ...

6: ... use MPI_Isend and MPI_Receive to exchange boundaries between distinct machines ...

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7: ... synchronize all machines ...

8: end-for

9: end-main

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Conclusion

• The EP benchmark is modeled using the following equation:

$$T_{total} = \frac{size}{R_p} + I \times N_{op} \times \log_2 P \times (L_d + \frac{M}{B_d} + o_d), \tag{1}$$

- size is the size of the problem
- R_p is the relative computing power
- I is the number of iterations
- N_{op} is the number of communication operations per iteration
- L_d is the latency
- o_d is the overhead
- B_d represents the bandwidth
- $\blacksquare\ P$ is the number of processors used in the experiments and
- $\blacksquare~M$ is the message size

FT and SP

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• The FT benchmark is modeled using the following equation:

$$T_{total} = I \times (R_p + N_{op} \times (P - 1) \times (L_d + \frac{M}{B_d} + o_d))$$
(2)

• The SP benchmark is modeled using the following equation:

$$T_{total} = I \times (R_p + N_{op} \times (L_d + \frac{M}{B_d} + o_d))$$
(3)

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• The HIS benchmark is modeled using the following equation:

$$T_{total} = I \times (R_p + T_{ij}), \tag{4}$$

where

$$T_{ij} = (L_d + \frac{M}{B_d} + o_d) \tag{5}$$

Experimental environment

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Sixteen machines

- Two distinct CPUs
 - Intel E5620 dual quad-core processors
 - AMD 6272 dual sixteen-core processors

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- One process per machine
- Three distinct GPUs
 - Tesla C1060
 - Tesla M2050
 - Tesla M2075
- Two distinct networks
 - Gigabit ethernet
 - InfiniBand

Results

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Table: Results for the EP, FT and HIS using 2 AMD processors. All times are in seconds.

		Ethernet		Infiniband			
	Real	Estimated	Error	Real	Estimated	Error	
EP	295.6	295.8	0.1%	297.2	295.8	0.5%	
FΤ	95.0	96.3	1.5%	66.1	69.4	5.0%	
HIS	213.4	219.1	2.7%	102.7	109	6.1%	

SP code requires a square number of processors

Results

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Table: Results for the EP, FT and SP kernel using both Intel and AMD processors (half of each), Ethernet network. All times in seconds.

	4 Nodes			8(9) Nodes*			16 Nodes		
	Real	Estimated	Error	Real	Estimated	Error	Real	Estimated	Error
EP	118.0	110.5	6.4%	52.0	55.2	6.3%	28.6	28.6	0.0%
FT	71.4	72.0	0.9%	67.0	68.1	1.8%	65.8	64.1	2.7%
SP	442.3	445.7	0.6%	265.9	267.7	1.0%	343.7	345.4	0.5%

*For SP, we used 9 nodes (4 AMDs and 5 Intels) since the code requires a square number of processors

Results

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Table: Results for HIS using both GPUs and CPUs. All times in seconds.

	Real	Estimated	Error
1	47.2	42.5	10.0%
2	59.8	54.2	9.2%
3	107.8	95.0	12.0%

- Configuration number 1: 2 CPUs (1 AMD and 1 Intel) and 2 GPUs (M2075 and C1060)
- Configuration number 2: 4 CPUs (2 AMDs and 2 Intels) and 4 GPUs (2 M2075 and 2 C1060)
- Configuration number 3: 7 CPUs (5 AMDs and 2 Intels) and 7 GPUs (3 M2075, 2 M2050 and 2 C1060)

Conclusion

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- New model that generalizes the LogP model in order to deal with heterogeneous parallel environments
- Model can predict the total computation time of applications with distinct characteristics, running on distinct devices and interconnected by different network types
- Errors found during the estimation of the total execution time were below 6.4% in all experiments
 - Except for the HIS simulator, where the error was about 12% when distinct CPUs and GPUs were used in the simulation

Future works

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- Better understand the causes of the higher error found in HIS
 - Evaluate the model with more applications
 - Use the model to choose the data partition and work assignment that minimizes the execution time of an application

Thank you!

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