## The SIMGRID Project

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Inria

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Algorille INRIA Simulation for Large-Scale Distributed Computing Research

## The Accuracy vs. Speed tradeoff

Common Belief in 2008: Simulation as a toy methodology



## The Accuracy vs. Speed tradeoff

- ► Common Belief in 2008: Simulation as a toy methodology
- ► Consensus in 2012: SimGrid as a scientific instrument (w/ Grid'5000)



#### How did we turn **Simulation** into a **Reliable and Versatile Scientific Instrument** for Distributed Computing Research?

- A Performant et Versatile Simulation Kernel (high-performance simulation for computer science)
- Simulating Real MPI Applications (beyond prototypes)
- Toward a Coherent Workbench for Distributed Applications (when simulation is not enough)

### Layered Infrastructure for a Versatile Tool

### SimGrid: strictly layered and built bottom-up

#### SimGrid Functional Organization

- Models: Actions get mapped onto resources Resource sharing and termination dates
- Activities: Processes interact and synchronize
- User interfaces: User-friendly syntaxic sugar

#### SimGrid user APIs

- SimDag: heuristics as DAG of (parallel) tasks
- MSG: heuristics as CSP (Java/Lua/Ruby bindings)
- SMPI: simulate MPI codes



### **Models: Resource Sharing between Actions**

#### How to Model the Platform?



- $x_1 \leqslant Power_-CPU_1$  (1a)
- $x_2 + x_3 \leqslant Power_-CPU_2$  (1b)
- $y_1 + y_2 \leqslant Power\_link_1$  (1c)
- $y_1 + y_3 \leqslant Power_link_2$  (1d)

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#### Production-grade Implementation

- ▶ Efficiency: Sparse structure; Cache oblivious; Lazy evaluation
- Realism: Several fairnesses can be expressed this way (or NS3 bindings)

# Putting the Models in Use



Bob			
(some code)			
Receive from Alice			
(other code)			
Send "blah" to Alice			



#### SimGrid Internal Main Loop

- 1. Run every ready user process in row
  - Each wants to consume resources
  - Assign actions on resources
- 2. Compute share for actions
- 3. Get earliest finishing action; update clock

#### Production-grade Implementation

- Scalability: Contextes instead of threads; Hierarchical networks
- Speed: Context switches in assembly; Futexes; Original parallelisation schema
- Other: Resource availability changes and failures; Dynamic Formal Verification



How big and how fast? (1/3 - Grid)

#### Size of platform description files

Community	Scenario	Size
P2P	2,500 peers with Vivaldi coordinates	294KB
VC	5120 volunteers	435KB + 90MB
Grid	Grid5000: 10 sites, 40 clusters, 1500 nodes	22KB
HPC	1 cluster of 262144 nodes	5KB
HPC	Hierarchy of 4096 clusters of 64 nodes	27KB
Cloud	3 small data centers $+$ Vivaldi	10KB

#### Speed of Grid Scenario

A master distributes 500,000 fixed size jobs to 2,000 workers (round robin)

	GridSim	SimGrid
Network model	delay-based model	flow model
Topology	none	Grid5000
Time	1h	14s
Memory	4.4GB	165MB

# How big and how fast? (2/3 - P2P)

- ► Scenario: Initialize Chord, and simulate 1000 seconds of protocol
- Arbitrary Time Limit: 12 hours (kill simulation afterward)



- SIMGRID orders of magnitude more scalable than state-of-the-art P2P simulators
- $\blacktriangleright$  Precise model incurs a  $\approx 20\%$  slowdown, but accuracy is not comparable

How big and how fast? (3/3 - HPC)

#### Simulating a binomial broadcast



Model:

- ► SIMGRID: contention + cabinets hierarchy
- ► LogGOPSIM: simple delay-based model

#### Results:

- SIMGRID is roughly 75% slower
- SIMGRID is about 20% more fat (15GB required for 2<sup>23</sup> processors)

The genericity of SIMGRID data structures comes at the cost of a slight overhead BUT scalability does not necessarily comes at the price of realism

# **Conclusion**

#### SimGrid is ready to ground your Research

- ▶ Versatile: Grid, P2P, HPC, Volunteer Computing, Clouds, ...
- ► Valid: Accuracy limits studied and pushed further for years
- ► Scalable: 3M chord nodes; 1000× faster than other (despite sound models)
- ▶ Usable: Tooling (generators, runner, vizu); Open-souce, Portable, ...



#### But a simulation kernel is not sufficient

- Users need love (Coming: Simulating MPI applications)
- Simulation is no universal solution (Coming: coherent workbench)

### Single online simulation with SMPI

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Simulation for Large-Scale Distributed Computing Research

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  - Validity of these models for MPI applications
- Folding of the parallel program processes onto a single node
  - Serialization of computations
  - Single address space
  - Requires to reduce
    - Memory footprint (scalability)
    - Simulation time (speed)

### **Reworked Network Model**

- Simple Model:  $T(S) = L + \frac{S}{B}$
- ► Improved model:  $T(S) = \alpha \cdot L + \frac{S}{\min(\beta \cdot B, \frac{\gamma}{2 \cdot L})}$ 
  - $\blacktriangleright \ \alpha$  accounts for TCP slow-start
  - $\beta$  accounts for the overhead induced by TCP/IP headers (e.g 92%)
  - $\blacktriangleright~\gamma$  enables the modeling of the TCP window induced behavior
  - $\blacktriangleright$  Model valid for S  $\geq$  100 KiB, does not address a lot of message sizes found in MPI applications
- ▶ Need for a new, accurate network model when S < 100 KiB



Experimental measurement using SKAMPI







- packet size < MTU,
- eager/rendezvous switch limit



# **Collectives and Contention**

Scatter: 16-processes test



- Comparison SMPI/OpenMPI: error 5.3%
- Taking contention into account is important

- Idea: Share arrays between processes
- Implemented as optional macros

```
double* data = (double*)SMPI_SHARED_MALLOC(...);
...
SMPI_SHARED_FREE
(data);
```

# **Reducing the Memory**



- Average reduction by factor of 11.9 (maximum 40.5x)
- Class C can now be simulated

# **Reducing the Simulation Time**

- Idea: Do not execute all the iterations
- Use sampling instead
  - LOCAL: each process executes a specified number of iterations
  - GLOBAL: a specified number of samples is produced by all processors
- Remaining iterations are replaced by average of measured values
- Implemented as optional macros

```
for(i = 0; i < n; i++) SMPI_SAMPLE_LOCAL( 0.75*n , 0.01 ) {
}
...
for(j = 0; j < k; j++) SMPI_SAMPLE_GLOBAL(0.5*k,0.01) {
...
}</pre>
```

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# Wrap-up

SMPI is a functional simulation tool

- Open Source and freely available
- Reproducible simulation of unmodified MPI application
- On a single node
- Main issues addressed:
  - scalability and speed through macros,
  - accuracy through extensions of the network model

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However, microscopic behaviors are difficult to capture, e.g:

- network communication jitters,
- network catastrophes,
- cache effects,
- ▶ ...

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  - ▶ ...

And hence, simulation must be used in conjunction with other experimental approaches: emulation or experimentation in the real environment.

#### Simulator



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- Team Focus on emulation and orchestration of experiments
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#### Key role in both projects

#### Simulator



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#### **Complementary solutions:**

- ③ Work on algorithms
- © More scalable, easier

Towards an unified workbench

- © Work on applications
- © Perceived as more realistic











#### Goal: convergence of methodologies

# **Challenges and opportunities**

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  - Frameworks for data analysis and vizualisation

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 $\neq$  scientific instruments implementing  $\neq$  scientific methodologies Towards an unified workbench

# **Challenges and opportunities**

- Share experimental methods and software
  - Infrastructure for Design of Experiment
  - Frameworks for data analysis and vizualisation
- Design better models and better testbeds using the common expertise e.g. network or power consumption modelling vs instrumentation
- Attack the same goals together, from both sides Reproducibility, trustworthiness, Open Science

We are in a unique position to address those challenges