













High Performance Computing as a Combination of Machines and Methods and Programming



*Soutenance en vue de l'obtention du Diplôme
d'Habilitation à Diriger les Recherches*

Claude Tadonki
Mines-ParisTech

Université Paris-Sud / 16 mai 2013

- ▶ Fundamental Aspects of Algorithms and Complexity
(Formalism, proof, quantification, classification, ...) *"A vouloir toujours aller au fond des choses, on court le risque d'y rester."*
- ▶ Graph Theory and Applications (characterisation, modeling, scheduling)  Claude Berge
- ▶ Discrete Dynamical Systems (Chip Firing Game, Periodicity, Garden of Eden, Invariants, ...)  Charles Leiserson
- ▶ Parallel Scheduling (Theory, SPMD, SIMD, Systolic, ...)  Quinton  Charles Leiserson  Richard Karp
- ▶ Polyhedral Model (Recurrences Equations and Scheduling)
- ▶ Efficient Parallel Programming (Hybrid Parallel Computing, Scalability, Load Balancing, Data Transfers & Exchanges)
- ▶ Automatic Code Generation and Transformations (High-Level Specification \Rightarrow Program; Program + Annotations \Rightarrow Program)
- ▶ Power Aware Computing (Energy Efficient Programming: Models and Methods, Cloud Computing)  Schrijver  Vial  Maculan
- ▶ Applied Mathematics and Operation Research (Mathematical modeling, Optimization, Linear Algebra, Matrix Computation)  Saad  Vial



"A vouloir toujours aller au fond des choses, on court le risque d'y rester."



Claude Berge



Quinton



Charles Leiserson



Richard Karp



Schrijver



Vial



Maculan



Saad

Places where I have been (study/work/conference/cooperation/teaching)

University of Yaoundé (Cameroon) → University of Rennes/IRISA → University of Geneva (HEC/CUI)
→ European Laboratory of Molecular Biology → University Paris-Sud → Ecole des Mines de Paris

92: CARI

97: CEPAMOQ

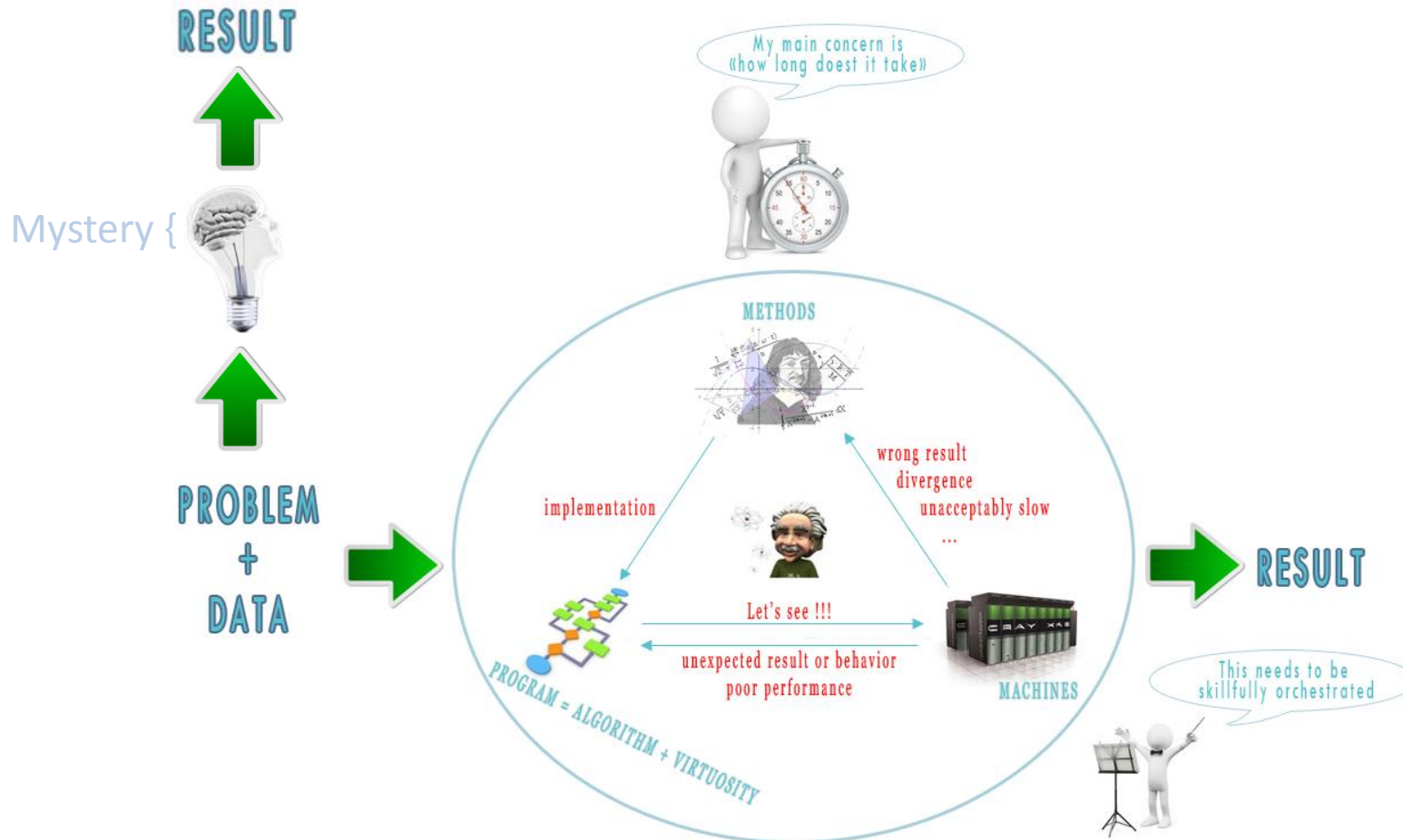


«science without border »
is the hallmark of my route.



France	Italy	Spain	Argentina	Austria
Switzerland	Germany	Hungary	Madagascar	Brazil
Romania	Portugal	Finland	Senegal	UK
Danmark	Greece	USA	Morroco	Ireland
Canada	Mexico	Cuba	Egypt	India
Madagascar	China	Japan	Cameroon	





- **Significant advances** have been achieved in each of the aforementioned aspects.
- A **skillfull combination** of all HPC components is really the key to **absolute efficiency**.
- This **expected pluridiscipline interaction** should be better done at the **earliest**.
- This is the **main point of my defense** and the motivation behind **my future plans**.

▶ TRAVELING SALESMAN PROBLEM

Given a list of cities and their pairwise distances, the task is to find a **shortest tour** that visits each city exactly once.

$$\left\{ \begin{array}{l} \text{minimize} \quad \sum_{i=1}^n \sum_{j=1}^n c_{ij} x_{ij} \\ \text{subject to} \quad \sum_{j=1}^n x_{ij} = 1 \quad i = 1, \dots, n, i \neq j \quad (2.4) \\ \sum_{i=1}^n x_{ij} = 1 \quad j = 1, \dots, n, i \neq j \\ x \in \{0, 1\}^{n \times n} \quad + \text{Subtour breaking constraints} \end{array} \right.$$



NP-complete Problem



Method

(Super)Computer

The TSP has an a priori $n!$ complexity. Solving any instance with $n = 25$ using the current world fastest supercomputer (TITAN-CRAY XK7) might require 25 years of calculations.

TITAN in 1min \approx 6 billion people calculating

"Tu quieres celeste que te cueste."

Applications: 24h/24h during 300 years



Transportation, logistic, genome sequencing, benchmark for optimization methods, ...

TSP heros: Applegate, Bixby, Chvatal, and Cook

- ▶ As difficult as the HAMILTONIAN CYCLE PROBLEM, which is **NP-COMplete** (Karp)
- ▶ Please, forget about brute force approach! (16 cities \Rightarrow 653 837 184 000 possibilities)
- ▶ Modern optimization method have shown optimistic performances on practical instances
- ▶ Competitive solutions are **parallel implementation of powerful optimization methods**
- ▶ TSP (VLSI-Bell Labs) of size 85 900 solved in 1.5 year (2004-2006) using a cluster of 96 2.8 GHz Intel Xeon and 32 2.4 GHz AMD Opteron connected with 100 MB ethernet.

● Approximation algorithm is also an interesting pragmatic way to go



- **Quicksort** is a worst-case n^2 algorithm, but is still preferred to the $n \log n$ **heapsort**.
- LP can be solved in a **polynomial time**, but the (**exponential**) **SIMPLEX** is still preferred.

▶ The simplex has led to the leading solver **CPLEX**



▶ We have implemented a nice interface between **CPLEX** and **MATLAB** in cooperation with **David Musicant** (Carleton College) and **Travis Johnson**.

Our interface (started in 2004) has been used and cited in several mathematical programming papers, including ours (*P-median, Portfolio, Energy minimization*).

Can be downloaded at

http://www.omegacomputer.com/staff/tadonki/using_cplex_with_matlab.htm

General Linear Programming

```
[obj, x, lambda, status, colstat, it] = lp_cplex(c, A, b, l, u, le, ge, maxIter, optimizer)
```

```
min c' * x with linear constraints
```

optimizer is a variable for optimizer choice, which can be one of the following

```
0 : primal simplex
1 : dual simplex
2 : network
3 : primal - dual barrier
```

Mixed Integer Programming

```
[obj, x, lambda, status, colstat, it] = mip_cplex(q, c, A, b, l, u, le, ge, binvar, genvar, maxIter, verbose)
```

```
min c' * x with linear constraints and integer variables ( binary or general )
```

Indices of binary variables are listed in *binvar*, and indices of general integer variables are listed in *genvar*

- **High-precision Lattice Quantum ChromoDynamics** simulations.
- The ANR project **PetaQCD** was targeting **256×128³** lattices.
- One evaluation of the *Dirac operator* on a **256×128³** lattice involves **256 × 128³ × 1500 ≈ 10¹²** (stencil) floating-point operations

With **10,000 cores**, we can roughly perform **500 × 10³ × 10⁶ = 5×10⁹ fps**



Our **256×128³** lattice would then require **200 seconds ≈ 3 minutes** for each evaluation of the *Dirac operator*.

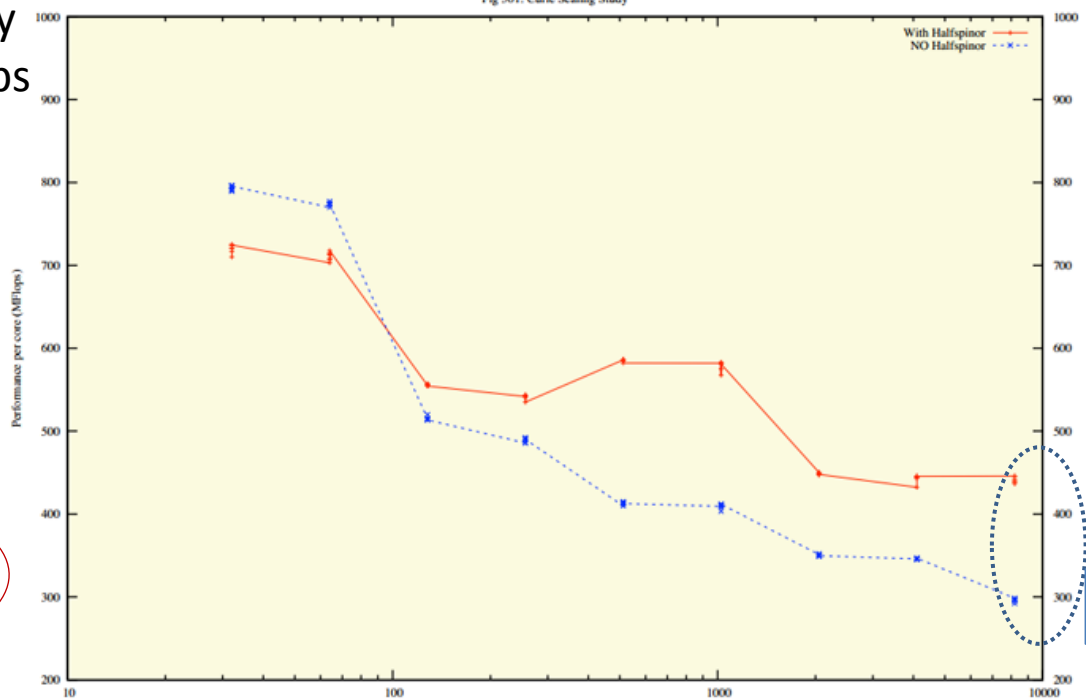
10 days !!!



Now, image that we have to do it 5000 times to solve One Dirac linear system !!!

Curie Fat performance (weak scaling)

Page 500: Version 7.0 G.Grosdidier 07/03/11 - TGCC-Curie (Last Update Time-stamp: <11/03/07 15:08:22 gridfs>)
Fig 501: Curie Scaling Study



G.Grosdidier, « Scaling stories », PetaQCD Final Review Meeting, Orsay, Sept. 27th – 28th 2012 **500 Mflops/core**

Important Facts about Supercomputers

➤ The (peak/sustained) performance of supercomputers is increasing significantly ($\approx \times 10^5$ since 1993).

The following characteristics are becoming a standard

- Several cores (2012 list: 84% ≥ 6 cores and 46% ≥ 8 cores)
- Vector units (with larger vector registers)
- Accelerated (mainly GPU: 62 systems of the 2012 including the #1)

● The gap between peak and sustained performances on real-life applications is clearly questionable.

☞ HPC investigations should focus on this.

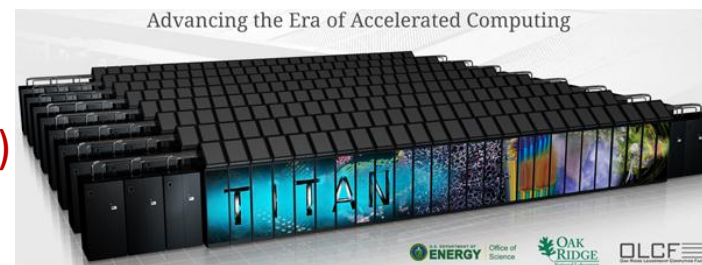
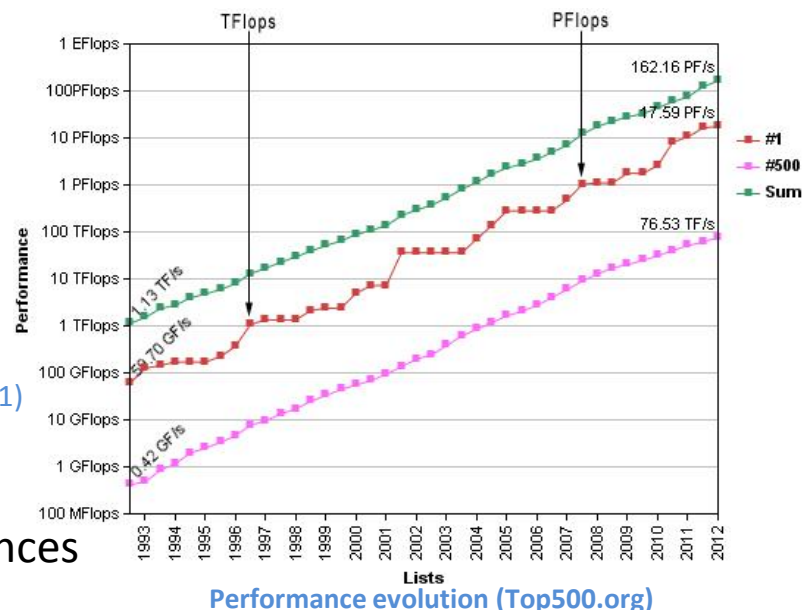
● When it comes to accelerators, data transfer is critical. ☞ Compromise and overlap (scheduling)

Chris Gregg and Kim Hazelwood, *Where is the Data? Why You Cannot Debate CPU vs. GPU Performance Without the Answer*, International Symposium on Performance Analysis of Systems and Software (ISPASS), Austin, TX. April 2011.

● Taking advantage of all aspects of a computing nodes requires a complex hybrid/heterogeneous programming. ☞ Code generation and transformation

● Energy is major concern.

☞ Power aware programming and scheduling.



TITAN CRAY-XK7 the (2012) world fastest supercomputer

- 299 008 CPU cores (16-cores AMD Opteron 6274)
- 18 688 NVIDIA Tesla K20 GPUs
- Peak: 27.11 PFlop/s.
- Sustained: 17.59 PFlop/s (Linpack)





Let's have a look on some
of my achievements

"Only those who attempt the absurd...will
achieve the impossible." M. C. Escher

Given a graph $G = (X, A)$ and a transition function φ such that $\varphi(G) = (X, A')$ with $\text{card}(A) = \text{card}(A')$. We study the evolution $G_0 = G, G_1 = \varphi(G_0), \dots, G_{n+1} = \varphi(G_n)$.

Remark: on a finite graph, such an evolution is ultimately **periodic**. Any state of the graph before the period is called **transcient**.

Interesting questions are

- ▶ the set of period lengths for a given graph
- ▶ bounds of period lengths
- ▶ garden of eden (graph with no transcient length)
- ▶ relation between the structure of the graph and periodicity
- ▶ applications (games, simulation, graph scheduling)

This work gave an inspiration to derive a parallel scheduling methodology for acyclic graphs with some recurrent structure.

René Ndoundam, Claude Tadonki, Maurice Tchuenta: *Parallel Chip Firing Game Associated with n- cube Edges Orientations*. International Conference on Computational Science (2004). , [CoRR abs/1007.0381](#) (2010)

A system of recurrence equations (SRE) defining a variable X has the following form

$$X(z) = \begin{cases} D_i^X & : g_i(\dots X(f(z)) \dots) \\ \vdots & \\ \vdots & \end{cases} \quad F(i, j, k) = \begin{cases} \{i, j, k \mid k = 0\} & : a_{i,j} \\ \{i, j, k \mid i = j = k\} & : F(i, j, k - 1) * \\ \{i, j, k \mid i = k \neq j\} & : F(k, k, k) \otimes F(i, j, k) \\ \{i, j, k \mid j = k \neq i\} & : F(i, j, k - 1) \otimes F(k, k, k) \\ \{i, j, k \mid i \neq k; j \neq k\} & : F(i, j, k - 1) \oplus \\ & (F(i, k, k) \otimes F(k, j, k - 1)) \end{cases}$$

- Algebraic and syntactic transformations (polyhedral model framework, compiler, scheduler, symbolic analysis, ...)
- Graph based approaches easily apply to SRE through the underlying dependence graph \Rightarrow systematic synthesis !

Scheduling a system of recurrence equations onto p processors is the task of finding a valid *timing function* T (i.e. $T(z) > T(f(z))$) and an allocation function A such that $[T(z_1) = T(z_2)] \Rightarrow [A(z_1) \neq A(z_2)]$.

Theorem 1. Let $G_1 = (X_1, F_1)$ and $G_2 = (X_2, F_2)$ two isomorphic graphs, and let φ an isomorphism from G_2 to G_1 . If (t, a) is a valid schedule of G_1 , where t is a timing function and a an allocation function, then $(t \circ \varphi, a \circ \varphi)$ is a valid schedule of G_2 .

Theorem 2. Let $G = (X, A)$ be a directed acyclic graph. If G is self isomorphic with the decomposition $X = X_1 \cup X_2 \cup \dots \cup X_\gamma$ and the set of isomorphisms $\{\varphi_k, k = 2, \dots, \gamma\}$, then the schedule specified by the timing function t and the allocation function a given by (6-7) is valid.

Applications: sorting, APP, Cholesky, tensor product of matrices

Our scheduling method (already published) opens the following perspectives

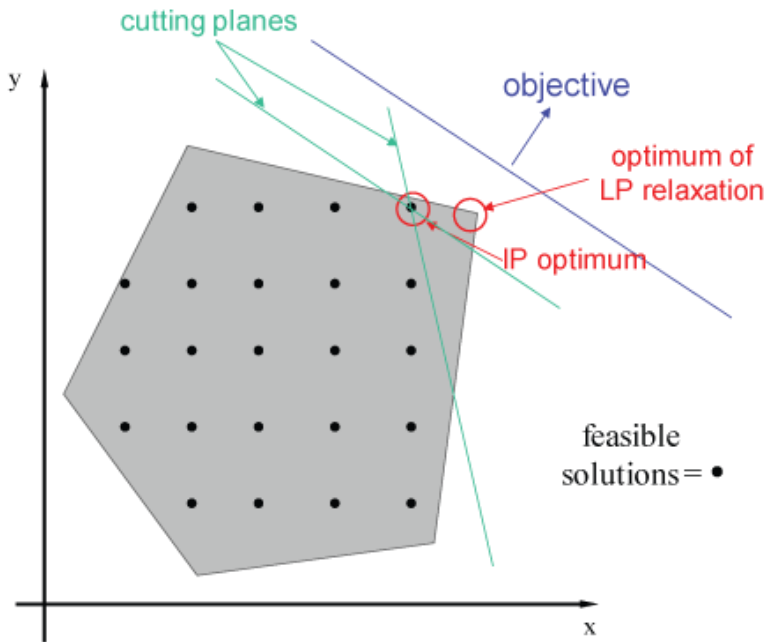
- Implementation as a scheduling module into a polyhedral model based framework.
- Study of an appropriate graph clustering (hybrid parallelism or modularity)
- Inclusion of hardware parameters



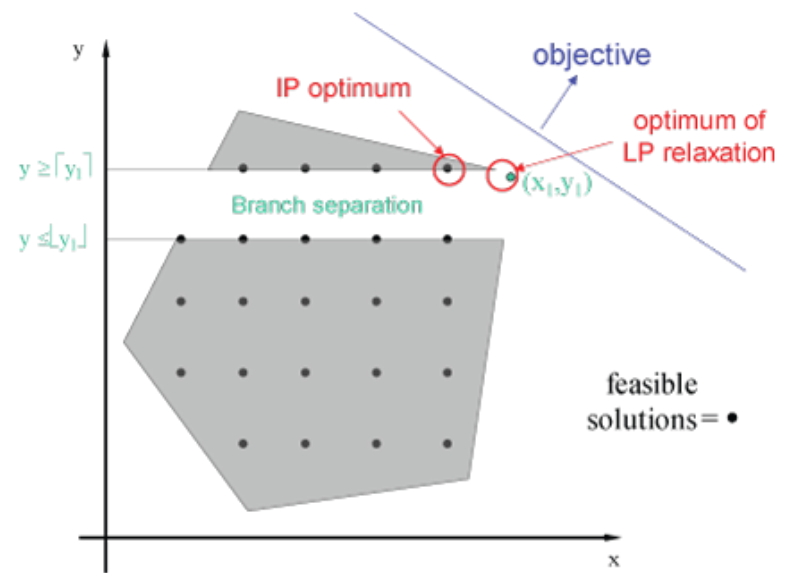
Inside the LOGILAB, Operation Research Laboratory
founded by Pr Jean-Philippe Vial and Pr Alain Haurie
at the University of Geneva (Switzerland)

*Dealing with modern, efficient, and somehow pragmatic approaches
for modeling and solving difficult real-world problems*

Think about gaussian pivoting vs iterative methods



Integer Program: Cutting Planes



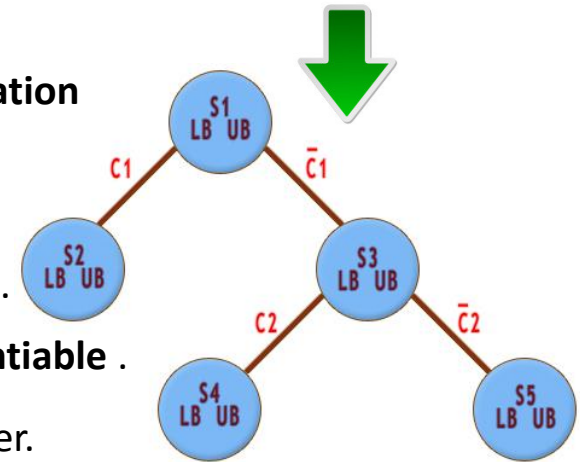
Integer Program: Branch and Bound
(or Divide and Conquer)

➤ Practical instances of **discrete** (pure or mixed) **optimization problems** are better solved through a skillfull combination of **continuous optimization** techniques and **branch&bound-like mechanisms**.

➤ For a pure discrete problem, a **relaxation** is used.
For a mixed formulation, a **decomposition approach** can be considered.

➤ In number of cases, the objective function is (or becomes) **non differentiable**.

➤ We then need a good non differentiable optimization method and solver.



$$\min\{f(y) = f_1(y) + b^T y + f_2(y) \mid y \in C\}, \quad (1)$$

where

- f_1 and f_2 are convex functions on R^n defined over $C \subset R^n$
- f_1 is nonsmooth and f_2 is twice differentiable
- C is a convex domain
- $b \in R^n$ is a constant vector

Remark 1 *The nonsmooth function f_1 is often a positively weighted sum of p nonsmooth functions*

$$f_1(y) = \sum_{i=1}^p \pi_i f_{1i}(y). \quad (2)$$

This property can be exploited in the method.

Remark 2 *For a given convex function f , there are many combinations of the form $f(y) = f_1(y) + b^T y + f_2(y)$. The appropriate choice is left to the user convenience.*

Project funded by the Swiss National Science Foundation

In order to accomplish the optimization task, *ProxAccpm* runs in conjunction with user defined *oracles*.

An *oracle* is a user defined routine (black box for ProxAccpm), which computes and returns informations about feasibility and/or optimality.

We consider tree types of oracle:

Feasibility oracle: Check if the current point y belongs to the optimization domain. Otherwise, its returns a so-called *feasibility cut*, we denote $(\xi, \tau) \in R^n \times R$, which satisfies

$$\xi^T(y' - y) + \tau \leq 0, \forall y' \in C. \quad (3)$$

First order oracle: Given a feasible point y , it returns $f_1(y)$ and one subgradient $u \in \partial f_1$. This yields the following relation:

$$u^T(y' - y) + (f_1(y') - f_1(y)) \leq 0, \forall y' \in C. \quad (4)$$

Second order oracle: Given a feasible point y , it returns $f_2(y)$, $f'_2(y)$, and $f''_2(y)$.

Localization Set: At a given step, the *localization set* is defined by

$$\mathcal{L}_{\bar{\theta}} = \{(y, z, \zeta) \mid A^T y - E^T z \leq c, \pi^T z + b^T y + \zeta \leq \bar{\theta}, f_2(y) \leq \zeta, y \in Y_2\}, \quad (7)$$

where

- $\bar{\theta}$ is the current upper bound
- A is the matrix of cuts (columnwise)
- E is a binary banded matrix

Methods differ in

- the management of the localization set (objective cut, updating, spatial transformations, etc...)
- the selection of the query point from the localization set
- the lower bound and the termination criterion

Accpm and Proximal Accpm

The Analytic Center Method: The query point is obtained by minimizing (over the localization set) the weighted logarithmic barrier given by

$$F(\bar{s}) = - \sum w_i \log s_i - \omega \log \sigma, \quad (8)$$

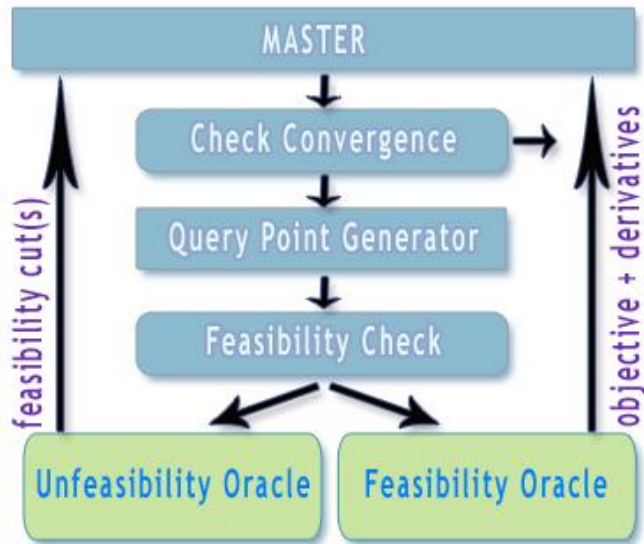
with $\bar{s} = (s_0, s, \sigma) > 0$ defined by

$$\begin{aligned} s_0 &= \bar{\theta} - (\pi^T z + b^T y + \zeta) \\ s_i &= c_i - (A^T y - E^T z)_j, \quad i \in I = \{1, \dots, m\}, \\ \sigma &= \zeta - f_2(y). \end{aligned}$$

The Proximal Analytic Center Method: The barrier function is augmented with a proximal term to yield the augmented barrier

$$\frac{1}{2}(y - \bar{y})^T Q (y - \bar{y}) + F(\bar{s}) \quad (9)$$

where Q is a positive definite matrix and \bar{y} the so-called *proximal center*..



The main concerns are

- force the global convergence
- minimize the number of oracle calls
- optimize the computation time and memory space
- improve the computation accuracy
- take care of numerical stability

Babonneau, F., Beltran, C., Haurie, A., Tadonki, C. and Vial, J.-P., *Proximal-ACCPM: a versatile oracle based optimization method*, In *Optimisation, Econometric and Financial Analysis*, E. J. Kontogiorghes editor, vol. 9 of *Advances in Computational Management Science*, 2006.

Achievements

- Implementation of the method in a **complete framework** (<https://projects.coin-or.org/OBOE>).
- Implementation as a query point generator for **connection with other packages** (*branch and bound*)
- Several case studies **published in journal and conferences**

Some perspectives

- Deep investigation of the **parallelization** on supercomputers (*scheduling, scalability, load balance,...*)
- Study how to deal with **updates** instead of performing matrix computations from scratch at each step
Sherman-Morrison formula
- Investigate on numerical issues with large-scale ill-conditioned systems
- Approximation algorithms

Energy Minimization

Power consumption is a crucial concern with embedded systems and supercomputers

Fujitsu K-Computer \approx US\$10 millions/year for electricity

Pr Jose Rolim

Theoretical Computer Science Laboratory (TCS Lab)

Centre Universitaire Informatique

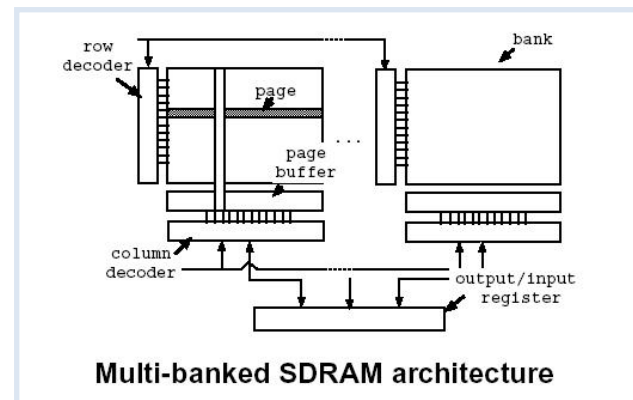
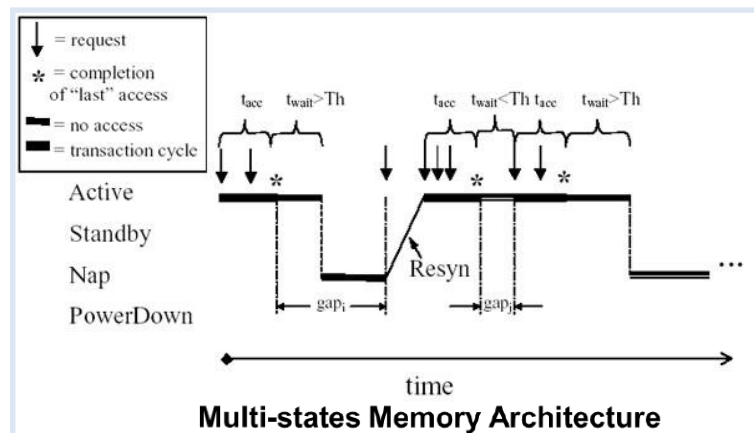
University of Geneva (Switzerland)



My work on Power Aware Computing and Distributed Algorithms were initiated from my stay at the TCS-Lab.

- We focus on memory energy
- The memory is partitioned into several banks
- Each bank can be put into a specific power mode
- We assume regular (uniform) power state transitions
- We formulate and solve the optimization problem

$$\begin{aligned} & \min x^T H^T Q y^T + R Q y^T \\ & \text{subject to} \\ & 1. \quad x \in \mathbb{N}^q, \\ & 2. \quad y \in \mathbb{R}^q, \\ & 3. \quad y_1 + y_2 + \dots + y_q = p, \\ & 4. \quad y_0 \geq \mu C, \\ & 5. \quad x_1 + x_2 + \dots + x_q = \rho C, \\ & 6. \quad y \leq \varphi x. \\ & 7. \quad y \geq \eta x. \end{aligned}$$



Good collaboration with Mitali Singh and Viktor Prasanna (University of Southern California)

Input

Hardware parameters : $W, Q, H, p, q,$ and δ

Program parameters : R, C

Memory management parameters : ρ, φ, η

Model and Optimization



Output

Optimal Energy E

Optimal transition repartition X

Optimal time repartition Y

Power aware
program design
and monitoring

Some perspectives are:

- Design a **methodology** that will use the output of our work for a systematic synthesis of **energy efficient policies**
- Extend and adapt our model to **current and future memory systems** (*multilevel and shared*)
- Use a similar formal approach to analyse the energy issue on the **cloud systems** (user and provider)
- Investigate on other approaches (dynamic scheduling, compilation, ...)

*One of our proposal on this topic received an important grant from the
Swiss National Science Foundation*

We have been also investigating on:

- Distributed algorithm in sensors networks (*localization and information retrieval*)
- Dual-power management problem (*mathematical programming approach and heuristics*)
- Algorithms for the web (*search engines and social networks*)

PetaQCD – Overview



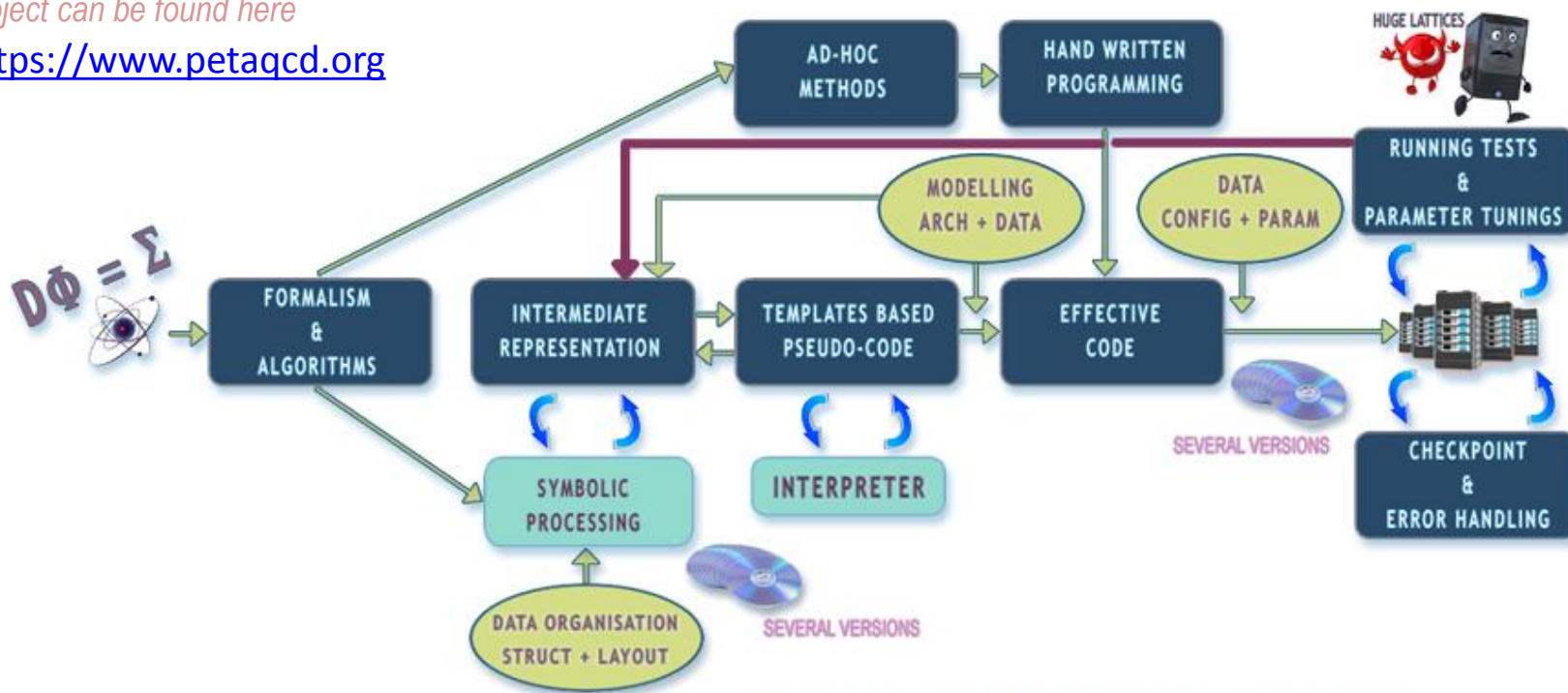
- ANR Project (HPC & Particles Physics)
- Origin of the universe (matter)
- Good cooperative effort
- Multidisciplinary collaboration
- HPC & numerical challenges

National: G. Grosdidier (*coordinator*), P. Roudeau, O. Pène, C. Tadonki, K. Petrov, D. Barthou, M. Kruse, C. Eisenbeis, B. Blossier, L. Morin, F. Bodin, F. Touze, O. Brand-Foissac, J. C. Angles d'Auriac, A. Sez nec, A. Cohen, C. Bastoul, and more.

Abroad: K. Urbach, K. Jansen, L. Scorzatto, D. Pleiter, R. Tripiccione

Lot of materials and outputs of this project can be found here

<https://www.petaqcd.org>



PETAQCD WORKING CHAIN

- Hybrid extension of the reference package (tmLQCD) using *Pthread* library
- Accelerated algorithm & implementation of the **Dirac operator** and other linear algebra kernels

Illustrative results

- A 32×16^3 configuration solved using the CGR algorithm on the PPU in **138 seconds**
- The same configuration and algorithm on the (PPU + 8 SPEs) double precision in **4.58 seconds**

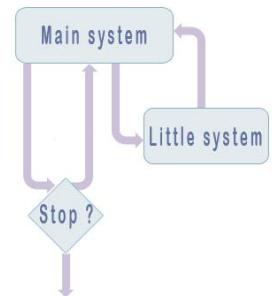
CGR: 57 iterations in 4.58 s (QS20) 3.68 s (QS22) 28.34s s (Intel 2.83Ghz)
 CG: 685 iterations in 51.70 s (QS20) 38.80 s (QS22) 362.45 s (Intel 2.83Ghz)

- Compare to the PPU, a speedup around **30** was obtained (also valid per iteration)

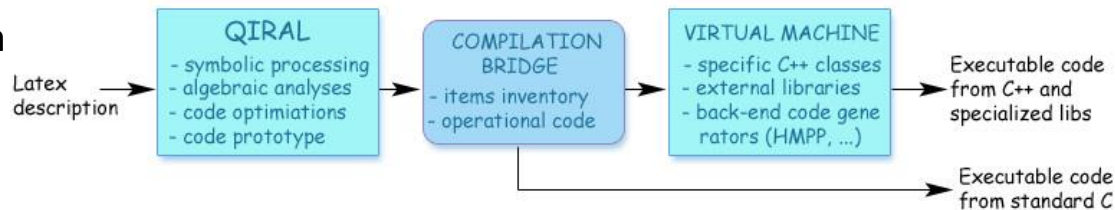
C. Tadonki, G. Grosdidier, and O. Pene, « *An efficient CELL library for Lattice Quantum Chromodynamics* », ACM SIGARCH Computer Architecture News, vol 38(4) 2011.

- Block decomposition C. Tadonki, «Strengthening deflation implementation» for large scale LQCD inversions, Orsay, Sept. 27th – 28th 2012

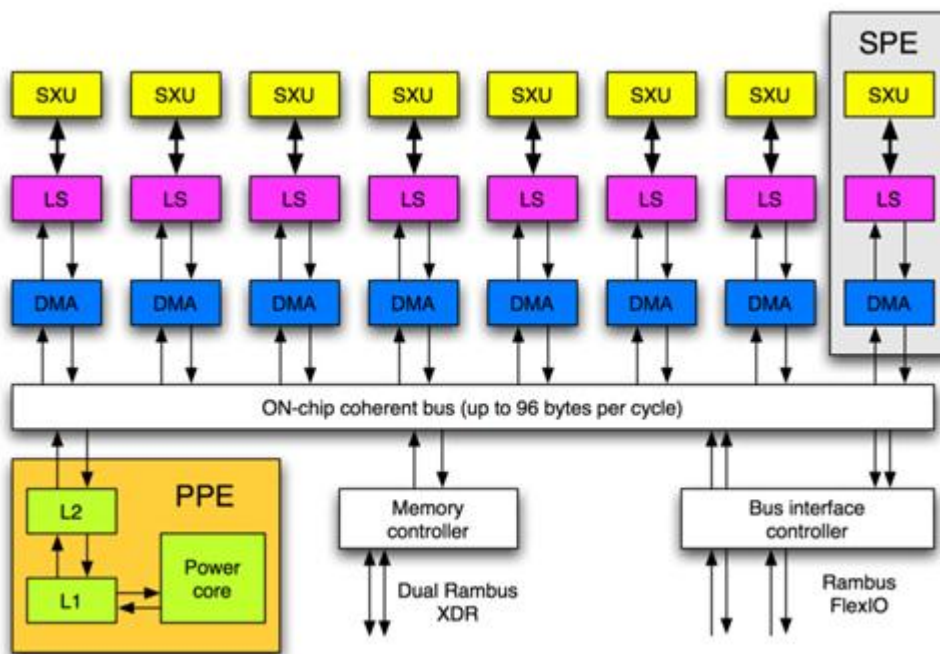
We have implemented a generic block decomposition (**multidimensional** and with no restriction on the **number of blocks per axis**) within the tmLQCD package. The aim is to increase the basis of the deflation method (to solve large ill-conditioned systems) without



- Code generation



D. Barthou, G. Grosdidier, M. Kruse, O. Pene and C. Tadonki, « *QIRAL: A High Level Language for Lattice QCD Code Generation* », Programming Language Approaches to Concurrency and Communication-centric Software ([PLACES'12](#)) in conjunction with the European joint Conference on Theory & Practice of Software ([ETAPS](#)), Tallinn, Estonia, March 24-April 1, 2012.



*I start investigating on this topic from my collaboration with Dr **Lionel Lacassagne** (ANR project *Ocelle* 2007-2009), also SIMD and Image Processing stuffs.*

*Good collaboration with Dr **Joel Falcou**, **Tarik Saidani**, **Khaled Hamidouche**, and Pr **Daniel Etiemble***

T. Saidani, J. Falcou, C. Tadonki, L. Lacassagne, and D. Etiemble,
« Algorithmic Skeletons within an Embedded Domain Specific Language for the CELL Processor »,
Parallel Architectures and Compilation Techniques (PACT), [PACT09](#), Raleigh, North Carolina (USA), September 12-16, 2009.

DMA issues related to tiling

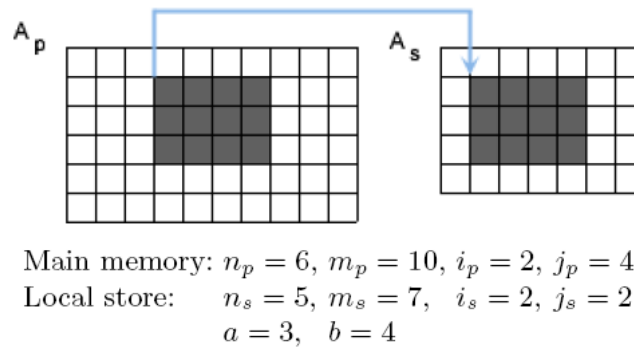


Fig. 4. Generic DMA pattern

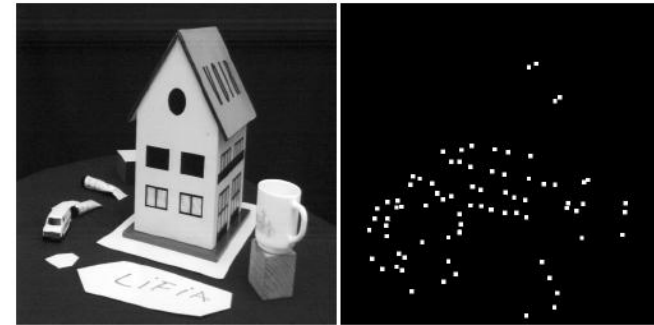
Performing the transfer expressed in figure 4 raises number of problems:

- *the region to be transferred is not contiguous on memory, thus list DMAs are considered*
- *the address of one given row is not aligned, thus the global list DMA is not possible*
- *the (address, volume) pair of a row does not match the basic DMA rules (the above two ones), thus the entire list DMA cannot be carried out*
- *misalignment could come from both sides (main memory and/or local store)*
- *the target region on the local store might be out of the container limits*

We have designed and implemented a routine which performs this task very efficiently

The Harris-Stephen algorithm is

- a corner (point of interest) detection algorithm
- an improved variant of the original algorithm by Moravec
- used in computer vision for feature extraction like
 - motion detection
 - image matching
 - tracking
 - 3D reconstruction
 - object recognition



Technically, the Harris algorithm is based on a pixelwise autocorrelation S given by

$$S(x, y) = \sum_{u, v} w(u, v) [I(x, y) - I(x - u, y - v)]^2$$

where (x, y) is the location of the pixel and $I(x, y)$ its intensity (grayscale mode).

$tile_h$	$tile_w$	total time(s)
8	512	0.0494
16	256	0.0598
32	128	0.0485
64	64	0.0345
128	32	0.0517
256	16	0.0699
512	8	0.0734

Table 1. Timings on a 512×512 image

$tile_h$	$tile_w$	total time(s)
8	512	0.198
16	256	0.238
32	128	0.187
64	64	0.110
128	32	0.180
256	16	0.218
512	8	0.352

Table 2. Timings on a 2048×512 image

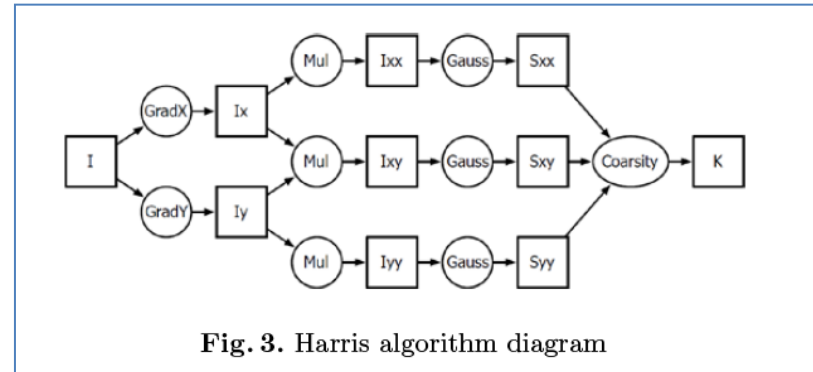


Fig. 3. Harris algorithm diagram

$tile_h$	$tile_w$	total time(s)
5	1200	0.494
10	600	0.360
20	300	0.264
40	150	0.235
80	75	0.183
160	37	0.247
320	18	0.275

Table 3. Timings on a 1200×1200 image

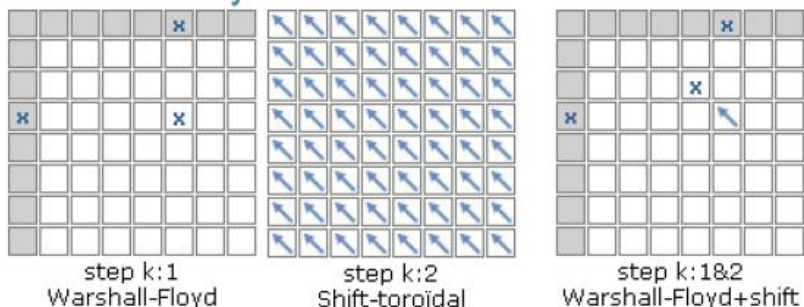
$tile_h$	$tile_w$	total time(s)
8	512	0.985
16	256	0.726
32	128	0.643
64	64	0.438
128	32	0.692
256	16	0.866
512	8	1.422

Table 4. Timings on a 2048×2048 image

C. Tadonki, L. Lacassagne, E. Dadi, M. Daoudi,
 «Accelerator-based implementation of the Harris algorithm»,
 5th International Conference on Image Processing ([ICISP 2012](#)),
 June 28-30, 2012.

We observe **50% improvement** between square tiles and full row tiles.

Warshall-Floyd + Shift-toroidal

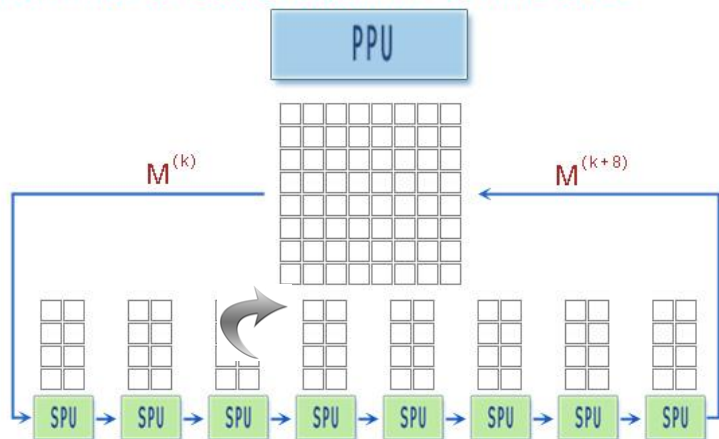


Analytical expression of the shift-toroidal



$$m_{i-1,j-1}^{(k)} = m_{ij} \oplus (m_{ik}^{(k-1)} \otimes m_{kj}^{(k-1)})$$

MAPPING OF OUR ALGORITHM ON THE CELL



- › PPE-DMA is issued only by the first and the last processor
- › Inner SPEs communicate and synchronize locally
- › Computation-communication overlap occurs for all communications
- › Can run on more SPEs or CELL Blades by natural extension

	1 SPE		2 SPEs		8 SPEs	
Tile	t(s)	t(s)	σ	t(s)	σ	
1	6.67	3.28	2.03	1.60	4.16	
4	5.01	2.50	2.00	0.62	7.99	
8	4.79	2.39	2.00	0.60	7.95	
12	4.70	2.32	2.02	0.58	7.98	
16	4.72	2.36	2.00	0.60	7.79	

(c) Performance with a 1024x1024 matrix

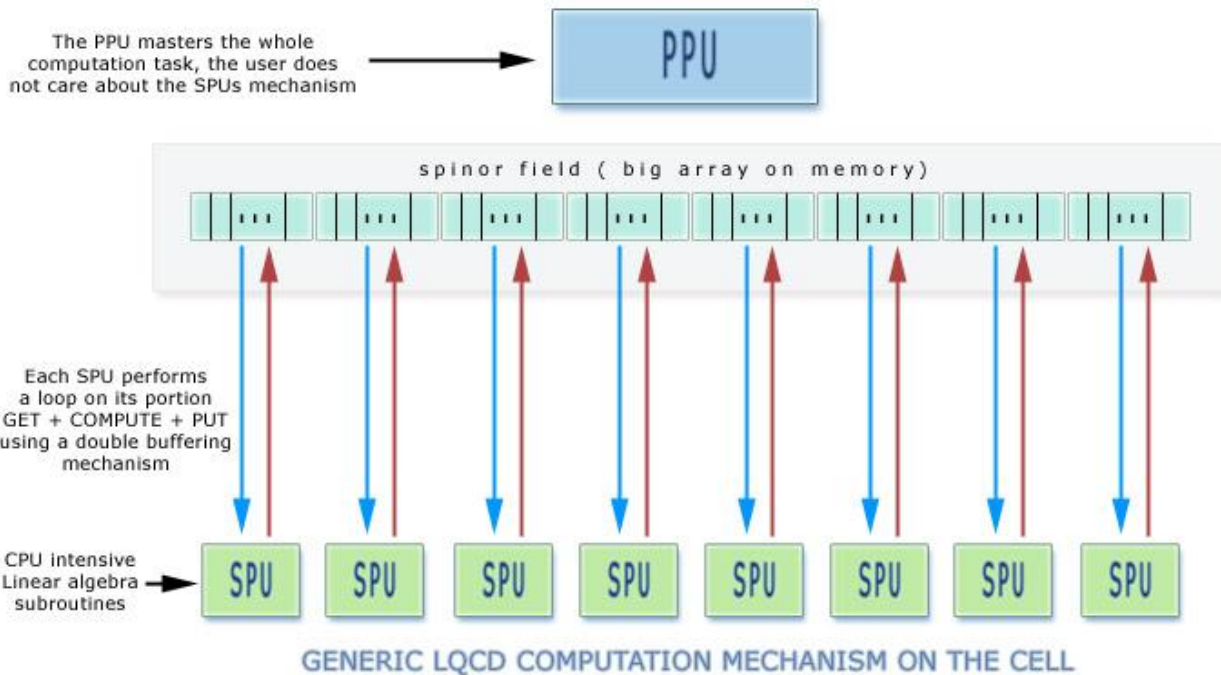
DMA timings (1024x1024 matrix)

	1 SPE		2 SPEs		8 SPEs	
Tile	t(s)	t(s)	σ	t(s)	σ	
1	1.87	1.20	1.56	0.53	3.49	
4	0.39	0.47	0.83	0.11	3.70	
8	0.19	0.12	1.64	0.06	3.78	
12	0.12	0.08	1.65	0.03	4.02	
16	0.09	0.06	1.63	0.03	3.78	

Benefits of the shift-toroidal

- ✓ The pivot row (column) remains on top
- ✓ Shift + computation can be done simultaneously
- ✓ All steps become rigorously identical
- ✓ The overhead of the shift can be hidden
- ✓ Does not change the $O(n^3)$ complexity

C. Tadonki, « Ring pipelined algorithm for the algebraic path problem on the CELL Broadband Engine », [WAMMCA 2010](#) - [SBAC PAD 2010](#), Petropolis, Rio de Janeiro, Brazil, October 27-30, 2010. (IEEE digital library)



- SIMD implementation of the basic operators (intensive use of SPU intrinsics)
- double buffering technique to overlap DMA and computation (ideal in double precision)
- optimal list DMA organisation by an algorithm similar that used for the *bin packing* problem (significant latency reduction)

C. Tadonki, G. Grosdidier, and O. Pene, «An efficient CELL library for Lattice Quantum Chromodynamics», [HEART](#) - ACM/ICS, Epochal Tsukuba, Tsukuba, Japan, June 1-4, 2010.

1 Performance results

All our experimentations are performed with *double precision* data. We used a 32×16^3 lattice (hence 131 072 sites)

#SPEs	time(s)	speedup	GFlops
1	0.109	1.00	0.95
2	0.054	2.00	1.92
3	0.036	3.00	2.89
4	0.027	3.99	3.85
5	0.022	4.98	4.73
6	0.018	5.96	5.78
7	0.015	6.93	6.94
8	0.013	7.88	8.01

Figure 1: 32×16^3 Wilson-Dirac on a QS20

#SPEs	time(s)	speedup	GFlops
1	0.0374	1.00	2.76
2	0.0195	1.91	5.31
3	0.0134	2.79	7.76
4	0.0105	3.56	9.90
5	0.009	4.81	13.27
6	0.0081	5.75	15.87
7	0.0076	6.84	18.88
8	0.0075	7.82	21.59

Figure 2: 32×16^3 Wilson-Dirac on a QS22

Without SSE		With SSE	
1 core	4 cores	1 core	4 cores
0.0820	0.0370	0.040	0.0280

Figure 3: 32×16^3 Wilson-Dirac timings (seconds) on an INTEL i7 quadcore 2.83 Ghz

UNIVERSAL REPORT: A Universal Source Code Analysis and Documentation Software

```

{
  double s[3], x, y, z, yy, zz, r;
  int i;
  r = 0.0;

  for( i = 0; i < 3; i++ )
  {
    x = pp[i];
    s[i] = x;
    r += x * x;
  }

  r = sqrt(r);

  if( ofdate )
  {
    precess( s, J, -1 );
    epsilon(J);
  }

  if( prtflg == 0 )
  {
    return(0);
  }

  printf( "edictic long" );
  dms( yy );
  printf( " lat" );
  dms( zz );
  printf( " rad %.9E\n", r );
  return(0);
}
    
```

Basic, C, C++, COBOL, Fortran, Java, Javascript, Matlab, Pascal, Visual Basic, Borland C++ Builder, Delphi, Kylix, Perl, PL1, Python, Visual C++, Visual Basic .Net, Visual C#, Visual J++, and more





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- Provides several options and parameters
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- Automatic screenshots and database reporting
- Automatic changes analysis and multi-reporting
- Minimum user effort and real-time processing
- Special character sets (arabic, chinese, hebrew, japanese, korean, etc ...)
- HTML FlowChart for each routine (example)
- Translation of comments (Google engine) **NEW!!!**

Features overview: recommend universal report

People involved with programming activities know maintaining (large/complex) codes is tedious. Programs that you have written yourself are difficult enough to understand when you come back to them later. Things are more difficult when the program have been written by someone else or a big team, and has been modified and added so many times that the existing documentation no longer accurately reflects the program code and structure. This is a nightmare awaiting those involved with program maintenance, comprehension, audit, or analysis. Therefore, tools that automate the process of code analysis and documentation clearly help to overcome this problem painlessly.

UNIVERSAL REPORT is a high quality code analysis and documentation software. Its goal is to analyse and generate a structured and well formatted documentation of a given program. Universal Report can handle various programming languages (Basic, C, C++, COBOL, Fortran, Java, Javascript, Matlab, Pascal, Visual Basic, Borland C++ Builder, Delphi, Kylix, Perl, PL1, Python, Visual C++, Visual Basic .Net, Visual C#, Visual J++, and more), and generates html, latex, and plain text documentations using state-of-the-art pattern matching algorithms and compilation techniques. The tool is parameterizable both in the behavior point of view and the quality/layout of the outputs. [Features]

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Claude Tadonki , « Universal Report: A Generic Reverse Engineering Tool »,
12th IEEE International Workshop on Program Comprehension, [IWPC 2004](#) , University of Bari, Bari, Italy , June 2004 ,

- Cooperation with Brazil
 - **BioCloud** project of the STIC-AmSud
 - Sandwich PhDs (*Brazilia/Rio/Niteroi*)
 - Papers, visitings, events, courses
 - Other Latin America partners (*Chile, ...*)
 - + INRIA and Paris-Sud University
- Cooperation with Morocco (*Oujda*)
 - Sandwich PhDs
 - Papers, visitings, seminars
- Cooperation at Mines ParisTech
 - TIMC project (multi-target images)
 - CMM (Math Morphology) and CAOR



BioCloud Partners at Orsay

A screenshot of the CPOCC workshop website. The page features a navigation menu with buttons for HOME, IMPORTANT DATES, SUBMISSION, COMMITTEES, REGISTRATION, CAMERA READY, PROGRAM, and NEWS. The main content area includes a title 'CPOCC' and a subtitle '1st Workshop on Cost and Performance Optimization in Cloud Computing'. Below this, there is a paragraph of introductory text about cloud computing and the workshop's goals. A 'TOPICS' section lists various areas of interest like power consumption models, cost models, federated cloud models, etc. A 'WORKSHOP ORGANIZATION' section lists the names and affiliations of the organizers. On the right side, there is an image of a globe and server racks. At the bottom, there is a graphic with the text 'Workshop on Cost and Performance Optimization in Cloud Computing' and 'CLOUD COMPUTING' in large letters, accompanied by icons of a laptop and a tablet.



In memory of Jean Tadonki (1939-2001)
« You are the greatest »

Thanks to all of you



Family, friends, colleagues, collaborators, advisors, examiners, referees, administration, technicians, audience, you, ...