

Modeling Power and Energy of the Task-Parallel Cholesky Factorization on Multicore Processors

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Motivation

- High performance computing:
 - Optimization of algorithms applied to solve complex problems
- Technological advance $\Rightarrow$ improve performance:
 - Higher number of cores per socket (processor)
- Large number of processors and cores $\Rightarrow$ **High energy consumption**
- Tools to analyze performance and power in order to detect code inefficiencies and **reduce energy consumption**

Outline

- 1 Introduction
- 2 Task-parallelism in the Cholesky factorization
 - Algorithm specification
 - Parallelization
 - SMPs operation
- 3 Power model
 - Formulation
 - Environment setup
 - Component estimation
 - Power/energy model testing
- 4 Experimental results
 - Energy model evaluation
 - Power model evaluation
- 5 Conclusions and future work

Introduction

- **Parallel scientific applications**

- Examples for dense linear algebra: Cholesky, QR and LU factorizations

- **Tools for power and energy analysis**

- Power profiling in combination with performance/tracing tools for HPC

Parallel applications + Power profiling



Is it possible to predict power/energy consumption?

- **Objective:** Power modeling

- Predict power consumed by applications without power measurement devices.
- Estimations are needed to determine how to address the power-challenge for energy-efficient hardware and software



Energy savings

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Is it possible to predict power/energy consumption?

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- Predict power consumed by applications without power measurement devices.
- Estimations are needed to determine how to address the power-challenge for energy-efficient hardware and software



Energy savings

Algorithm specification

Cholesky factorization:

$$A = U^T U$$

$A \in \mathbb{R}^{n \times n}$ symmetric definite positive (s.p.d.) matrix

$U \in \mathbb{R}^{n \times n}$ unit upper triangular matrix

$\Rightarrow$ Consider a partitioning of matrix A into blocks of size $b \times b$

```

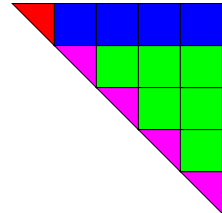
for k = 1, 2, ..., s do
   $A_{kk} = U_{kk}^T U_{kk}$ 
  for j = k + 1, k + 2, ..., s do
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  for i = k + 1, k + 2, ..., s do
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```

CHOL: CHOLESKY FACTORIZATION

TRSM: TRIANGULAR SYSTEM SOLVE

SYRK: SYMMETRIC RANK- b UPDATE

GEMM: MATRIX-MATRIX PRODUCT



Iteration 1

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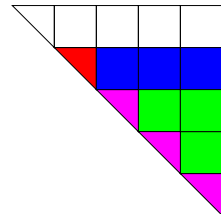
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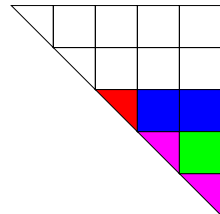
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Iteration 3

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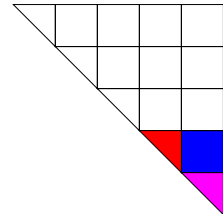
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```

```
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```
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```

```
  end for
```

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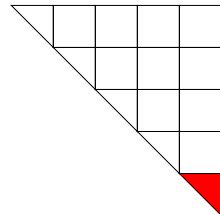
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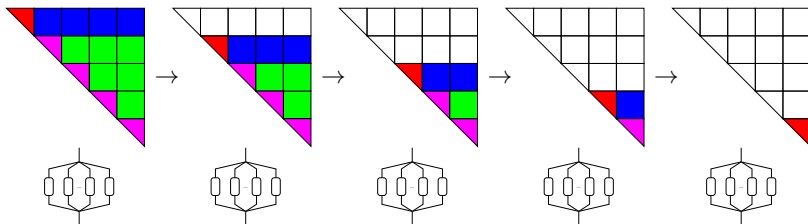
Iteration 5

Parallelization $\Rightarrow$ Not trivial at code level!

Parallelization

Option 1: Use **multi-threaded BLAS**

- Straightforward approach towards LAPACK-level parallelization
- Highly tuned multi-threaded kernels: Intel MKL, AMD ACML or IBM ESSL,...
- Fork/join approach: parallelism is not fully exploited



Parallelization

Option 2: Use a **runtime task scheduler**

- We use SMPs runtime-compiler framework to exploit task-parallelism
- Functions in code are annotated as tasks using OpenMP-like pragmas `#pragma cxx task`
- Operations are not executed in the order they appear in the code but respecting data dependencies
- SMPs easily obtains performance traces which can be analyzed using *Paraver* (Performance analysis tools from Barcelona Supercomputing Center)

SMPs proceeds in two stages:

- 1 A symbolic execution produces a DAG containing dependencies
- 2 DAG dictates the feasible orderings in which task can be executed

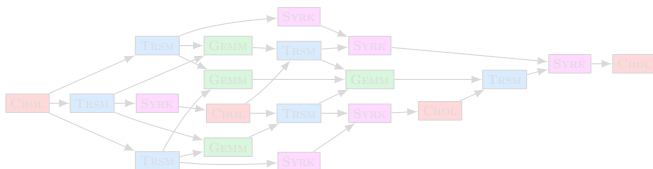


Figure: Right-looking Cholesky DAG with a matrix consisting of 4×4 blocks

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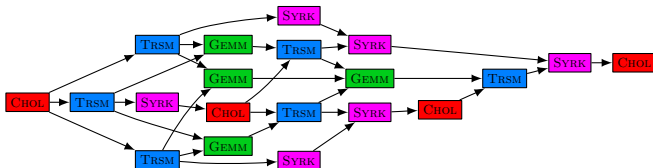


Figure: Right-looking Cholesky DAG with a matrix consisting of 4×4 blocks

Cholesky factorization with SMPs pragma annotations

```
void dpotrf_smpss( int n, int b, double *A, int Alda, int *info ){
    for (k=1; k<=n; k+=b) {
        dpotrf_u( b, &A-ref(k,k), Alda, info );
        if ( k+b <= n ) {
            for (j=k+b; k<=n; k+=b)
                dtrsm_lu1n( b, &A-ref( k, k ), &A-ref( k, j ), Alda );
            for (i=k+b; i<=n; i+=b) {
                dsyrk_ut( b, &A-ref( k, i ), &A-ref( i, i ), Alda );
                for (j=i+b; j<=n; j+=b)
                    dgemm_tn( b, &A-ref( k, i ), &A-ref( k, j ), &A-ref( i, j ), Alda );
            }
        }
    }
}

void dpotrf_u( int b, double A[], int ldm, int *info ){
    dpotrf( "Upper", &b, A, &ldm, info );
}

void dtrsm_lu1n( int b, double A[], double B[], int ldm ){
    double done = 1.0;
    dtrsm( "Left", "Upper", "Transpose", "Non-unit", &b, &b, &done, A, &ldm, B, &ldm );
}

void dsyrk_ut( int b, double A[], double C[], int ldm ){
    double dmone = -1.0, done = 1.0;
    dsyrk( "Upper", "Transpose", &b, &b, &dmone, A, &ldm, &done, C, &ldm );
}

void dgemm_tn( int b, double A[], double B[], double C[], int ldm ){
    double dmone = -1.0, done = 1.0;
    dgemm( "Transpose", "No_transpose", &b, &b, &b, &dmone, A, &ldm, B, &ldm, &done, C, &ldm );
}
```

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            }
        }
    }
}

#pragma css task input( b, ldm ) inout( A[1], info[1] )
void dpotrf_u( int b, double A[], int ldm, int *info ){
    dpotrf( "Upper", &b, A, &ldm, info );
}

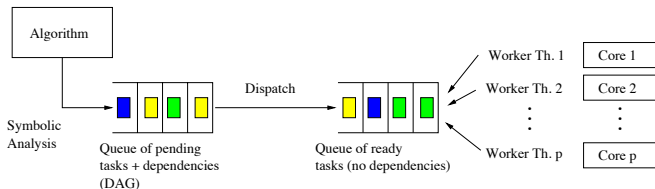
#pragma css task input( b, A[1], ldm ) inout( B[1] )
void dtrsm_lu1n( int b, double A[], double B[], int ldm ){
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#pragma css task input( b, A[1], ldm ) inout( C[1] )
void dsyrk_ut( int b, double A[], double C[], int ldm ){
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    dsyrk( "Upper", "Transpose", &b, &b, &dmone, A, &ldm, &done, C, &ldm );
}

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}
```

SMPs operation

SMPs runtime:



Basic scheduling:

- ① Initially only one a task in *ready queue*
- ② A thread acquires a task of the *ready queue* and runs the corresponding job
- ③ Upon completion checks tasks which were in the *pending queue* moving to *ready* if their dependencies are satisfied.

Power model formulation

Power model:

$$P = P^{C(PU)} + P^{S(Y)stem} = P^{S(tatic)} + P^{D(ynamic)} + P^{S(Y)stem}$$

$P^{C(PU)}$ Power dissipated by the CPU: $P^{S(tatic)} + P^{D(ynamic)}$

$P^{S(Y)stem}$ Power of remaining components (e.g. RAM)

Considerations:

- Study case: **Cholesky factorization**. It exercises CPU+RAM and discards other power sinks (network interface, PSU, etc.)
- We assume P^Y and P^S are constants!
- P^S grows with the temperature inertia till maximum! $\Rightarrow$ We consider a "hot" system!

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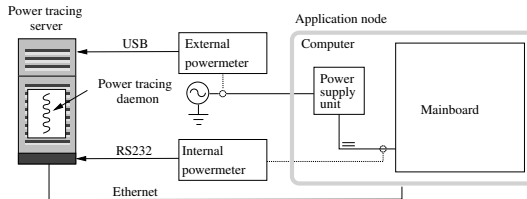
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Environment setup

Setup:

- Intel Xeon E5504 (2 quad-cores, total of 8 cores) @ 2.00 GHz with 32 GB RAM
- Intel MKL 10.3.9 for sequential `dpotrf`, `dtrsm`, `dsyrk` and `dgemm` kernels
- SMPSs 2.5 for task-level parallelism
- Performance and tracing modes are enabled
- Power measurements: `pmlib` library

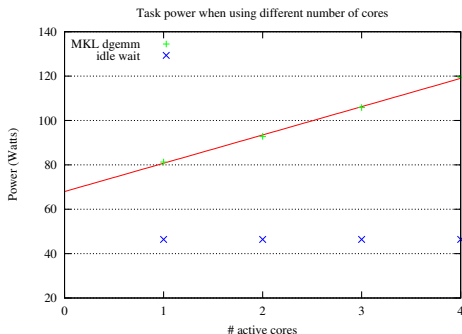


- Internal power meter:
 - ASIC-based powermeter (own design!)
 - LEM HXS 20-NP transducers with PIC microcontroller
 - Sampling rate: 25 Hz

System and static power

Obtaining $P^{S(Y)stem}$ and $P^{S(tatic)}$ components:

- P^Y directly obtained measuring idle platform: $P^Y = 46.37$ Watts
- P^S obtained by executing **dgemm** kernel using 1 to 4 cores and adjusting via linear regression:



Linear regression: $P_{\text{dgemm}}(c) = \alpha + \beta \cdot c = 67.97 + 12.75 \cdot c$

$P^S \approx \alpha - P^Y = 67.97 - 46.37 = 21.6$ Watts

Dynamic power

Dynamic power of kernels of the Cholesky factorization:

- To obtain P_K^D we continuously invoke the kernel K until power stabilizes and then sample this value. Example for dgemm:

$$P_G^D = P_{\text{dgemm}} - P^S - P^Y = P_{\text{dgemm}} - 67.97 \text{ Watts}$$

Task	1 kernel mapped to 1 core				2 kernels mapped to 2 cores of different sockets			
	Block size, b				Block size, b			
	128	192	256	512	128	192	256	512
P_P^D (dpotrf)	10.26	10.35	10.45	11.28	9.05	9.09	9.28	10.44
P_T^D (dtrsm)	10.12	10.31	10.32	10.80	9.45	9.57	9.60	11.08
P_S^D (dsyrk)	11.22	11.47	11.67	12.60	10.42	10.63	10.82	11.80
P_G^D (dgemm)	11.98	12.54	12.72	13.30	10.90	12.16	11.28	11.96
P_B^D (busy)	7.62	7.62	7.62	7.62	7.62	7.62	7.62	7.62

- Power increases linearly with the number of threads, from 1 to 4 mapped to a single core
- When two sockets are used, linear function changes, so we take into account this issue:

$$P_G^D = \frac{P_{\text{dgemm}} - 67.97}{2}$$

Power/energy model testing

Power model:

$$P_{Chol}(t) = P^Y + P^S + P_{Chol}^D(t) = P^Y + P^S + \sum_{i=1}^r \sum_{j=1}^c P_i^D N_{i,j}(t)$$

r stands for the number of different types of tasks, ($r=5$ for Cholesky)

c stands for the number of threads/cores

P_i^D average dynamic power for task of type i

$N_{i,j}(t)$ equals to 1 if thread j is executing a task of type i at time t ; equals 0 otherwise

Energy model:

$$\begin{aligned} E_{Chol} &= (P^Y + P^S)T + \int_{t=0}^T P_{Chol}^D(t) \\ &= (P^Y + P^S)T + \sum_{i=1}^r \sum_{j=1}^c P_i^D \left(\int_{t=0}^T N_{i,j}(t) \right) = (P^Y + P^S)T + \sum_{i=1}^r \sum_{j=1}^c P_i^D T_{i,j} \end{aligned}$$

$T_{i,j}$ total execution time for task of type i onto the core j

Experimental model evaluation:

- Matrix sizes: $n = 4096, 8192, \dots, 32768$
- Block sizes $b = 128, 192, 256, 512$
- Cores/threads $c = 2, 3, \dots, 8$

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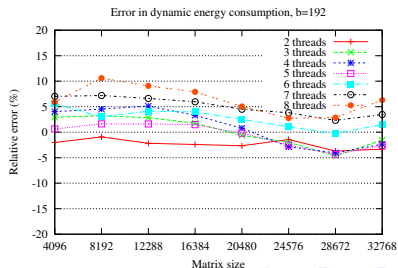
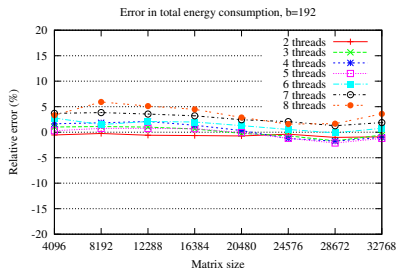
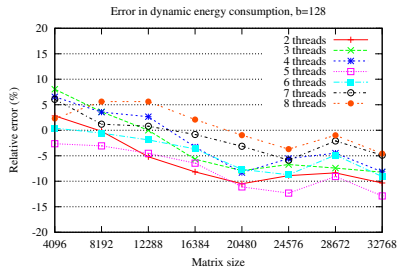
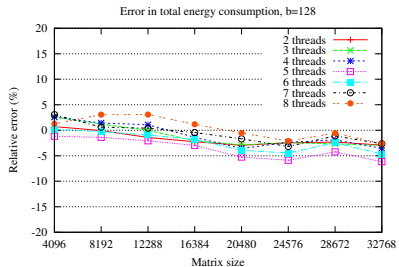
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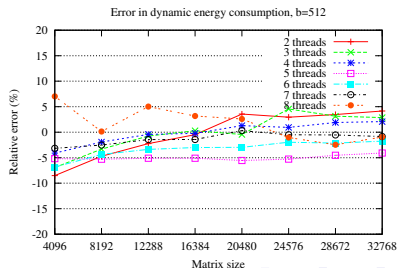
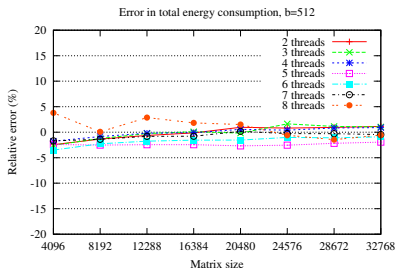
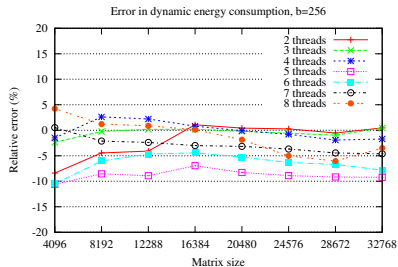
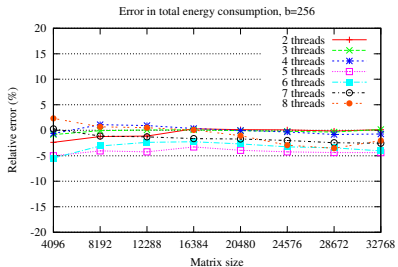
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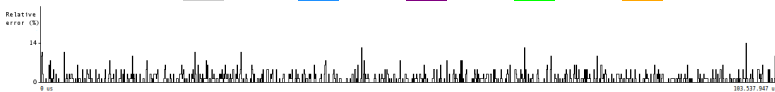
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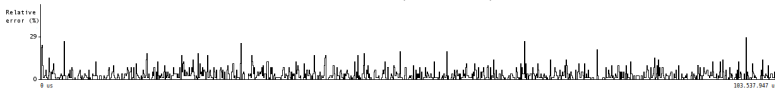
Power model evaluation

Reconstruction of power profile using the power model $\Rightarrow$ Performance trace is needed!

Trace of Cholesky factorization of order $n = 20,490$ and block size $b = 512$, using 4 cores



Relative error in the estimated total power (reconstruction); Average error of 2.92%



Relative error in the estimated dynamic power (reconstruction); Average error of 6.85%

Conclusions and future work

Conclusions:

- Elaboration and validation of an hybrid analytical-experimental model to estimate power/energy for the Cholesky factorization
- Experimental results reveal the accuracy of the model:
 - Energy consumption estimation: $\pm 5\%$ and $\pm 15\%$ of error for the total and dynamic energy, respectively
 - Power profile estimation: relative average error of 2.92% and 6.85% for total and dynamic power, respectively
- However, it is easier to obtain an energy estimation than a power profile estimation due to inaccuracy of power meter (around $\pm 5\%$)!

Future work:

- Predict power/energy even without executing the code!
- Initial step towards more ambitious goal $\Rightarrow$ Development of models for the functionality of LAPACK
- Model extension to task-parallel procedures for distributed-memory platforms

Thanks for your attention!

Questions?